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**HUMAN HEALTH RISK ASSESSMENT  
FOR THE FORMER GULF STATES CREOSOTING FACILITY,  
HATTIESBURG, MISSISSIPPI**

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## **Executive Summary**

A baseline human health risk assessment (HHRA) was conducted for the Former Gulf States Creosoting facility in Hattiesburg, Mississippi. The HHRA was performed in accordance with: Mississippi Commission on Environmental Quality's (MCEQ's) *Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi* (1999); US EPA's *Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A)* (1989); US EPA Region 4 guidance entitled *Technical Services Supplemental Guidance to RAGS, Region 4 Bulletins* (1995); and other relevant US EPA guidance documents.

Creosoting constituents of potential health concern include polycyclic aromatic hydrocarbons, of which benzo(a)pyrene is the predominant contributor to potential risks. Much of the former creosoting process area is currently covered with asphalt or large building structures. Potential future exposure scenarios included a construction worker, a maintenance worker, an infrequent Site visitor, and off-Site residents. Media of concern included soils, sediment, and surface water.

Hazards posed by chemical constituents in soils, sediment, and surface water for health effects other than an increased risk of cancer were well below a threshold of possible concern for each receptor evaluated in this risk assessment. Cancer risks for all exposure scenarios were within or below the US EPA's acceptable target risk range of  $1 \times 10^{-6}$  to  $1 \times 10^{-4}$  (*i.e.*, one in one million to one in ten thousand) with the exception of maintenance worker exposure to soils in EU4 and off-site resident exposure to sediments in EU6. The added lifetime cancer risk conservatively estimated for a maintenance worker was  $1 \times 10^{-3}$  for the entire Site, while that for the off-site resident was  $2 \times 10^{-4}$  for the entire Site. The potential risk for a construction worker was estimated to be  $5 \times 10^{-5}$  for the entire Site. The estimated potential risk for an adolescent Site visitor was  $2 \times 10^{-5}$  for the entire Site. For the Site visitor, maintenance worker, and construction worker scenarios, oral contact with benzo(a)pyrene in sediment and soils drove the cancer risk level. For the off-Site resident scenario, oral contact with benzo(a)pyrene in sediment drove the cancer risk level.

Risk levels are mainly attributable to residual concentrations of carcinogenic polycyclic aromatic hydrocarbons (cPAH) in EUs 4, 5, and 6. Remedial actions currently planned for these areas, including deed restrictions, will result in incomplete exposure pathways thereby resulting in acceptable levels of risks to potential receptors. Proposed remediation activities to address impacted media in EUs 4, 5, and 6 include the following:

- Conduct in-situ biological treatment of impacted soils in the unpaved area between the former Process Area and the Southern railroad tracks (EU4);
- Attempt to recover free product from targeted areas within the former Process Area to address continuing sources (EU5);
- Remove impacted sediments from the northeast drainage ditch and install a culvert to provide for surface drainage (EU6);

- Establish deed restrictions limiting the use of property to non-residential (i.e., “restricted”) purposes (EU4 and EU5); and
- Include in the deed restrictions provisions for maintaining pavement to preclude contact with impacted media left in place (EU5).

Constituent concentrations in surface soils at two isolated locations within EU2 also resulted in maintenance worker risk levels slightly greater than  $1 \times 10^{-6}$ . Because these locations are within a densely wooded area where no maintenance activities currently occur and remediation would require significant clearing, no remediation activities are planned to address surface soils at these locations. Deed restrictions limiting the use of properties within EU2 to non-residential purposes will be established.

## **1.0 Introduction**

Environmental Standards, Inc. (Environmental Standards) was retained by Kerr-McGee Chemical Corporation (Kerr-McGee) to perform a human health risk assessment (HHRA) to evaluate hazards and risks potentially posed by residual levels of chemicals present at the Former Gulf States Creosoting facility (Site). The Site, located near the intersection of US Highways 49 and 11 in Hattiesburg, Mississippi, was formerly a wood treating facility that operated between the early 1900s and 1960. In the early 1960s, the Site was redeveloped for commercial and light industrial uses (Michael Pisani & Assoc., 1997). The land on which the Site is located is a portion of the Sixteenth Section land owned by the Hattiesburg Public School District and leased to the current tenants under a 99-year lease, granted on July 7, 1947. At the time of this report, the Site, with the exception of the grassy and wooded areas in the south and southwest, respectively, was primarily used for automobile dealerships. There are no residential or institutional (*i.e.*, schools) uses of the Site (Michael Pisani & Assoc., 1997).

Operations at the Site consisted of a small-scale wood preserving process using creosote. The creosoting process was primarily confined to a 2.5-acre area in the northeast corner of the Site; this is known as the former Process Area and is currently occupied by Courtesy Ford. During the redevelopment of the Site in the early 1960s, construction debris (*e.g.*, broken concrete, asphalt, etc.) appears to have been relocated to the southwestern corner of the Site along Gordon's Creek. This area is known as the Fill Area and currently remains undeveloped.

This assessment has been conducted as a result of an agreement between Kerr-McGee, the Mississippi Department of Environmental Quality (MDEQ), and the Mississippi Commission on Environmental Quality (MCEQ) pursuant to the Uncontrolled Site Voluntary Evaluation Program. The MDEQ Office of Pollution Control, Uncontrolled Sites Section has been providing oversight and review of investigations and reports relating to the former Gulf States Creosoting facility.

This report will address the potential for on-Site exposures to human receptors and off-Site exposures to humans along the northeast drainage ditch.

The primary guidance used to develop this risk assessment was the MCEQ *Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi* (1999). US EPA Region 4's *Technical Services Supplemental Guidance to RAGS: Region 4 Bulletins* (1995) were also referred to for guidance. Additional US EPA guidance documents cited herein include:

- *Guidance for Remediation of Uncontrolled Hazardous Substance Sites in Mississippi (MDEQ, 1990);*
- *Risk Assessment Guidance for Superfund, Volume 1, Human Health Evaluation Manual/ Part A (RAGS/Part A)* (US EPA, 1989);
- *Human Health Evaluation Manual, Supplemental Guidance: "Standard Default Exposure Factors"* (US EPA, 1991);
- *Exposure Factors Handbook* (US EPA, 1997);
- *Guidelines for Exposure Assessment* (US EPA 1992);
- *Dermal Exposure Assessment: Principles and Applications* (US EPA, 1992);

These documents are not listed in a hierarchical manner; other US EPA guidance documents and peer-reviewed technical papers may have also been referenced in this risk assessment report.

## **2.0 Hazard Identification and Conceptual Site Model**

As a result of the historical wood preservation process, residual levels of creosote-related chemicals are present in soils in the former Process Area. Sediment and surface water in a drainage ditch along the southeast border of the former Process Area also contain chemical residuals. These Site-related chemicals, mostly polycyclic aromatic hydrocarbons (PAHs) are also present in the Fill Area. Residual levels of PAHs have been found in soil in the Fill Area and in Gordon's Creek surface water and sediment.

PAH residuals have also been detected in shallow groundwater underlying the Site. Currently, there are no private water wells located on-Site that access this shallow groundwater for potable purposes. The results of a door-to-door survey conducted by Michael Pisani and Associates on October 3, 2000 indicated no private uses of shallow groundwater downgradient of the Site. For these reasons, the groundwater exposure pathway, both on- and off-Site, was considered incomplete and not evaluated in this assessment.

A conceptual site model (CSM) was developed for the Site to aid in determining the potential receptors and exposure units to be evaluated under current and future potential land use (Figure 1). These receptors were identified as infrequent Site visitors, maintenance workers, construction workers, and off-Site residents.

Under current land use assumptions, Site visitors may potentially contact residual chemicals in Gordon's Creek surface water and sediment, and/or surface soils in the Fill Area and surrounding woods, the grassy field southeast of the Fill Area, and/or the drainage ditch along side of the former Process Area. Visitors may also potentially contact surface soil, surface water, and sediment along the former Process Area drainage ditch. The remaining affected areas of the Site are covered with either buildings or pavement precluding casual direct contact with surface soils. As a conservative measure, however, visitor exposure to soils from these paved areas was also assessed.

Under both current and future land use assumptions, a maintenance worker may contact surface soils in the Fill Area and surrounding woods, the grassy field southeast of the Fill Area, and/or the former Process Area and surrounding affected areas, including the drainage ditch located to the southeast of the former Process Area. Although most of the former Process Area and vicinity are paved, maintenance activities may involve some shallow digging; therefore, direct contact with shallow soils in this area was assessed.. As a conservative measure, exposure to surface water and sediment in Gordon's Creek was assessed. The remainder of the Site was relatively unaffected by historical creosoting activities.

Although there are currently no major construction activities at the Site, these types of activities may occur at some time in the future. As with the maintenance worker scenario, construction activities could potentially occur in the Fill Area and vicinity, the grassy field southeast of the Fill Area, and the former Process Area and vicinity. Construction workers may be exposed to both surface and subsurface soils (down to the water table). Construction worker exposure to surface water and sediment in Gordon's Creek was assessed as a conservative measure. The remainder of the Site was relatively unaffected by historical creosoting activities.

Areas of the Site affected by historical creosoting activities will be deed restricted prohibiting future residential development. Off-Site areas along the northeast Drainage Ditch, currently a residential neighborhood, were assessed for residential exposures to soil, sediment, and surface water.

### **3.0 Data Evaluation**

To characterize potential exposures to Site-related chemicals, the former Gulf States Creosoting facility was divided into six exposure units (EUs). Each exposure unit outlines potentially affected areas of the Site and adjacent on-Site locales that may be frequented by individuals accessing the Site for recreational or occupational purposes. The use of EUs is encouraged by the US EPA Region 4 (1995), which defines an EU as “an areal extent of a receptor’s movements during a single day... .” Each of these exposure units is depicted on Figure 2 and is discussed below.

A sixth EU was created for off-Site residential exposures to surface water and sediment along the northeast Drainage Ditch. This EU is delineated on Figure 3.

#### **3.1 Exposure Unit Delineation**

The following EUs were delineated based upon the presence of residual chemicals and the potential for receptors to contact those chemicals. Areas of the Site most affected were included in at least one of the five EUs while areas with relatively low or non-detectable concentrations of residuals were not included in an EU. By limiting Site-wide exposures to the EUs most affected by historical activities at the Site, worst-case scenarios were created.

##### **3.1.1 Exposure Unit 1**

EU1 outlines the on-Site areas in, adjacent to, and downstream of the Fill Area along Gordon’s Creek (Figure 2). EU1 includes exposures to surface water and sediment by an infrequent Site visitor, future maintenance worker, and future construction worker. Although US EPA Region IV guidance indicates that “In most cases it is unnecessary to evaluate human exposures to sediments covered by surface water,” (US EPA, 1995) dermal and oral surface water exposures were conservatively assessed herein at the request of the MDEQ (2000). Sediment samples included in EU 1 were SD07 and SD08. Surface water samples included in were SW-07 and SW-08.

Soil samples from this area were considered part of EU2 and exposures were assessed accordingly.

### 3.1.2 Exposure Unit 2

EU2 delineates the upland areas of the Fill Area and adjacent woody and grassy areas (Figure 2). Surface soils from zero to one foot and zero to six feet below ground surface [bgs] in this area were evaluated for potential visitor and future hypothetical maintenance worker scenarios, respectively. Surface and subsurface soils were also evaluated for a hypothetical future construction worker scenario. Available data for subsurface soils for a construction scenario were evaluated from the surface to the water table (approximately 10 feet bgs) as recommended by the MDEQ (2000). Soil samples included in EU2 are presented in the table below:

| Soils (0-1' bgs)  | GEO-13/0-1' | SS-1         | SS-2        | SS-3        | SS-4        |
|-------------------|-------------|--------------|-------------|-------------|-------------|
| SS-5              | SS-6        | SS-7         | SS-8        | SS-9        |             |
| SS-10             | SS-11       | SS-12        | SS-13       |             |             |
| Soils (0-6' bgs)  | GEO-03/2-3' | GEO-03/5-6'' | GEO-10/2-3  | GEO-10/5-6  | GEO-13/0-1' |
| GEO-13/2-3'       | GEO-13/5-6' | GEO-44/5-6'  | SS-1        | SS-2        |             |
| SS-3              | SS-4        | SS-5         | SS-6        | SS-7        |             |
| SS-8              | SS-9        | SS-10        | SS-11       | SS-12       |             |
| SS-13             |             |              |             |             |             |
| Soils (0-10' bgs) | GEO-03/2-3' | GEO-03/5-6'  | GEO-10/2-3  | GEO-10/5-6' | GEO-13/0-1' |
| GEO-13/2-3'       | GEO-13/5-6' | GEO-43/7-8'  | GEO-44/5-6' | GEO-45/7-8' |             |
| SB-03/8-9.3       | SB-05/4-9   | SB-07/5-7    | SS-1        | SS-2        |             |
| SS-3              | SS-4        | SS-5         | SS-6        | SS-7        |             |
| SS-8              | SS-9        | SS-10        | SS-11       | SS-12       |             |
| SS-13             |             |              |             |             |             |

### 3.1.3 Exposure Unit 3

In the southwest corner of the Site there exists a grassy field east of West Pine Street between Henson Auto Sales and Eagan Cars and Trucks. This grassy area has been defined as EU3 for

purposes of this risk assessment (Figure 2). Similar to EU2, surface soil from zero to one foot and zero to six feet bgs were evaluated in EU2 for visitor and hypothetical future maintenance worker scenarios, respectively. Surface and subsurface soils in this EU were evaluated for a hypothetical future construction worker scenario. Available data for subsurface soils for a construction scenario were evaluated from the surface to the water table(approximately 20 feet bgs) as recommended by the MDEQ (2000). Soil samples included in EU3 are presented in the table below:

|                            |             |             |             |             |       |
|----------------------------|-------------|-------------|-------------|-------------|-------|
| Soils (0-1' bgs)           | SS-15       | SS-16       | SS-17       |             |       |
| Soils (0-6' and 0-20' bgs) | GEO-16/2-3' | GEO-16/5-6' | GEO-17/2-3' | GEO-17/5-6' | SS-15 |
|                            | SS-16       | SS-17       |             |             |       |

### 3.1.4 Exposure Unit 4

EU 4 encompasses the grassy drainage ditch area along the fenceline behind Courtesy Ford in the northeast corner of the Site and continues parallel to the railroad tracks, and west through EU 3 and EU 2 (Figure 2). EU 4, along the southeast side of the former Process Area, has been widened to include soil data from that area. Receptors associated with EU 4 included Site visitor exposures via casual contact with surface soil, sediment, and surface water. Maintenance worker and construction worker scenarios were also evaluated for exposures to surface water and sediment in EU 4 as well as soils in EU 4 near the former Process Area. Soils down to six feet bgs were evaluated for maintenance workers while soils down to the water table (approximately20 feet bgs) were evaluated for construction workers in this EU as requested by the MDEQ (2000). Sediment, surface water, and soil samples included in EU4 are presented in the following table:

|                   |              |              |             |             |             |
|-------------------|--------------|--------------|-------------|-------------|-------------|
| Sediment          | SD-02        | SD-12        | SD-18       | SD-19       | SD-20       |
|                   | SD-21        | SD-22        | SD-23       |             |             |
| Surface Water     | SW-02        |              |             |             |             |
| Soils (0-1' bgs)  | GEO-19/0-1'  | GEO-20/0-1'  | GEO-21/0-1' | GEO-46/0-1' | GEO-47/0-1' |
|                   | GEO-48/0-1'  |              |             |             |             |
| Soils (0-6' bgs)  | GEO-19/0-1'  | GEO-19/2-3'  | GEO-19/5-6' | GEO-20/0-1' | GEO-20/2-3' |
|                   | GEO-20/5-6'  | GEO-21/0-1'  | GEO-21/2-3' | GEO-21/5-6' | GEO-46/0-1' |
|                   | GEO-46/2-3'  | GEO-46/5-6'  | GEO-47/0-1' | GEO-47/2-3' | GEO-47/5-6' |
|                   | GEO-48/0-1'  | GEO-48/2-3'  | GEO-48/5-6' |             |             |
| Soils (0-20' bgs) | GEO-19/0-1'  | GEO-19/2-3'  | GEO-19/5-6' | GEO-20/0-1' | GEO-20/2-3' |
|                   | GEO-20/5-6'  | GEO-20/9-10' | GEO-21/0-1' | GEO-21/2-3' | GEO-21/5-6' |
|                   | GEO-21/9-10' | GEO-46/0-1'  | GEO-46/2-3' | GEO-46/5-6' | GEO-47/0-1' |
|                   | GEO-47/2-3'  | GEO-47/5-6'  | GEO-47/7-8' | GEO-48/0-1' | GEO-48/2-3' |
|                   | GEO-48/5-6'  |              |             |             |             |

### 3.1.5 Exposure Unit 5

EU5 outlines the former Process Area and the historical drip track and treated wood storage areas of the former Gulf States Creosoting facility (Figure 2). Surface soils from zero to six feet bgs were evaluated in EU5 for a hypothetical maintenance worker scenario. Available data for soils down to the water table (approximately 20 feet bgs) were evaluated in EU5 for a hypothetical future construction worker scenario. Soil samples included in EU5 are presented in the table below:

|                   |             |             |                 |             |             |
|-------------------|-------------|-------------|-----------------|-------------|-------------|
| Soils (0-1' bgs)  | GEO-28/0-1' | GEO-29/0-1' | GEO-30/0-1'     | GEO-31/0-1' | GEO-32/0-1' |
|                   | GEO-33/0-1' | GEO-59/0-1' | GEO-60/0-1'     |             |             |
| Soils (0-6' bgs)  | GEO-28/0-1' | GEO-28/2-3' | GEO-28/5-6'     | GEO-29/0-1' | GEO-29/2-3' |
|                   | GEO-29/5-6' | GEO-30/0-1' | GEO-30/2-3'     | GEO-30/5-6' | GEO-31/0-1' |
|                   | GEO-31/2-3' | GEO-31/5-6' | GEO-32/0-1'     | GEO-32/2-3' | GEO-32/5-6' |
|                   | GEO-33/0-1' | GEO-33/2-3' | GEO-33/5-6'     | GEO-59/0-1' | GEO-59/2-3' |
|                   | GEO-59/5-6' | GEO-60/0-1' | GEO-60/2-3'     | GEO-60/5-6' |             |
| Soils (0-20' bgs) | GEO-28/0-1' | GEO-28/2-3' | GEO-28/5-6'     | GEO-29/0-1' | GEO-29/2-3' |
|                   | GEO-29/5-6' | GEO-30/0-1' | GEO-30/2-3'     | GEO-30/5-6' | GEO-31/0-1' |
|                   | GEO-31/2-3' | GEO-31/5-6' | GEO-32/0-1'     | GEO-32/2-3' | GEO-32/5-6' |
|                   | GEO-33/0-1' | GEO-33/2-3' | GEO-33/5-6'     | GEO-59/0-1' | GEO-59/2-3' |
|                   | GEO-59/5-6' | GEO-60/0-1' | GEO-60/2-3'     | GEO-60/5-6' | GEO-60/7-8' |
|                   | SB-01/8-10  | SB-02/9-11  | SB-05/10.5-12.5 | SB-06/6-10  | SB-07/14-16 |

### 3.1.6 Exposure Unit 6

EU6 outlines a stretch (approximately 2700 feet in length) of the northeast drainage ditch that leads from the Site into the neighboring residential area. EU6 exposures include oral and dermal exposures by off-Site residents to sediment and surface water along the northeast drainage ditch. Soil exposures were not assessed in this area for lack of soil data. Also, it was anticipated that sediment exposures in this area represent a more conservative estimate of exposure in that chemical concentrations in the exposed sediment along the drainage ditch are likely to be greater than concentrations in the surrounding soils. Sediment and surface water samples included in EU6 are presented in the table below:

|               |       |       |       |       |
|---------------|-------|-------|-------|-------|
| Sediment      | SD-03 | SD-04 | SD-05 | SD-13 |
|               | SD-14 | SD-15 | SD-16 | SD-17 |
| Surface Water | SW-03 | SW-04 |       |       |

average concentration as the exposure-point concentration by explaining that toxicity criteria for both carcinogenic and non-carcinogenic effects are based on lifetime average exposures and that the “average concentration is most representative of the concentration that would be contacted at a site over time” (*Supplemental Guidance to RAGS: Calculating the Concentration Term*, 1992). Other US EPA guidance states that “...in most situations, assuming long-term contact with the maximum concentration is not reasonable” (*Risk Assessment Guidance for Superfund, Part A*, 1989). US EPA Region 4 also states that, generally, it is reasonable to assume that soil data are distributed lognormally (1995). In keeping with these guidances, the lognormal 95% UCL was considered in the screening process where the data distribution for a compound could not be defined as specifically normal or lognormal.

If the 95% UCL (or lognormal 95% UCL where appropriate) of a constituent was less than the Tier 1 TRG, then that constituent was eliminated from further quantitative analysis. If the 95% UCL (or lognormal 95% UCL where appropriate) of a constituent in soil exceeded the Tier 1 TRG, then that constituent was retained for quantitative analysis in the Site-specific risk assessment (Tier III).

MCEQ guidance (1999) does not specify screening levels for constituents in sediment or surface water; therefore, Region 4 was referred to for guidance (1995). Sediment is only found on the Site in drainage ditches that contain little to no water most of the time. US EPA Region 4 guidance states that sediments in an intermittent stream (or ditch) should be considered as surface soil for the portion of the year the stream is without water. Based on these factors and comments provided by the MDEQ (2000), the maximum detected constituent concentrations in sediment was compared to MCEQ unrestricted Tier 1 TRGs. The screening process then followed the same procedure as mentioned above for other soils.

For surface water, the maximum detected concentration of a constituent in an EU was compared to the US EPA Human Health Water Quality Standard (WQS) for consumption of water and organisms in accordance with US EPA Region 4 guidance (1995). If the maximum

### 3.2 Statistical Evaluation

Environmental samples undergo laboratory analyses that are designed to quantitate the concentrations of constituents in the various environmental media. As a result of the analytical procedures, a constituent may be detected and its concentration measured, detected but not able to be quantitated, or not detected at all in a sample. The data set for the Site contains a number of nondetections for some chemicals of potential concern (COPCs) in various samples. Assuming that the COPC is present in these samples at the achieved detection limit is biased because the chemical may be absent altogether. Assuming a concentration of zero is also flawed because the chemical could be present at a level below laboratory capabilities to detect and quantify the concentration. Consequently, in the event that an analyte identified at least once in a given medium was not detected in a given sample, it was conservatively assumed for the risk assessment purposes to be present at a concentration equivalent to one-half of the sample quantitation limit (SQL). In addition, samples labeled with an "R" (rejected) qualifier were not included in the data analysis because those data were deemed unreliable and, therefore, unusable. Constituents that were not detected in any sample from a particular medium were eliminated from further consideration in accordance with US EPA guidelines (1989).

Site analytical data used in this assessment were collected during the Phase I (1997) and Phase II (1998) remedial investigations as well as the additional investigation conducted in 2000 at the request of the MDEQ. These data were fully validated by qualified technical professionals using standard data validation protocols, as required by the MCEQ (1999).

Previous investigations at the Site have been conducted since 1990. These investigations included the following:

- 1990 soil gas and soil sampling by Roy F. Weston
- 1991 MDEQ Site inspections and Phase II report
- 1994 Phase II Site investigation by Environmental Protection Systems (EPS)
- 1994 Site investigation by Bonner Analytical Testing Company (BATCO)
- 1994 preliminary subsurface investigation by BATCO

- 1995 three-dimension resistivity surveys by American Remediation Technology
- 1996 investigation by McLaren/Hart
- 1996 investigation by Kerr McGee Chemical Corporation

Data acquired from these historical (pre-1997) investigatory activities were not used in this assessment as they were not validated by qualified chemists and sampling locations for some of the data could not be accurately established. These historical data were not considered valid and were, therefore, not appropriate to use in this assessment of risks. Only validated data that were considered to be representative of Site conditions with a reasonable level of confidence were used for this assessment.

The validated laboratory data from 1997, 1998, and 2000 investigations were compiled into data sets representing areas of potential exposure (EUs) for each potential receptor. Each data set was analyzed statistically using SiteStat®, a commercially available software package, to calculate the minimum, maximum, arithmetic mean, logarithmic mean, standard error of the mean, and the 95% upper confidence limit of the mean concentration (95% UCL) for each constituent based on distributional analysis of the data (*i.e.*, utilizing goodness-of-fit statistical tests to determine whether the data are distributed normally or lognormally). The data qualifier associated with the minimum and maximum detected concentrations as well as the location of the maximum detected concentration for each EU were also determined. Results of the quantitative and statistical analyses for each of the EUs discussed above are presented in Tables 1 through 18.

Standard sampling protocol requires the collection of duplicate field samples used to ensure the quality of a laboratory analysis (*i.e.*, to ensure that analytical results can be replicated). As such, duplicate sample results were provided as part of the database for the Hattiesburg Site. In accordance with US EPA guidance (1989), duplicate sample results were averaged (for any sample containing duplicates) and the average concentration was used as a single concentration for that sample in the calculation of summary statistics as discussed below.

Soils down to one foot deep were assumed to be representative of surface soils at the Site for infrequent visitor exposures. A depth of 0 to 6 feet was used to define surface soils for maintenance worker exposures. These assumptions were recommended by the MDEQ (2000). The groundwater table was considered the extent of subsurface soils as recommended by MDEQ (2000). This value (depth-to-groundwater) varies significantly across the Site and, as such, the extent of subsurface soil was EU-specific as follows:

EU2 – soils down to 10 feet

EU3 – soils down to 20 feet

EU4 – soils down to 20 feet

EU5 – soils down to 20 feet

This risk assessment focuses mainly on environmental data collected from the former Process and Fill Areas and any other portions of the Site that were affected by former creosoting operations. Virtually unaffected areas (e.g., the developed area north of West Pine Street) as delineated using historical data were not considered to contribute significantly to risk levels and, therefore, were excluded from this risk assessment.

### 3.3 Determination of Exposure-Point Concentrations

Exposure-point concentrations were determined to be the 95% UCL or the maximum concentration of a COPC in an EU, whichever was lower. This methodology is in accordance with US EPA guidance (1989). If the distribution of the concentration data was determined to be lognormal, then the lognormal 95% UCL was compared to the maximum concentration to determine the exposure-point concentration. In the event that the distribution of a chemical in any given medium could not be confidently labeled as normal or lognormal, it is termed either “unknown” or “normal/lognormal.” In these cases, the lognormal 95% UCL was compared to maximum concentration when determining the exposure-point concentration. It should be noted, however, that in cases where the distribution is “unknown,” the normal and lognormal 95% UCLs could not be reliably predicted. Assuming a lognormal distribution of the data increases

the uncertainty associated with this step of the risk assessment process; however, hazard and risk estimates are likely to be less uncertain than if the maximum concentrations were used.

Exposure-point concentrations are provided on the statistical summary tables, Tables 1 through 18.

### 3.4 COPC Selection

Soils (both surface and subsurface) were screened according to MCEQ (1999) guidance. The first tier of the screening process compared maximum concentrations of a constituent in an EU with the Restricted Tier 1 target remediation goal (TRG) for maintenance worker and construction worker scenarios. Restricted TRGs were used because the Site is not currently used for residential purposes and the current commercial/industrial land-use is anticipated to remain into the future as a result of the implementation of deed restrictions on the impacted areas of the Site. If a maximum concentration of a constituent was less than the Restricted Tier 1 TRG, then that constituent was eliminated from further quantitative assessment.

Surface soil data (zero to one foot bgs) for the visitor scenario were screened using Unrestricted Tier 1 TRGs at the request of MDEQ (2000). If a maximum concentration of a constituent was less than the Unrestricted Tier 1 TRG, then that constituent was eliminated from further quantitative assessment. Conversely, if the maximum concentration of a constituent exceeded the Tier 1 TRG, that constituent was retained for quantitative analysis.

If the maximum concentration of a constituent in an EU exceeded the Tier 1 TRG, then the 95% UCL of the constituent was compared to the Tier 1 TRG (Restricted or Unrestricted, depending on the exposure scenarios as described above) as part of the Tier II screening process. In the event that the concentrations of a chemical were distributed lognormally, the lognormal 95% UCL of that constituent was compared to the Tier 1 TRG. If the distribution of data of a chemical could not be positively identified as either normal or lognormal, the lognormal 95% UCL was used in the screening process. In these cases, either the maximum concentration or the lognormal 95% UCL can be conservatively used. The US EPA, however, justifies the use of an

concentration of a constituent in surface water was less than the WQS, then that constituent was eliminated from quantitative analysis. If the maximum concentration of a constituent in surface water exceeded the WQS, then that constituent was retained for quantitative analysis.

At the request of MDEQ (2000), if any single carcinogenic polycyclic aromatic hydrocarbon (cPAH) was retained as a COPC in a medium, then all cPAHs were also retained as COPCs in that medium. This guidance refers to the following chemicals: benzo(a)anthracene, benzo(b)fluoranthene, benzo(k)fluoranthene, benzo(a)pyrene, chrysene, dibenzo(a,h)anthracene, and indeno(1,2,3-cd)pyrene. To establish an exposure point concentration for undetected cPAHs retained as COPCs in an EU, one-half the maximum detection limit was used.

The results of the screening process are presented on the statistical summary tables, Tables 1 through 18. The screening process eliminated detected constituents from the subsurface soil dataset down to 20 feet bgs and surface soil dataset down to 6 feet bgs in EU3. For this reason, construction worker and maintenance worker exposures to soils in EU3 were not evaluated quantitatively in this assessment.

## **4.0 Exposure Assessment**

Currently, a majority of the Site is used for commercial and light industrial purposes and is paved for roads and parking lots. Unpaved areas are limited to Gordon's Creek (EU 1), the wooded portion in and around the Fill Area (EU2) and the grassy field outlined by EU 3, and the drainage ditches and surrounding area delineated by EU 4 (Figure 2). Since the developed and undeveloped areas of the Site vary considerably with respect to both residual chemical concentrations and land use, the Site was divided into five EUs for the exposure assessment. A sixth EU was created to assess off-Site residential exposures. Chemical data from each EU were combined with EU-specific exposure parameter values and receptor scenarios to determine the chemical intake for each receptor potentially accessing an EU for occupational, recreational, or residential purposes.

### **4.1 Receptor Identification**

The following exposures pathways (indicated with an "X") have been selected for this risk assessment as reasonable and realistic scenarios under current and future land-use assumptions:

| EU/Media:         | EU1  |             | EU2  |      | EU3  |      | EU4         |      | EU5  |             | EU6  |             |
|-------------------|------|-------------|------|------|------|------|-------------|------|------|-------------|------|-------------|
| Receptor/Route:   | Sed. | Surf. Water | Soil | Soil | Soil | Sed. | Surf. Water | Soil | Sed. | Surf. Water | Soil | Surf. Water |
| Visitor           |      |             |      |      |      |      |             |      |      |             |      |             |
| Dermal            | X    | X           | X    | X    | X    | X    | X           |      |      | X           |      |             |
| Oral              | X    | X           | X    | X    | X    | X    | X           |      |      | X           |      |             |
| Inhalation        |      |             |      |      |      |      |             |      |      |             |      |             |
| Maint. Worker     |      |             |      |      |      |      |             |      |      |             |      |             |
| Dermal            | X    | X           | X    | X    | X    | X    | X           |      |      | X           |      |             |
| Oral              | X    | X           | X    | X    | X    | X    | X           |      |      | X           |      |             |
| Inhalation        |      |             |      |      |      |      |             |      |      |             |      |             |
| Const. Worker     |      |             |      |      |      |      |             |      |      |             |      |             |
| Dermal            | X    | X           | X    | X    | X    | X    | X           |      |      | X           |      |             |
| Oral              | X    | X           | X    | X    | X    | X    | X           |      |      | X           |      |             |
| Inhalation        |      |             | X    | X    | X    |      |             |      |      | X           |      |             |
| Off-Site Resident |      |             |      |      |      |      |             |      |      |             |      |             |
| Dermal            |      |             |      |      |      |      |             |      | X    | X           |      |             |

| EU/Media:       | EU1  |             | EU2  | EU3  | EU4  |      |             | EU5  | EU6  |             |
|-----------------|------|-------------|------|------|------|------|-------------|------|------|-------------|
| Receptor/Route: | Sed. | Surf. Water | Soil | Soil | Soil | Sed. | Surf. Water | Soil | Sed. | Surf. Water |
| Oral            |      |             |      |      |      |      |             |      |      | X           |
| Inhalation      |      |             |      |      |      |      |             |      |      |             |

Surface water present on-Site is either ephemeral or very shallow and is conducive only to wading-type activities. Ingestion of Site surface water was considered an insignificant exposure pathway since on-Site drainage ditches "contain little or no water most of the time" (MDEQ, 2000). In addition, US EPA IV guidance indicates that "In most cases, it is unnecessary to evaluate human exposures to sediments covered by surface water" (1995). At the request of MDEQ (2000), however, dermal and oral exposures to surface water were assessed for visitors, maintenance workers, and construction workers in EU's 1 and 4. Surface water exposures were also assessed for residents in off-Site EU 6.

Each of the potential receptors is discussed below.

#### 4.1.1 Infrequent Site Visitor

Since the Site is not currently fenced or guarded, the general public has access to most areas of the Site at any given time. It is possible, though unlikely, that an individual may use some areas of the Site, such as EU1, EU2, or EU3, for recreational purposes. For this reason, sediment and surface water exposures to visitors in EU1, and surface soil exposures in EU2 and EU3 were assessed for the visitor scenario. The vast majority of the remainder of the Site (EU5) is covered with either buildings or pavement, precluding direct contact with surface soils; however, a small exposed area encompassing a drainage ditch exists along side of the former Process Area (EU4). Although this area is not attractive for recreational purposes, it is possible that an individual traversing the Site may contact surface soils, sediment, or surface water in this EU; therefore, these potential exposures were assessed. Sediment exposures in EU1 and EU4 were addressed in accordance with US EPA Region 4 guidance that recommends evaluating sediment exposures in intermittent streams. At the request of MDEQ (2000), soil exposures were assessed for visitors

in EU5 regardless of the existence of buildings and pavements precluding almost all potential direct contact with soils in this area.

#### 4.1.2 Maintenance Worker

Currently, maintenance activities are most likely limited to the developed portions of the Site. Of these, the former Process Area and adjacent former drip track and treated wood storage areas (EU5) were most affected by historical wood preserving processes. Although these areas are mostly paved or built upon, it is possible that maintenance activities may require some shallow digging in unpaved areas; therefore, exposures to surface soils in EU5 were assessed. As a conservative measure, surface soil data from sample locations located in paved areas were evaluated in conjunction with surface soil data from exposed areas in EU5. If the currently undeveloped portions of the Site (EU2 and EU3) become developed in the future, similar maintenance activities may be required and, therefore, exposures to surface soils in EU2 and EU3 were also assessed. The drainage ditch encompassed by EU4 requires periodic maintenance; therefore, exposures to soil, sediment, and surface water in this area were assessed. At the request of MDEQ (2000), maintenance worker exposures to surface water and sediment in EU 1 were also assessed.

#### 4.1.3 Construction Worker

Although there are currently no major construction activities at the Site, such activities may hypothetically occur in the future. Thus, exposures to surface water and sediment in EUs 1 and 4, and exposures to soil in EUs 2 through 5 were assessed herein. Construction workers may be exposed to both surface and subsurface soils during activities such as excavating. Subsurface soils, for purposes of this assessment, were defined as those soils at the water table and shallower. Since the depth to the water varies significantly across the Site, so does the definition of "subsurface" soils. Accordingly, subsurface soils were evaluated down to 10 feet for EU2 and 20 feet for EUs 3, 4, and 5.

#### 4.1.4 Future On-Site Residents

The affected areas of the Property (the Site) are currently zoned for industrial or light-commercial use, and, at the time of this report, there were no plans to develop the Site for residential housing. In fact, deed restrictions preventing residential development are in the process of being implemented for the impacted areas on Site. Because of these deed restrictions, it is reasonable and realistic to assume that the Site will remain commercial/industrial in the future; therefore, on-Site residential exposures were not addressed in this risk assessment.

#### 4.1.5 Off-Site Residential Exposures

The northeast drainage ditch extends from the former Process Area to the northeast into a nearby residential community. Surface water and sediment data from areas along the northeast drainage ditch (EU6, Figure 3) were evaluated for off-Site residential exposures. For purposes of exposure assessment, a child resident between the ages of 1 and 6 years and an adolescent/adult resident between the ages of 7 and 30 years were evaluated. Hazards and risks for these two receptors were then combined (summed) to reflect the exposures incurred by a single individual living off-Site in the vicinity of the northeast drainage ditch for 30 years.

### 4.2 General Intake Equation

Chemical exposure/intake is expressed as the amount of the agent at the exchange boundaries of an organism (*i.e.*, skin, lungs, gut) that is available for systemic absorption. An applied dose is defined as the amount of a chemical at the absorption barriers such as skin, lung, digestive tract, available for absorption and is (usually expressed in milligrams, or mg) absorbed per unit of body weight of the receptor (usually expressed in units of kilogram, or kg). Absorbed dose can be defined as the amount of chemical that penetrates the exchange boundaries. If the exposure occurs over time, the total exposure can be divided by the time period of interest to obtain an average exposure rate (*e.g.*, mg/kg-day). The general equation, as defined by US EPA, for estimating a time-weighted average intake is:

$$\text{Intake (mg/kg - day)} = \frac{C \times IR \times EF \times ED}{BW \times AT} \quad [\text{Equation 1}]$$

where:

|    |   |  |
|----|---|--|
| C  | = | chemical concentration at the exposure point (e.g., mg/m <sup>3</sup> air);        |
| IR | = | intake rate (e.g., m <sup>3</sup> /hr);  |
| EF | = | exposure frequency (days/year);  |
| ED | = | exposure duration (years);   |
| BW | = | body weight of exposed individual (kg); and  |
| AT | = | averaging time (period over which exposure is averaged, usually measured in days). |

Additional parameters (e.g., skin surface area) were incorporated into the above general equation to evaluate the different potential exposure routes (dermal, oral, inhalation).

Table 19 presents the general and pathway-specific exposure parameters utilized for the intake equations in this assessment.

#### 4.2.1 General Exposure Parameters

Although some of the parameters used to calculate potential exposure are pathway- or route-specific, exposure frequency (EF), exposure duration (ED), averaging time (AT; determined separately for carcinogenic and non-carcinogenic exposures), and body weight (BW) are present in each intake model. These general parameters remain consistent throughout the intake calculations for each specific receptor.

##### 4.2.1.1 Exposure Frequency

The exposure frequency (EF) describes the number of times per year an event is likely to occur. It is most often expressed in units of days/year or events/year, depending on the scenario. Variables such as weather, vacations, sick days, and institutional controls often aid in determining reasonable and realistic exposure frequencies.

The EF for an adolescent visitor was extracted from US EPA *Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A) Interim Final* (1989). This EF value of 12 days/year per EU is a reasonable estimate that assumes an adolescent would most likely be engaged in outdoor activity on the unpaved areas of the Site for one day a week during

the three warmest months of the year. This value was used for soil, sediment, and surface water exposures.

Typical construction projects, especially at industrial complexes, generally involve several phases of activity prior to completion. The EF parameter used for oral exposure in construction workers, therefore, was subdivided into two exposure events. The first event hypothetically lasts for 10 days (used in relevant exposure model calculations under "Exposure Level A") and would involve earth-moving activities such as foundation. The second exposure event to the same individual hypothetically lasts for 70 days (for a total of 80 days at the Site for an individual; this value was used in relevant exposure model calculations under "Exposure Level B") and included remaining construction activities such as building framing, plumbing installation, electrical installation, and roofing. Generally, to complete each of these phases, a different team of specialized contractors is employed to perform the tasks for which they are most qualified. As a result, an individual may only remain at the construction site for a few days or weeks until his/her task has been completed and the next phase has begun. This is especially true for those activities involving direct contact with soil such as excavating and foundation pouring. Individuals performing these tasks are not usually qualified or employed to continue with the actual building processes. For dermal and inhalation exposures, however, an 80-day EF was used and accounted for an individual to be involved in construction activities for four entire months of the year (assuming five-day work weeks).

For surface water and sediment exposures to construction workers, an EF value of 8 days/year was used. This value represents 1/10<sup>th</sup> of the time a worker may be on-Site for construction-type activities and is conservative in that it is unlikely that construction workers would be exposed at all to Site surface water or sediment.

The EF value used for the maintenance worker scenario was 150 days/year for surface soil exposures in EUs 2, 3, and 5. This is also a conservative assumption in that the currently developed areas of the Site are covered with buildings or pavement. Maintenance activities in

these areas would require little contact with the obscured surface soils. The undeveloped areas of the Site currently require little or no maintenance as they are only occasionally mowed or allowed to grow naturally. Should these areas become developed, they will most likely take on the appearance of the remainder of the Site, including industrial/commercial buildings and paved roads or parking lots. Once again, extensive direct contact with surface soils would be minimal for a maintenance worker.

For maintenance worker sediment and surface water exposures in EUs 1 and 4 and surface soil exposures in EU 4, an EF value of 30 days/year was used. Historically, the northeast drainage ditch has been maintained on an as-needed basis (less than annually). Maintenance worker exposures to sediment and surface water in these areas were assessed at the request of the MDEQ (2000). An EF value of 30 days/year is amply conservative in that both Gordon's Creek (EU 1) and the northeast drainage ditch (EU 4) are currently maintained less than annually.

For residential soil exposures, an exposure frequency of 350 days/year was used in accordance with Region IV guidance. This value assumes that 15 days/year are spent away from home (US EPA, 1991).

Sediments along the bank of the northeast drainage ditch are not comparable to surface soils comprising a yard with respect to exposure. Typically, yard soils include relatively large areas where children frequently play and where surface soils are tracked into the home to become part of the household dust that can be ingested, particularly by crawling infants, on a daily basis. These are the assumptions that underlie the standard residential soil exposure algorithm and parameter values. However, it is not realistic to assume that infants, children, or adults will directly contact a relatively small area of sediments on the banks of a drainage ditch on a daily basis. A more realistic exposure scenario for this unique area under an assumption of residential land use is for a resident child to play on occasion in the drainage ditch that traverses the residential property. An exposure frequency of 40 days/year, two hours per exploring event, is conservatively plausible.

#### 4.2.1.2 Exposure Duration

The ED parameter represents the number of years during which an event is likely to occur. Factors affecting this parameter include variables such as age of receptor, population mobility, and occupational mobility. Exposure durations of less than seven years typically correspond to subchronic exposures while those greater than seven years are typically considered chronic exposures (US EPA, 1989). Toxicity indices are selected based on subchronic or chronic exposure durations.

The future construction worker scenario used an ED of one year because it is highly unlikely that a future construction worker would remain on one site for more than a year. Often, two months is considered the maximum amount of time a construction worker may reasonably remain at the same site.

The future maintenance worker ED, on the other hand, is based on occupational mobility studies. The ED of 25 years was obtained from US EPA (1991) which recommends a 95th percentile value of 25 years based on a study by the Bureau of Labor Statistics as of 1987. US EPA Region 4 also recommends a default value of 25 years for worker scenarios (1995).

The adolescent visitor scenario used an ED of 10 years. An adolescent was defined in this assessment as an individual aged seven to 16 years in accordance with US EPA Region 4 (1995); therefore, an exposure duration of 10 years was most appropriate.

An ED of 30 years (US EPA Region 4, 1995) was used for off-Site residents. This value assumes an individual spends 6 years as a child and 24 years as an adolescent/adult in the same location.

#### 4.2.1.3 Averaging Time

The averaging time (AT) parameter is the time period over which exposure is averaged. For human health cancer risk calculations, the AT<sub>c</sub> value prorates a total cumulative dose over a

lifetime. As a conservative approach, the AT<sub>c</sub> value for each receptor is the product of a 365-day year and a 70-year life span, equaling 25,550 days.

The AT<sub>n</sub> used for non-carcinogenic effects is the product of a 365-day year and the exposure duration (*i.e.*, AT<sub>n</sub> = 365 days × ED). Because the ED parameter changes for each receptor, the AT<sub>n</sub> changes as well. The AT<sub>n</sub> values used for each receptor are summarized below:

Future Construction Worker - 365 days  
Maintenance Worker - 9125 days  
Adolescent Visitor - 3650 days  
Off-Site Child Resident – 2,190 days  
Off-Site Adult Resident – 8,760 days

#### 4.2.1.4 Body Weight

The body weight used for the adult exposures (future construction worker and maintenance worker) analyzed in this assessment was the current US EPA default value of 70 kg (US EPA, 1989; US EPA Region 4, 1995). This value was also used for the adolescent/adult off-Site resident scenario. The adolescent body weight used for the visitor scenarios was 45 kg. This value was extracted from US EPA Region 4 guidance (1995). For the child resident scenario, a body weight of 15 kg was used as recommended by US EPA (1991).

#### 4.2.2 Route-Specific Exposure Parameters

The general intake equation discussed above (Equation 1) was modified by including route-specific exposure parameters in order to calculate route-specific intake values. For dermal exposures, skin surface area, adherence factor, exposure time (surface water exposures only), and absorption factor parameters were included in the intake equation. For ingestion exposures, an ingestion rate and a matrix effect were included in the intake calculation. For inhalation exposures, an inhalation rate and a retention factor for fugitive dusts were included in the intake equation. Also, for inhalation exposures, an additional paradigm was necessary to convert soil concentrations to concentrations in air available for intake.

#### 4.2.2.1 Dermal Exposure Parameters

##### Skin Surface Area

The total skin surface area used for adult receptors in this assessment was 20,000 cm<sup>2</sup>. This is a US EPA default value extracted from the *Exposure Factors Handbook* (1997). For adolescent exposures, a value of 12,768.3 cm<sup>2</sup> was used for total skin surface area. This was a mean value calculated based on the distributions of total skin surface areas for males and females between the ages of 7 and 16 as presented in *Exposure Factors Handbook* (1997). For the off-Site child resident scenario, a skin surface area of 7,213 cm<sup>2</sup> was used. This value was based on skin surface area data for male and female children provided in *Exposure Factors Handbook* (1997).

For purposes of exposure, it was assumed that only portions of the body would be exposed to the affected media on the Site. For the construction worker scenario, it was assumed that the hands, forearms, lower legs, and face would be exposed to Site soils. These body parts comprise 27.8% of the total skin surface area, or 5560 cm<sup>2</sup>.

For maintenance worker exposures to Site soils, it was assumed that the hands, forearms, and face would be exposed. These body parts comprise 15 percent of the total skin surface area, or 3000 cm<sup>2</sup>.

For surface water and sediment exposures, exposed body parts for construction and maintenance workers included hands, forearms, and face or 3000 cm<sup>2</sup> (15% of the total skin surface area).

The visitor and off-Site resident scenarios assumed that the hands, forearms, and lower legs would be exposed for contact with Site soils. These body parts comprise 23.9% of the total skin surface area, or 3052 cm<sup>2</sup> for adolescent visitors, 1724 cm<sup>2</sup> for child residents, and 4780 cm<sup>2</sup> for adult residents. For exposures to surface water and sediment, hands, forearms, lower legs, and feet were assumed exposed for adolescent visitor and off-Site resident scenarios. These body parts comprise 30.9 % of the total skin surface area or 3945 cm<sup>2</sup> for adolescent visitors, 2229 cm<sup>2</sup> for child residents, and 6180 cm<sup>2</sup> for adult residents.

### Soil Adherence Factor

Until recently, the US EPA-recommended default for soil adherence on skin ranged from 0.2 to 1.0 mg/cm<sup>2</sup> for the entire exposed surface area, without consideration of the type of activity (US EPA, 1992). However, the data from which that range was derived were primarily the result of indirect measurements, artificial activities, and sampling of hands only. A more recent study has presented the results of direct measurement of soil loading on skin surfaces before and after normal occupational and recreational activities that might result in soil contact (Kissel *et al.*, 1996). A five-order of magnitude range (roughly 10<sup>-3</sup> to 10<sup>+2</sup> mg/cm<sup>2</sup>) was reported for observed activity-related hand loadings. That report indicated that hand loadings within the range of 0.2 to 1.0 mg/cm<sup>2</sup> were produced by activities in which there was vigorous soil contact (*e.g.*, rugby, farming); but for activities in which there was less soil contact (*e.g.*, soccer, professional grounds maintenance), loadings substantially less than 0.2 mg/cm<sup>2</sup> were found on hands and other body parts. Kissel *et al.* (1996) concluded that, because non-hand loadings attributable to higher contact activities exceeded hand loadings resulting from lower contact activities, hand data from limited activities cannot be used as a conservative predictor of loadings that might occur on other body surfaces without regard to activity. Furthermore, because exposures are activity-dependent, dermal exposure to soil should be quantified using data describing human behavior (*e.g.*, type of activity, frequency, duration, including interval before bathing, clothing worn, etc.).

The most recent version of the *Exposure Factors Handbook* (1997) states:

In consideration, of these general observations and the recent data from Kissel *et al.* (1996, 1997), this document recommends a new approach for estimating soil adherence to skin. First use Table 6-12 [Summary of Field Studies, Kissel *et al.*, 1996a] to select the activity which best approximates the exposure scenario of concern. Next, use Table 6-13 [Mean Soil Adherence by Activity and Body Region, Kissel *et al.*, 1996a] to select soil loadings on exposed skin surfaces which correspond to the activity of interest. This table contains soil loading estimates for various body parts. The estimates were derived from soil adherence measurements of body parts of individuals engaged in specific activities described in Table 6-12. These results provide the best estimate of central loadings, but are

based on limited data. Therefore, they have a high degree of uncertainty such that considerable judgment must be used when selecting them for an assessment.

In another study that assessed the percentage of skin coverage in several soil contact trials in a greenhouse and an irrigation pipe laying trial, Kissel *et al.* (1996) concluded that adjusted loadings may be two to three orders of magnitude larger than average loadings if average loadings are small.

The activity-specific soil adherence factor for exposures to a maintenance worker was calculated based on data presented by Kissel *et al.* (1996) for grounds keepers, as presented below:

|                    |                         | Soil Adherence Factor by Body Part (mg/cm <sup>2</sup> ) |                |                 |               |
|--------------------|-------------------------|--|----------------|-----------------|---------------|
| Receptor           | Representative Activity | Hands  | Arms           | Lower Legs      | Face          |
| Maintenance Worker | Grounds Keepers         | 0.030 - 0.15   | 0.0021 - 0.023 | 0.0008 - 0.0012 | 0.0021 - 0.01 |

Data for the grounds keepers were used for the maintenance worker estimates because the activities of a grounds keeper best mimic those of a maintenance worker.

Soil adherence factors were calculated by normalizing each body part-specific soil adherence value (using the mid-points of the ranges tabulated above) with regard to the percentage of total body surface area represented by the respective body part (extracted from the US EPA *Dermal Exposure Assessment: Principles and Applications* [US EPA, 1992]). The maintenance worker adherence factor for soil was calculated based upon exposure to the hands, forearms and face. Surface area percentages for the hands, forearms, and face are 5.2, 5.9, and 3.9 percent, respectively (US EPA, 1997). Those body parts comprise 15 percent of the total body surface area. The normalized values for all body parts of interest were added, and the sum was divided by the total percentage of body surface area occupied by the parts. For example, the soil and sediment adherence factors for maintenance worker soil exposures (0.038 mg/cm<sup>2</sup>) were calculated as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.09 \times 0.052) + (0.0126 \times 0.059) + (0.006 \times 0.039)}{0.15} = 0.038$$

The construction worker adherence factor was also calculated in this fashion. This exposure scenario assumed that the hands, forearms, lower legs, and face would be exposed to Site soils. Soil loadings for the upper torso (chest and back) were not measured by Kissel *et al.* (1996) for construction workers because this body area is generally covered. However, to account for exposure to the upper torso during the very hot months of the year, the total area of the forearms, legs, hands, and face were assumed to be completely exposed. The hands, forearms, legs, and face comprise 5.2%, 5.9%, 12.8%, and 3.9% of the total skin surface area, respectively (with the face comprising one-third the surface area of the head), for a total of 27.8% exposed surface area. The construction worker soil adherence factor was based on data from Kissel *et al.* (1996) for construction workers as follows:

|                     |                         | Soil Adherence Factor by Body Part ( $\text{mg/cm}^2$ ) |       |            |       |
|---------------------|-------------------------|---|-------|------------|-------|
| Receptor            | Representative Activity | Hands   | Arms  | Lower Legs | Face  |
| Construction Worker | Construction Worker     | 0.24  | 0.098 | 0.066      | 0.029 |

The soil adherence factor for the construction worker scenario was calculated as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.24 \times 0.052) + (0.098 \times 0.059) + (0.066 \times 0.128) + (0.029 \times 0.039)}{0.278} = 0.1$$

For sediment exposures, the soil adherence factor was calculated for the construction worker scenario using adherence data from Kissel *et al.* (1996) for construction workers (as tabulated above) for the hands, forearms, and face. The hands, forearms, and face comprise 5.2, 5.9, and

3.9 percent of the total skin surface area, respectively (totaling 15 percent). Thus, the adherence factor for construction workers exposed to sediment ( $0.13 \text{ mg/cm}^2$ ) was calculated as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.24 \times 0.052) + (0.098 \times 0.059) + (0.029 \times 0.039)}{0.15} = 0.13$$

The adherence factor for visitor and off-Site resident exposures to soil assumed that the forearms, hands, and lower legs would be exposed to soil or sediment. The data used in these calculation were based on data by Kissel *et al.* (1996) for soccer players (exposed to a playing field of roughly one-half grass and one-half bare earth in a light mist) as presented below:

|                               |                         | Soil Adherence Factor by Body Part ( $\text{mg/cm}^2$ ) |              |                |
|-------------------------------|-------------------------|---|--------------|----------------|
| Receptor                      | Representative Activity | Arms  | Hands        | Lower Legs     |
| Visitor and Off-Site Resident | Soccer Players          | 0.0029 – 0.011  | 0.019 – 0.11 | 0.0081 – 0.031 |

The forearms, hands, and lower legs comprise 5.9%, 5.2%, and 12.8% of the total skin surface area, respectively, for a total of 23.9% (US EPA *Exposure Factors Handbook*, 1997). The adherence factor was then calculated for visitor and off-Site resident dermal exposures to soil as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.00695 \times 0.059) + (0.0645 \times 0.052) + (0.0196 \times 0.128)}{0.239} = 0.026$$

A value of  $0.026 \text{ mg/cm}^2$  was used as the soil adherence factor for visitors to the Site and off-Site residents.

Soil adherence factors for sediment exposures to Site visitors and off-Site residents were calculated using adherence data for the hands, forearms, lower legs, and feet. Adherence data for

reed gatherers were used for these exposures to best mimic activities that may incur sediment exposures. The reed gatherers studied by Kissel *et al.* (1996) periodically visited tidal flats to collect raw materials for basket weaving. The data from Kissel *et al.* (1996) presented in *Exposure Factors Handbook* (US EPA, 1997) were as follows:

|                                 |                         | Soil Adherence Factor by Body Part (mg/cm <sup>2</sup> ) |       |            |      |
|---------------------------------|-------------------------|--|-------|------------|------|
| Receptor                        | Representative Activity | Hands  | Arms  | Lower Legs | Feet |
| Visitors and Off-Site Residents | Reed Gatherers          | 0.66   | 0.036 | 0.128      | 0.63 |

The hands, forearms, lower legs, and feet comprises 5.2, 5.9, 12.8 and 7.0 percent of the total skin surface area, respectively (totaling 30.9 percent). Thus, the adherence factor for visitors and off-Site residents exposed to sediment (0.33 mg /cm<sup>2</sup>) was calculated as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.66 \times 0.052) + (0.036 \times 0.059) + (0.16 \times 0.128) + (0.63 \times 0.07)}{0.309} = 0.33$$

### Exposure Time

To estimate intakes as a result of dermal exposure to surface water, an exposure time (ET) parameter was included in the intake formula for Site visitors and off-Site residents. The parameter value of 1.0 hour/day was estimated using best professional judgement. This value represents the amount of time a Site visitor or off-Site resident may spend exposed to surface water in any one EU.

### Dermal Permeability Constant

The permeability constant, Kp, accounts for the movement of a constituent dissolved in water through the skin, across the stratum corneum, and into the blood stream. Kp values for the constituents examined in this assessment for surface water exposures were obtained from US EPA *Dermal Exposure Assessment: Principles and Applications* (1992). For values not available in

US EPA *Dermal Exposure Assessment* (1992), the K<sub>p</sub> value were calculated using the equations provided by the US EPA in the same document.

#### Dermal Absorption Factor

The final parameter included in the dermal intake paradigm was a dermal absorption factor. In general, the skin provides an effective barrier to environmental toxins. For example, certain hair-coloring formulations which are vigorously rubbed onto the scalp on a daily basis contain lead acetate at concentrations up to 200,000 ppm, yet lead toxicity does not appear to result. Moore *et al.* (1980) determined that the rate of lead absorption from <sup>203</sup>Pb labeled lead acetate in cosmetic preparations containing six mmol Pb acetate/L in male volunteers over 12 hours was 0.06% during normal use of such preparations. For most inorganic salts, percutaneous (skin) absorption is considered insignificant relative to incidental ingestion (for example, US EPA, 1986). On the other hand, some drugs (*e.g.*, nicotine) are effectively administered and absorbed into the blood stream from dermal "patches."

Most dermal bioavailability data for impacted soil have been obtained in laboratory animals or in vitro test systems. This introduces a significant source of uncertainty for predicting the human response. Safety factors have sometimes been applied to dermal absorption data obtained in animals to conservatively estimate the upper-bound of likely human percutaneous uptake of a certain constituent from skin exposure. This is usually unnecessary because human skin has generally been shown, for a diverse group of constituents, to be about 10-fold less permeable than the skin of typical animal species, such as rabbits and rats (Bartek and LaBudde, 1975; Shu *et al.*, 1988).

US EPA Region III evaluated available data concerning the dermal absorption of specific constituents and classes of constituents and provided several recommendations (US EPA Region 3, 1995). For semivolatile compounds, such as *bis*(2-ethylhexyl)phthalate, the US EPA recommends a range of 1% to 10% (US EPA, 1995). Kao *et al.* (1985) reported 2.7 percent for absorption of topically applied pure benzo(a)pyrene by human skin *in vitro*. The US EPA

Region 3 recommends using 10% as a conservative assumption based on the Ryan *et al.* study (1987). In addition, US EPA Region 4 guidance (1995) states that a soil dermal absorption factor “of 1.0% for organics and 0.1% for inorganics should be used as defaults in determining the uptake associated with dermal exposure” (see the Dermal Contact subsection of Exposure Assessment section of the 1995 guidance). For the purpose of this risk assessment, an ABS of 3% for cPAHs and of 10% for other SVOCs were conservatively assumed for dermal absorption, in keeping with US EPA Region 3’s and MDEQ’s recommendations.

#### 4.2.2.2 Ingestion Exposure Parameters

##### Ingestion Rate

US EPA’s *Exposure Factors Handbook* (1997) discusses three adult soil ingestion studies with results ranging from 10 mg/day to 480 mg/day. Hawley’s (1985) value of 480 mg/day (as recommended by the MDEQ) was “derived from assumptions about soil/dust levels on hands and mouthing behavior” (US EPA, 1997). Since no supporting measurements were made for Hawley’s study, the US EPA states that Hawley’s estimate “must be considered conjectural” (1997). As such, the US EPA goes on to suggest adult soil ingestion rates of 50 mg/day for industrial settings and 100 mg/day for residential and agricultural settings, although “50 mg/day still represents a reasonable central estimate of adult soil ingestion and is the recommended value...” (1997). Accordingly, a value of 100 mg/day for the maintenance worker and adult off-Site resident is amply conservative and was used in this assessment. In conjunction with the use of a two-tiered EF to reflect the different stages of potential future construction activities (see Section 4.2.1.1), the soil ingestion s for the construction worker scenario was also divided into two exposure levels for a single individual. A highly conservative ingestion rate of 480 mg/day (used in relevant exposure model calculations under “Exposure Level A”) was used for construction workers for the first 10 days of exposure to address direct contact with soil during earth-moving activities such as foundation excavating. A soil ingestion rate of 100 mg/day (used in relevant exposure model calculations under “Exposure Level B”) was used for the remainder of the construction worker exposure (70 days). Risks were then summed for both exposure levels to estimate the total potential risk posed to an individual construction worker

The ingestion rate used for the adolescent visitor scenario was 100 mg/day. The US EPA Region IV (1995) recommends a value of 200 mg/day as a mean ingestion rate for children under six years of age. This value was conservatively used in this assessment to estimate soil and sediment ingestion exposures for an off-Site resident child aged one to six years.

#### Gastrointestinal Matrix Effects of Soil

Incidental ingestion incorporates the matrix effect (ME; sometimes called the absorption adjustment factor [AAF]) into the general intake equation. When constituents are administered in solid vehicles such as food and soil, only a fraction of the ingested dose is extracted from the vehicle and subsequently absorbed through the gastrointestinal tract (US EPA *Estimated Exposure to Dioxin-like Compounds*, 1992). Gastrointestinal absorption of constituents sorbed onto such a medium is inhibited by physical-constituent bonding to the matrix (Hawley, 1985). This phenomenon is referred to as the gastrointestinal matrix effect of soil. Several studies referenced in the US EPA's *Estimated Exposure to Dioxin-like Compounds* (1992) have been performed to estimate the oral absorption factors of constituents from soil. At the request of MDEQ (2001), however, a gastrointestinal matrix effect of 1.0 was used in accordance with US EPA Region IV guidance (1995), although this approach is highly conservative and does not account for scientific studies that indicate the absorption of chemical constituents through the gastrointestinal tract is less than 100%.

#### 4.2.2.3 Inhalation Exposure Parameters and Paradigms

##### Inhalation Rate

The inhalation rate used for the construction worker scenario was 20 m<sup>3</sup>/day. This is a common US EPA default value and was recommended by US EPA Region 4 (1995).

##### Retention Factor

According to the International Commission on Radiological Protection (ICRP), 75 percent of respirable dust particles (PM<sub>10</sub>, or particles less than 10 microns in aerodynamic diameter) are retained when inhaled, the vast majority of which is potentially subsequently swallowed (ICRP,

1968). This 75% was included in the inhalation intake equation as the retention factor parameter (RF). This parameter applies only to non-VOC constituents entrained onto dust particles.

### Concentration in Air

To estimate airborne dust levels during hypothetical construction activities, an emission rate of suspendible particles of less than 15 microns in aerodynamic diameter ( $PM_{15}$ ) was calculated (grams/second); particles less than 10 microns were considered to be respirable. Considering particles of 15 microns or less in diameter in the emission rate calculation is a conservative assumption, inasmuch as only particles with an aerodynamic diameter of less than five to seven microns are inhaled into the lung.

The two types of construction activities at the Site that have the potential to emit fugitive dusts are vehicular movement over bare (unpaved or unvegetated) surfaces and the excavation of soil. Estimation of fugitive dust emissions caused by each activity were examined separately, as follows, and were derived from existing estimates of general construction exposure. The sum of the emissions from these two activities was multiplied by the concentration of constituent in the soil ( $C_s$ ) in order to derive the total emission rate ( $E_i$ ) for non-VOCs as follows:

$$E_i = C_s \times (PERv + PERe) \quad [Equation 2]$$

where:

- $E_i$  = Emission rate (mg/sec);  
 $C_s$  = Concentration in soil (mg/kg);  
 $PERv$  = Particulate emission rate for vehicular movement (lb/vehicle mile);  
and  
 $PERe$  = Particulate emission rate for excavation (lb/vehicle mile).

The following empirical expression (US EPA, 1988) was used to estimate the fugitive dust generated by vehicles during construction activities:

$$PERv (\text{lbs/vehicle mile}) = k \times 5.9 \times (s/12)(S/30) \times (mvw/3)^{0.7} \times (ww/4)^{0.5} \times ((365 - p)/365)$$

[Equation 3]

where:

- PERv = Vehicle particle emission rate (lb/vehicle mile traveled);  
s = Percent silt content (unitless);  
k = Particle size multiplier (unitless);  
S = Mean vehicle speed (mph);  
mvw = Mean vehicle weight (ton);  
ww = Mean number of wheels per vehicle (unitless); and  
p = Mean number of days with  $\geq$  0.01 inches of precipitation per year (unitless).

It was assumed that the vehicle travels during 40% of the 80-day exposure duration and 0.5 miles per day. The result is a value of 16 miles per construction event. Percent silt content was estimated to have a mean value of 50%, based on geotechnical data provided in the *Remedial Investigation Report* (Pisani & Assoc., 1997). US EPA default values were utilized and referenced for all other parameters. The particle size multiplier was assumed to be 0.50, corresponding to particles less than 15 microns (US EPA, 1996). Vehicle characteristics consist of the following: mean vehicle speed was assumed to be 15 mph, with mean vehicle weight assumed to be approximately 12.5 tons, for 8-wheeled vehicles (US EPA, 1988). The estimated mean number of days with precipitation equal to or greater than 0.01 inches per year is 110 (US EPA, 1988). Total resultant dust emissions for constituents during vehicular movement activities were estimated to be approximately 16.5 lbs/vehicle mile traveled, or 0.0001 kg/sec. Calculations are summarized in Table 20.

Future excavation may be performed by bulldozers, a backhoe, or other heavy construction equipment. The following estimate of particulate emissions, less than 15  $\mu\text{m}$  in diameter resulting from bulldozing activity, was based on the approach described in the US EPA *Compilation of Air Pollution Emission Factors* (1996), as developed from studies of emissions from uncontrolled open dust sources resulting from bulldozing at western surface coal mines.

$$PERe \text{ (lb/hour)} = \frac{1.0 \times s^{1.5}}{M^{1.4}} \quad [\text{Equation 4}]$$

where:

PERe = Excavation particle emission rate (lb/hr);  
s = Percent silt content (unitless); and  
M = Soil moisture content (unitless).

Percent soil moisture content was assumed to be 15.1%, an average of Site-specific soil moisture data and percent silt content 50%, as described above.

The resultant fugitive dust emission rate during excavation activities was 7.9 lbs/hr or 0.001 kg/sec. Table 20 summarizes these calculations.

Once the emission rate (Ei in Equation 2) was calculated, it was converted to a concentration in ambient air. Gaussian models are conventionally used to determine downwind ambient air concentrations, Ca, from the emission rate, Ei, estimated. However, in this scenario, such models have limited applicability when the receptor(s) is at or very near the source of emission. In this case, a bulldozer operator, for example, is situated directly within the area of ground emissions of vapors and dusts. Average ambient air concentrations in this circumstance are best estimated by use of a near-field box model (US EPA, 1988).

The near-field box model assumes uniform wind speed and uniform mixing throughout the box. The release and mixing of VOCs or respirable dusts in ambient air is estimated as follows:

$$Ca \text{ (mg/m}^3\text{)} = \frac{Ei}{W_b \times H_b \times V} \quad [\text{Equation 5}]$$

where:

|                |   |  |
|----------------|---|--|
| Ca             | = | Concentration of constituent in ambient air (mg/m <sup>3</sup> );                        |
| Ei             | = | Emission rate of constituent (mg/sec);   |
| W <sub>b</sub> | = | Width of box in crosswind dimension within the area of residual constituent in soil (m); |
| H <sub>b</sub> | = | Downwind height of box (m); and  |
| V              | = | Average wind speed through the box (m/sec).  |

The value of H<sub>b</sub> in this calculation is determined by the downwind distance and the atmospheric turbulence at ground level, which determines the trajectory of a release from the upwind edge of the source of vapor or dust emissions. For neutral atmospheric conditions, the height at the downwind boundary (H<sub>b</sub>) may be expressed by the following function (Pasquill 1975, Horst 1979):

$$z = 6.25 r [H_b/r \times \ln(H_b/r) - 1.58 H_b/r + 1.58] \quad [\text{Equation 6}]$$

where:

|                |   |  |
|----------------|---|--|
| H <sub>b</sub> | = | Downwind height of box (m);              |
| z              | = | Downwind distance to boundary (m); and   |
| r              | = | A terrain-dependent roughness height (m) |

H<sub>b</sub> (defined in Equation 5) is adjusted until the z parameter is equal to W<sub>b</sub> (defined in Equation 5). The resulting H<sub>b</sub> value is the height of the box. On any given workday, it is estimated that grading or excavation activities occur over the entire "workable" Site area (exposure unit) from which dusts are generated. This area is estimated to be 2,500 m<sup>2</sup>, with length of the box estimated to be 50 meters (downwind distance) and the width of the box (W) estimated to be 50 meters. The greater the roughness height, the greater the wind turbulence and constituent dilution (*i.e.*, the height of the box increases). For the purposes of this risk assessment, it is conservatively assumed that the roughness height is 0.20 meters, which corresponds to a terrain with grass, some small bushes, and occasional trees (US EPA *Rapid Assessment of Exposure to Particulate Emission from Surface Contamination Sites*, 1985). This

assumption is appropriate for the actual Site conditions. An annual average wind speed (4.69 m/sec) is obtained from the STAR data set, accessed through the Personal Computer Graphical Exposure Modeling System (PCGEMS), for STAR station 03940, Jackson/Thompson, MS for the period 1974-1978 (Table 21).

## **5.0      Toxicity Assessment**

The toxicity assessment involves the evaluation of available toxicity information to be utilized in the risk assessment process. Toxicity values derived from a dose-response relationship can be used to estimate the potential for the occurrence of adverse effects in individuals exposed to various constituent levels.

Exposure to a constituent does not necessarily result in adverse effects. The relationship between dose and response defines the quantitative indices of toxicity required to evaluate the potential health risks associated with a given level of exposure. If the nature of the dose-response relationship is such that no effects can be demonstrated below a certain level of exposure, a threshold can be defined and an acceptable exposure level derived. Humans are routinely exposed to naturally-occurring constituents and man-made constituents through the typical diet, air, and water, with no apparent adverse effects. However, the potential for adverse effects may occur if the exposure level exceeds the threshold in a variably sensitive population. This threshold applies primarily to constituents which produce non-carcinogenic (systemic) effects, although there is a growing body of scientific evidence which suggests that exposure thresholds may exist for certain carcinogenic constituents as well.

Adverse effects can be caused by acute exposure, which is a single or short-term exposure to a toxic substance, or by chronic exposure on a continuous or repeated basis over an extended period of time. "Acceptable" acute or chronic levels of exposure are considered to be without any anticipated adverse effects. Such exposure levels are commonly expressed as reference doses (RfDs), health advisories, etc. An acceptable exposure level is calculated to provide an "adequate margin of safety."

Chronic RfDs, which have been derived by the US EPA for a large number of constituents, were utilized to evaluate exposures lasting seven to 70 years (US EPA, 1989). Activities involving exposures of shorter duration to COPCs at the Site are anticipated to result in hazard and risk estimates that are lower than those associated with the long-term exposures. Identification of

subchronic toxicity values corresponding to shorter-term exposure scenarios (*i.e.*, less than seven years) are included in the risk assessment to ensure that both short-term and long-term risks can be addressed.

Currently, the US EPA has not developed toxicity values to be utilized in dermal exposure scenarios; however, the US EPA does provide the following guidance for dermal exposure:

No RfDs or slope factors are available for the dermal route of exposure. In some cases, however, non-carcinogenic or carcinogenic risks associated with dermal exposure can be evaluated using an oral RfD or oral slope factor, respectively. (US EPA, 1989).

Provisional dermal toxicity values were developed and utilized in the dermal exposure pathways considered in the human health risk assessment to provide a more accurate Site-specific risk assessment. These dermal RfD values were developed by multiplying the published oral RfD for a given constituent by the fraction of that constituent that can be absorbed through the gastrointestinal tract (stomach/intestine lining). The absorption fraction utilized was 50% for semivolatiles as extracted from US EPA Region 4 guidance (1995).

A number of sources of toxicity information exists, and these sources vary with regard to the availability and strength of supporting evidence. The following protocol has been established for the determination of toxicity indices; it defines a hierarchy of sources to be consulted and the methodology for the determination of toxicity values. This protocol has been developed in accordance with current US EPA methodology. Toxicity values for the COPCs at the Site were obtained with reference to the following hierarchy of sources developed in accordance with MCEQ guidance (1999):

- 1) Toxicity values were obtained from the *Integrated Risk Information System* (IRIS, 1999) database. This database contains the RfDs and Cancer Slope Factors (CSFs), which have been verified by the US EPA's RfD and Carcinogen Risk Assessment Verification Endeavor (CRAVE) workgroups, and is, thus, the

agency's preferred source for toxicity values. IRIS supersedes all other information sources.

- 2) For toxicity values which are unavailable on IRIS, the most current source of information is the Health Effects Assessment Summary Tables (HEAST, US EPA, 1997), published by the US EPA. HEAST contains interim, as well as verified RfDs and CSFs. Supporting toxicity information for verified values is provided in an extensive reference section of HEAST.
- 3) In cases where IRIS or HEAST could not provide toxicity values, US EPA Region III's Risk-Based Concentration (RBC) Tables were visited. These tables often provide toxicity values generated by reliable sources other than IRIS or HEAST. For example, in response to specific requests from risk assessors, the US EPA National Center for Environmental Assessment (NCEA) develops provisional RfDs or CSFs for chemicals not listed in IRIS or HEAST. Region III's RBC tables will list such provisional values. Also, RfDs or CSFs that have since been withdrawn from IRIS or HEAST may still be listed on the Region III RBC tables, although they are flagged with a "W." These toxicity values were no longer agreed upon by US EPA scientists; however, the Region III RBC tables continue to publish such values because risk assessors still need to quantify exposures to these chemicals. Lastly, the Region III RBC tables will list toxicity indices found in "other" US EPA documents. These values are flagged with an "O" on the tables.

The US EPA has derived carcinogenic slope factors for both oral and inhalation pathways, and these are utilized to quantitatively estimate risks. In the first step of the US EPA's evaluation, the available data are analyzed to determine the likelihood that the agent is a human carcinogen. The evidence is characterized separately for human studies and animal studies as sufficient, limited, inadequate, no data, or evidence of no effect. The characterizations of these two types of data are combined, and based on the extent to which the agent has been shown to be a carcinogen in experimental animals or humans, or both, the agent is given a provisional weight-of-evidence classification. The US EPA scientists then adjust the provisional classification upward or downward, based on other supporting evidence of carcinogenicity (see Section 7.1.3, US EPA, 1989). For a further description of the role of supporting evidence, see the US EPA guidelines (US EPA, 1986).

The US EPA classification system for weight of evidence is shown in the table below. This system is adapted from the approach taken by the International Agency for Research on Cancer.

| <b>US EPA WEIGHT-OF-EVIDENCE<br/>CLASSIFICATION SYSTEM FOR<br/>CARCINOGENICITY</b> |  |
|--|--|
| <b>Group</b>   | <b>Description</b>   |
| A  | Human carcinogen   |
| B1 or<br>B2  | Probable human carcinogen<br><br>B1 indicates that limited human data are available<br><br>B2 indicates sufficient evidence in animals and inadequate or no evidence in humans |
| C  | Possible human carcinogen  |
| D  | Not classifiable as to human carcinogenicity   |
| E  | Evidence of non-carcinogenicity for humans   |

(US EPA, 1989)

Table 22 summarizes the available toxicity values for the identified COPCs. COPCs lacking published toxicity values were not able to be quantitatively evaluated in this assessment in accordance with MCEQ guidance (1999). The MCEQ limits the use of toxicity values to those that have been published in IRIS, HEAST, ATSDR toxicity profiles, or other peer-reviewed reference sources or literature approved by the MCEQ (1999). The MDEQ (2001), however, requested that risks from dermal exposure to cPAHs be estimated using the oral cancer slope factor for benzo(a)pyrene, applying benzo(a)pyrene relative potency factors, and accounting for an absorption efficiency of 50%. This methodology was used accordingly.

## **6.0 Risk Characterization**

The objective of the risk characterization is to determine potential risk to receptors by combining the results of the exposure and toxicity assessments. Non-carcinogenic effects and carcinogenic risks are summarized in Table 23. Tables 24 through 78 provide algorithms and parameters for each pathway.

The estimated intakes calculated for each exposure pathway considered and each COPC were compared to RfDs for non-carcinogenic effects. The following formula was used to estimate the potential for non-carcinogenic health effects for each COPC.

$$HQ = ADI/RfD \quad [Equation 7]$$

where:

HQ = Hazard quotient - potential for noncancer health effects (unitless);  
ADI = Average daily intake of COPC (mg/kg-day); and  
RfD = Reference dose (mg/kg-day).

RfDs have been developed by the US EPA for chronic (*e.g.*, lifetime) and/or subchronic exposure to constituents based on the most sensitive non-carcinogenic effects. The chronic RfD for a constituent is an estimate of a lifetime daily exposure level for the human population, including sensitive subpopulations, that is likely to be without an appreciable risk of deleterious effects. The potential for noncancer health effects was evaluated by comparing the Site-specific exposure level with the RfD derived by the US EPA for a similar exposure period. This ratio of exposure to toxicity is called the hazard quotient (HQ). If the Site-specific exposure level exceeds the threshold (*i.e.*, the HQ exceeds a value greater than 1.0), there may be concern for potential noncancer effects.

To assess the overall potential for noncancer effects posed by multiple constituents, a hazard index (HI) is derived by summing the individual HQs. This approach assumes additivity of critical effects of multiple constituents. This is appropriate only for compounds that induce the

same effect by the same mechanism of action. This conservative approach significantly overestimates the actual potential for adverse health impacts.

In cancer risk assessment, the US EPA has required the use of the upper limit which produces an estimate of potential risk that has a 95% probability of exceeding the actual risk, which may, in fact, be zero. The following formula was utilized to estimate the upper bound excess cancer risk for each carcinogen (note that not all COPCs are carcinogens):

$$TR = CLDI \times SF \quad [Equation 8]$$

where:

- TR = Target risk - excess probability of an individual developing cancer (unitless);  
CLDI = Calculated lifetime average daily intake of carcinogenic COPC (mg/kg-day); and  
SF = Cancer slope factor (mg/kg-day)<sup>-1</sup>.

For exposures to multiple carcinogens, the upper limits of cancer risks are summed to derive a total cancer risk. The US EPA recognizes that it is not technically appropriate to sum upper confidence limits of the risk to produce a realistic total probability, but requires this approach be used.

Carcinogenic risk refers to the probability of developing cancer as a result of exposure to known or suspected carcinogens. The National Contingency Plan (NCP) endorses an acceptable risk range of  $10^{-4}$  to  $10^{-6}$  for exposure to multiple carcinogens. This range represents an incremental increase of 1 in 10,000 to 1 in 1,000,000 in the chance of developing cancer over a lifetime. The MCEQ (1999) indicates that the target risk level is  $1 \times 10^{-6}$  per individual carcinogen and an acceptable cumulative risk level is  $1 \times 10^{-4}$ . As such, risk levels totaled across oral, dermal, and inhalation pathways may exceed  $1 \times 10^{-6}$  and still be in compliance with MCEQ requirements (1999) as long as no single carcinogen exceeds  $1 \times 10^{-6}$  and the cumulative risk for a single receptor does not exceed  $1 \times 10^{-4}$ .

Table 23 provides a summary of the non-carcinogenic effects and carcinogenic risks associated with each of the pathways evaluated in this assessment.

The overall hazard index across the assessed pathways and EUs was 0.1 for the Site visitor scenario. This value is below the acceptable benchmark of 1.0. The highest hazard index associated with the Site visitor scenario was 0.07 corresponding to dermal exposure to sediment in EU4. The overall cancer risk for exposures to Site visitors was estimated to be  $2 \times 10^{-5}$  and is primarily attributable to oral and dermal exposure to benzo(a)pyrene and associated cPAHs in EU4 sediments. Oral exposure to the same constituents in EU4 and EU5 surface soils also contributed to the cancer risk estimate for the site visitor. Additional discussion regarding remediation goals for this scenario has been provided in section 8.0.

The overall hazard index for the maintenance worker scenarios was 0.1 and is below the acceptable benchmark of 1.0. The highest hazard index associated with the maintenance worker scenario was 0.1 corresponding to oral exposure to surface soil in EU4. The overall cancer risk for the maintenance worker scenario was  $1 \times 10^{-3}$  and was primarily attributable to dermal and oral exposure to benzo(a)pyrene and other cPAHs in surface soils in EUs 2, 4, and 5. Additional discussion regarding remediation goals for this scenario has been provided in section 8.0.

The overall hazard index for the hypothetical future construction worker was 0.003 and is well below the acceptable benchmark of 1.0. The highest hazard index associated with the construction worker scenario was 0.003 corresponding to oral exposure to soils in EU 5. The overall cancer risk for the hypothetical future construction worker scenario was  $5 \times 10^{-5}$  and is attributable to benzo(a)pyrene and associated cPAH oral exposure in EU4 sediment and oral and dermal exposure to EU4 and EU5 soils. Additional discussion regarding remediation goals for this scenario has been provided in section 8.0.

The off-Site resident scenario revealed a hazard index of  $1 \times 10^{-4}$ . This value is considerably below the acceptable benchmark of 1.0. The overall cancer risk for the resident exposure

scenario was estimated to be  $2 \times 10^{-4}$  and is attributable to oral and dermal exposure to benzo(a)pyrene and associated cPAHs in EU6.

## **7.0 Uncertainty Analysis**

Risk assessment uses a wide array of information sources and techniques. Even in those rare circumstances where constituent intake for an exposed individual may be measured relatively precisely, assumptions will still be required to evaluate the associated risk. Generally, data are not available for critical aspects of the risk assessment, and the use of professional judgment, inferences based on analogy, the use of default values, model estimation techniques, etc., result in uncertainty of varying degrees.

The expressions of risk in this assessment are not probabilistic; the expressions of risk are conditional, based on the conditions represented by the single-point values selected for the analysis. This section is intended to identify and qualitatively evaluate the more salient Site-specific uncertainties and their potential influence on the credibility of the estimated Site risks.

### **7.1 Uncertainty of Data Evaluation Factors**

Uncertainties in data analysis include analytical error, selection of COPCs, adequacy of sampling design, etc. Generally, there is far less uncertainty in this phase of the risk assessment process than other aspects contribute.

Laboratory analysis is extremely accurate relative to the potential error of "professional judgment" in exposure assessments. The uncertainty of analytical data is likely to be less than 25 percent, most of the time.

The adequacy of the sampling strategies to characterize Site conditions is a potentially large source of uncertainty. Because of the limited availability of resources, sample collection is generally limited. However, sampling (especially in multiple surveys) is not random, but is designed to locate the areas with the highest levels of constituents. Thus, test data are biased toward overestimation of average constituent levels. In addition, in most instances, the upper 95-percent confidence limit of the average concentration is utilized as an exposure-point

concentration in the risk assessment. The use of this value likely will result in an overestimation of risk, as the 95% UCL represents a value that will be greater than the true average 95% of the time.

Oftentimes, only a portion of detected constituents are carried through the risk assessment process because constituents are eliminated through COPC screening procedures (US EPA, 1989). This could result in an underestimation of risk, although the COPC selection process is intended to identify those constituents that account for the vast majority of potential risk. COPCs lacking published RfD values were not quantitatively evaluated and this may result in an underestimation of potential hazards (non-carcinogenic effects).

## 7.2     Uncertainty of Toxicity Values

The US EPA's IRIS states that the uncertainty associated with RfD values for non-carcinogenic endpoints of toxicity "span perhaps an order of magnitude." In fact, the uncertainty of extrapolating dose-response data from animals to humans with the application of multiple safety factors (100 to 10,000 or more) is likely to be several orders of magnitude. Current policies for deriving RfD values will often result in an overestimation of risk.

The uncertainty associated with the estimation of cancer risk contributes, by far, the major source of potential error and uncertainty. It is beyond the scope of this analysis to explore this toxicity assessment factor in any detail. However, a few salient points are noted below.

Some constituents classified as carcinogens have been shown to produce an increased incidence of cancer in mice but not rats, for example. If the mouse is not an adequate model for the rat, it may be wondered how reliable a model it is for human beings. The assumption of linearity and a non-threshold phenomenon in the dose versus risk relationship may not be valid and could result in a very large overestimation of actual cancer risk, if any even exist at low doses in humans.

The US EPA evaluated the uncertainty of cancer risk estimates from exposures to trichloroethene and several other related VOCs in public drinking water supplies (Cothorn *et al.*, 1984). These US EPA scientists concluded the following:

- The largest uncertainty in the calculations is due to the choice of the model [Multistage, Weibull, Logit, Probit, etc.] used in extrapolating risk to low doses in humans, and is 5 to 6 orders of magnitude;
- If a single model were chosen [assumed to be valid], the overall uncertainty in risk estimates would be 2 to 3 orders of magnitude;
- The exposure estimates contribute, at most, an order of magnitude to the uncertainty; and
- It would appear that until a particular compound's mechanisms of cancer are better known, it is likely that the uncertainty in the toxicity will not be improved.

### 7.3     Uncertainties in Assessing Potential Exposure

Ideally, Site-specific exposure values should be used when assessing potential intakes of chemicals at a Site. Oftentimes, however, Site-specific data are not available; therefore, the risk assessor must estimate values that most accurately reflect Site conditions. In doing so, US EPA or other regulatory default values were utilized in place of Site-specific data. These values may over- or under-estimate risks, depending on Site conditions and the percentile range in which the default values fall (*e.g.*, 50<sup>th</sup>, 95<sup>th</sup>).

Although a considerable amount of published data is available on the most common exposure parameters (*e.g.*, body weight, skin surface area), even these data contain uncertainties. Studies conducted by different scientists often provide differing levels of detail, statistics, and accuracy based on sample size, study design, geographic area, etc. Such discrepancies can increase uncertainty when the data are combined to derive a single-point default value. These data may be the best available; however, the reflection of reality may still be imprecise.

Where published exposure parameters were not available, best professional judgment had to be used, thereby increasing uncertainty. The default or estimated exposure parameters used in this assessment likely resulted in a moderate over-estimation of risk.

The intakes estimated for dermal absorption of PAHs adsorbed into soils adhering to skin may overestimate risks for a host of reasons. Early studies conducted by Falk and coworkers indicated that the carcinogenic effect of B(a)P on subcutaneous injection in mice could be markedly inhibited by the simultaneous administration of various non-carcinogenic PAHs (Falk *et al.*, 1964, as cited in ATSDR, 1988. In other subcutaneous injection and skin-painting studies with mice, it was shown that a combination of several non-carcinogenic PAH compounds, mixed according to the proportion occurring in auto exhaust, did not enhance or inhibit the action of two potent PAH carcinogens, B(a)P and dibenz(a,h)anthracene- (ATSDR, 1988).

The carcinogenic potency of B(a)P and other carcinogenic PAHs is generally determined by injecting solutions under the skin, painting the skin with the carcinogenic PAH dissolved in a solvent, or dissolved in corn oil in feeding studies. This vehicle or matrix affords a high level of bioavailability of the carcinogenic PAH compound. Recently, Krueger *et al.* (1999) conducted *in vitro* percutaneous absorption studies with contaminated soils and organic solvent extracts of contaminated soils collected at former manufactured gas plant (MGP) sites. The MGP tar-contaminated soils contained PAHs at levels ranging from 10 to 2400 mg/kg. The dermal penetration rates of PAH from the MGP tar-contaminated soils and soil solvent extracts were determined experimentally through human skin using tritium-labelled B(a)P as a surrogate. Results showed reductions of two to three orders of magnitude in PAH absorption through human skin from the most contaminated soils in comparison to the soil extracts. Reduction in PAH penetration was attributed to soil matrix properties. That is, PAH compounds adsorbed to organic carbon in a soil matrix are far less bioavailable for dermal flux than PAH compounds dissolved in a solvent. [No correction for such a profound soil matrix effect was applied in quantitatively estimating cancer risks due to dermal absorption of B(a)P and other carcinogenic PAHs in this assessment.]

## **8.0 Summary of Findings**

The results of the baseline human health risk assessment indicate potentially unacceptable risk levels for the following exposure scenarios:

| Potentially Exposed Population | Media           | EU      |
|--------------------------------|-----------------|---------|
| Site Visitor                   | Sediment        | 4       |
|                                | Surface Soil    | 4, 5    |
| Maintenance Worker             | Sediment        | 4       |
|                                | Surface Soil    | 2, 4, 5 |
| Construction Worker            | Sediment        | 4       |
|                                | Subsurface Soil | 4, 5    |
| Off-Site Resident              | Sediment        | 6       |

The risk levels associated with the above scenarios were driven by cPAHs, particularly benzo(a)pyrene. To determine the extent of remediation necessary to reduce these risks to acceptable levels, sediment and soil data for cPAHs in EUs 2, 4, 5, and 6 were closely examined.

The benzo(a)pyrene exposure-point concentration used to evaluate maintenance worker exposures to surface soil in EU2 was 5.2 mg/kg (sample location GEO-13/0-1'). This was the maximum benzo(a)pyrene concentration found in surface soil in EU2. The next highest concentration of benzo(a)pyrene in sediment was found at SS-10 (2.4 mg/kg). However, as previously noted, these samples were collected at locations within a densely wooded area. No remediation is planned to address surface soils at these locations for the following reasons:

- No maintenance activities are currently conducted in this area;
- Any remediation would require significant clearing; and
- Cancer risks associated with surface soils at these locations only slightly exceed  $1 \times 10^{-6}$  for two individual constituents, and the total cancer risk level is still less than  $1 \times 10^{-5}$ .

In EU4, the maximum concentration of benzo(a)pyrene was used as the exposure-point concentration for site visitor, maintenance worker, and construction worker exposure to sediment. The benzo(a)pyrene exposure-point concentration used to evaluate these in EU4 was 130 mg/kg (sample location SD-02, see Figure 2). The next highest concentration of benzo(a)pyrene in sediment was found at SD-12 (71 mg/kg). Implementing a remedy to remove, treat, or preclude contact with sediment at sample locations SD-02 and SD-12 would leave a concentration of 5.57 mg/kg (sample location SD-23) as the maximum concentration in sediment that could be potentially contacted by site visitors, maintenance workers, and/or construction workers in EU 4. Excluding samples SD-02 and SD-12 and using 5.57 mg/kg as the exposure-point concentration drops the risk level for dermal and oral contact with sediment by a visitor and oral contact with sediment by a maintenance worker or construction worker to within acceptable levels (i.e., no risk level associated with a single carcinogen exceeds  $1 \times 10^{-6}$ ; Tables 79 - 82).

In EU4, the maximum concentration of benzo(a)pyrene was also used as the exposure-point concentration for site visitor, maintenance worker, and construction worker soil exposures. Each of these receptors could potentially be exposed to soils at different depth ranges: visitor 0-1' bgs, maintenance worker 0-6' bgs, and construction worker 0-20' bgs. The sample locations and corresponding concentrations of benzo(a)pyrene that contributed to elevated risk estimates in the three exposure scenarios are presented in the table below:

| Sample Location | Benzo(a)pyrene Concentration<br>(mg/kg) |
|-----------------|---|
| GEO-48/0-1'     | 500                                     |
| GEO-21/0-1'     | 230                                     |
| GEO-21/2-3'     | 190                                     |
| GEO-19/0-1'     | 56                                      |
| GEO-46/0-1'     | 16                                      |

| Sample Location | Benzo(a)pyrene Concentration<br>(mg/kg) |
|-----------------|---|
| GEO-20/5-6'     | 11                                      |
| GEO-47/5-6'     | 9.6                                     |
| GEO-48/2-3'     | 6.1                                     |
| GEO-20/0-1'     | 3.2                                     |
| GEO-47/0-1'     | 3                                       |
| GEO-19/2-3'     | 2.4                                     |

Implementing a remedy to remove, treat, or preclude contact with the soils tabulated above would leave a maximum benzo(a)pyrene concentration of 0.29 mg/kg (sample location GEO-19/5-6'). Using the concentration of 0.29 mg/kg as the exposure-point concentration for estimating risk to site visitors, maintenance workers, and construction workers drops the risk levels to within acceptable levels (Tables 83 - 87). In situ biological treatment is proposed to address impacted soils within EU4. This will include clearing, tilling, application of inorganic nutrients, and, once soils are remediated to the extent practicable, placement of concrete cover. The area to be remediated will extend at least from Courtesy Ford to the edge of the railroad right-of-way, and may extend onto the railroad right-of-way with the permission of the Southern railway.

In EU5, the surface soil sample locations contributing most to elevated risk levels for the maintenance worker, construction worker, and site visitor scenarios were GEO-33/0-1', GEO-33/2-3', GEO-30/0-1', GEO-59/0-1, GEO-29/0-1', and GEO-28/0-1' (see Figure 2). All sample locations, with the exception of GEO-59/0-1', are located underneath paved areas in a parcel of land extending from Courtesy Ford to the southeast (Figure 2). Pavement in this area precludes direct contact with surface and subsurface soils; therefore, it is not anticipated that current or future maintenance workers or site visitors will have access to soils in or around these sample locations. In addition, a deed restriction will be implemented requiring the maintenance of the paved areas to ensure protection of human health in the future. Sample location GEO-59/0-1',

with a benzo(a)pyrene exposure point concentration is 6.1 mg/kg, however, is adjacent to West Pine Street in an unpaved area. Implementing a remedy to remove, treat, or preclude contact with surface soil at this location would leave a concentration of 0.37 mg/kg (GEO-60/0-1') as the maximum concentration in surface soil not covered by pavement that could potentially be contacted by any of the three receptors in this EU. Excluding sample GEO-59/0-1' and using 0.37 mg/kg as the exposure-point concentration drops the estimated exposures in EU5 to within acceptable levels (i.e., no risk level associated with a single carcinogen exceeds  $1 \times 10^{-6}$ ; Tables 88 - 91).

The benzo(a)pyrene exposure-point concentration used to evaluate adult and child resident exposures to sediment in EU6 was 49 mg/kg (sample location SD-03, see Figure 3). This was the maximum benzo(a)pyrene concentration found in sediments in EU6. Sample locations SD-04, SD-14, SD-13, SD-16, SD-15, and SD-17 (33, 12.2, 3.27, 2.8, 2.42, and 2.26 mg/kg, respectively) also contributed to elevated cancer risk estimates for both receptors. Implementing a remedy to remove, treat, or preclude contact with sediment at these sample locations would leave a concentration of 0.97 mg/kg (sample location SD-05). Using the benzo(a)pyrene concentration of 0.97 mg/kg as the exposure-point concentration for sediment exposure to adult and child residents reduces the risk estimate to within acceptable limits (i.e., no risk level associated with a single carcinogen exceeds  $1 \times 10^{-6}$ ; Tables 92 - 95). Remediation activities are proposed to remove impacted sediment and preclude contact with residuals in the northeast drainage ditch. These activities include removal and off-Site treatment and/or disposal of impacted sediments, installation of a storm water collection and conveyance pipe, backfilling around the culvert, and planting with native grass.

## Bibliography

- Andelman, J. B., and M. J. Suess. 1980. Polynuclear aromatic hydrocarbons in the water environment. Bull. WHO 43:479-508.
- Alexander, M. Aging, Bioavailability, and Overestimation of Risk from Environmental Pollutants. Environ. Sci. Technol, 2000, 34(20):4259.
- ATSDR (Agency for Toxic Substances and Disease Registry). Toxicological Profile for Benzo(a)Pyrene. Oak Ridge National Laboratory. 1988.
- ATSDR (Agency for Toxic Substances and Disease Registry). ATSDR's Toxicological Profiles on CD-ROM. Polycyclic Aromatic Hydrocarbons (PAHs), Update. CRC Press, 1999.
- Audere, A. K., Z. Y. Lindberg, G. A. Smirnov, and L. M. Shabad. 1973. Experiment in studying the influence of an airport located within the limits of a city on the level of environmental pollution by benzo(a)pyrene. Gig. Sanit. 38(9): 90-92.
- Bartek, M.J. and J.A. LaBudde. Percutaneous Absorption *in vitro*, in Animal Models in Dermatology. Ed. H.I. Maibach. New York: Churchill Livingstone, 1975. p. 103.
- Blumer, M. 1961. Benzopyrenes in soil. Science 134, 474-475.
- Blumer, M., W. Blumer, and T. Relch. 1977. Polycyclic aromatic hydrocarbons in soils of a mountain valley; correlation with highway traffic and cancer incidence. Environ. Sci. Technol. 11 (12), 1082-1084.
- Butler, J. D., V. Butterworth, C. Kellow, and H. G. Robinson. 1984. Some observations on the polycyclic aromatic hydrocarbon (PAH) content of surface soils in urban areas. Sci. Total Environ. 38, 75-85.
- Chu, M. M. L. and G. W. Chem. 1984. Evaluation and estimation of potential carcinogenic risks of polynuclear aromatic hydrocarbons. Paper presented at the Pacific Rim Risk Conference.
- Chung, N. and Alexander, M. Differences in Sequestration and Bioavailability of Organic Compounds Aged in Dissimilar Soils. Environ Sci. Technol. 32: 855.
- Coomes, R. M. 1981. Carcinogenic testing of oil shale materials. Twelfth Oil Shale Symposium Proceedings. Colorado School of Mines Pres.

Cothorn, C. R., W. Conniglio, W. Marcus. Techniques for the Assessment of the Carcinogenic Risk to the US Population due to Exposure from Selected Volatile Organic Compounds from Drinking Water via the Ingestion, Inhalation and Dermal Routes. NTIS PB84-213941. Office of Drinking Water. Washington DC: Environmental Protection Agency, 1984.

Edwards, C.A., Beck, S.D. and Lichtenstein, E.P., J. Econ. Entomol. 1957, 50: 622.

Edwards, N. T. 1983. Polycyclic aromatic hydrocarbons (PAHs) in the terrestrial environment – a review. J. Environ. Qual. 12 (4), 427-441.

Falk, H. L., and P. T. S. Kotin. Inhibition of carcinogenesis: The effects of polycyclic hydrocarbons and related compounds. Arch. Environ. Health Vol. 9 (1964):169-179.

Fritz, W. 1971. Extent and sources of contamination of our food with carcinogenic hydrocarbons. Ernaehrungsforschung 16(4), 547-557.

Health & Welfare Canada. Polycyclic Aromatic Hydrocarbons, Report No. 80-EHD-50, (1979) p. 38.

Horst, T. W. Langrangian Similarity Modeling of Vertical Diffusion for a Ground Level Source. Int. Applied Met, Vol. 18 (1979): 733-740.

HSDB (Hazardous Substances Data Bank), 1999. National Library of Medicine (NLM) On-Line Toxicological Network (TOXNET). Bethesda, MD.

ICRP (International Commission on Radiological Protection). Report of Committee IV on Evaluation of Radiation Doses to Body Tissues from Internal Contamination due to Occupational Exposure. ICRP Publication 10. New York: Pergamon Press, 1968.

Kao, J.K., F.K. Patterson, and J. Hall. Skin Penetration and Metabolism of Topically Applied Chemicals in Six Mammalian Species, Including Man: an in vitro Study with Benzo(a)pyrene and Testcaterone. Toxicol. Appl. Pharmacol, Vol. 81 (1985): 502-516.

Kelsey, J.W. and Alexander, M. Declining Bioavailability and Inappropriate Estimation of Risk of Persistent Compounds. Environ. Toxicol. Chem. 1997. 16(3): 582

Kelsey, J.W., Kottler, B.D. and Alexander, M. Environ. Sci. Technol. 1997, 31: 214.

Kissel, J., K. Richter, and R. Fenske. Field Measurements of Dermal Soil Loading Attributable to Various Activities: Implications for Exposure Assessment. Risk Analysis, Vol. 16, No. 1 (1996): 115-125.

Magee, B., P. Anderson, and D. Burmaster. Absorption Adjustment Factor (AAF) Distributions for Polycyclic Aromatic Hydrocarbons (PAHs). Human and Ecological Assessment: An International Journal. Vol. 2, No. 4 (December 1996): 841-873.

Menzi, C.A., B.B. Potocki and J. Santodonato. Exposure to Carcinogenic PAHs in the Environment. Environ. Sci. Technol. Vol. 26, No. 7, 1992.

Michael Pisani & Associates. Remedial Investigation Report, Former Gulf States Creosoting Site, Hattiesburg, Mississippi. New Orleans, Louisiana. 1997.

Michael Pisani & Associates. Phase II Remedial Investigation Report, Former Gulf States Creosoting Site, Hattiesburg, Mississippi. New Orleans, Louisiana. 1998.

Mississippi Commission on Environmental Quality (MCEQ). Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi. 1999.

Mississippi Department of Environmental Quality (MDEQ). Guidance for Remediation of Uncontrolled Hazardous Substance Sites in Mississippi. Office of Pollution Control. 1990.

Mississippi Department of Environmental Quality (MDEQ). Letter to Glen Pilie, Adams and Reese, from Tony Russell, MDEQ. August 2, 2000.

Mississippi Department of Environmental Quality (MDEQ). Letter to Glen Pilie, Adams and Reese, from Tony Russell, MDEQ. February 6, 2001.

Morrison, DE., Robertson, B.K. and Alexander, M. Bioavailability to Earthworms of Aged DDT, DDE, DDD, and Dieldrin in Soil. Environ. Sci. Technol. 2000 34: 709.

Moore, M.R., P.A. Meredith, W.S. Watson, D.J. Sumner, M.K. Taylor, and A. Goldberg. "The Percutaneous Absorption of Lead-203 in Humans From Cosmetic Preparations Containing Lead Acetate, as Assessed by Whole-Body Counting and Other Techniques." Food Cosmet. Toxicol. 18. (1980): 399.

Pancirov, R. J. and R. A. Brown. 1975. Analytical methods for polynuclear aromatic hydrocarbons in crude oil, heating oils, and marine tissues. In: Conference on prevention and control of oil pollution, San Francisco, CA, March, 1975. American Petroleum Institute, Wash., DC. pp 103-13.

Pao, E. M. *et al.* Home Economics Research Report No. 44. United States Department of Agriculture, Washington, DC. 1982.

Pasquill, I.. The Dispersion of Material in the Atmospheric Boundary Layer - The Basis for Generalization. In: Lectures on Air Pollution and Environmental Impact Analysis. Boston, MA: American Meteorological Society, 1975.

Ryan, E.A., E.T. Hawkins *et al.* "Assessing Risk From Dermal Exposure at Hazardous Waste Sites. in Bennett." Ed. G. and J. Bennett. Superfund '87: Proceedings of the Eighth National Conference. Washington, DC, 16-18 November 1987. The Hazardous Material Control Research Institute. p.166-168.

Santodonato, J., P. Howard, and D. Basu. Health and Ecological Assessment of Polynuclear Aromatic Hydrocarbons. Pathotox Publishers, Inc., Park Forest South, IL. 1981.

Shabad, L. M. 1980. Circulation of carcinogenic polycyclic aromatic hydrocarbons in the human environment and cancer prevention. *J. Natl. Cancer Inst.* 64(3): 405-410.

Shu, H.P., P. Teitelbaum, A.S. Webb, L. Marple, B. Brunck, D. Dei Rossi, F.J. Murray, and D.J. Paustenbach. "Bioavailability of Soil Bound TCDD: Dermal Bioavailability in the Rat." *Fundam. Appl. Toxicol.*, Vol. 10 (1988): 648-654.

Smirnov, G. A. 1970. The study of benzo(a)pyrene content in soil and vegetation in the airfield region. *Vopr. Onkol.* 16(5): 83-86.

State of Mississippi. Mississippi Code 1972 Annotated. Title 29 Public Lands, Buildings, and Property, Chapter 3, Sixteenth Section and Lieu Lands in General. 1998.

Suess, M. J. 1976. The environmental load and cycle of polycyclic aromatic hydrocarbons. *Sci. Total Environ.* 6:239-250.

Ta, Roy *et al.*, Studies Estimating the Dermal Bioavailability of Polynuclear Aromatic Hydrocarbons from Manufactured Gas Plan Tar-Contaminated Soils. *Env. Sci. Tech.* 32(20). 1998. 3113-3117.

Tang, W.C., White, J.C., and Alexander, M. *Appl. Microbiol. Biotechnol.* 1998, 49: 117.

Tang, J and Alexander, M. *Environ. Toxicol. Chem.* 1999, 18: 2711.

Thomas, J.F., M. Mukai, and B.D. Teggens. Fate of airborne benzo(a)pyrene. *Environ. Sci. Technol.* 2:33-39, 1968.

United States Department of Agriculture (USDA). Nationwide Food Consumption Survey: Continuing Survey of Food Intakes by Individuals, Men 19-50 years, 1 Day, 1985; United States Department of Agriculture. Human Nutrition Information Service. Nutrition Monitoring Division; Washington, DC, Report No. 86-1. 1986.

US EPA (United States Environmental Protection Agency). Rapid Assessment of Exposure to Particulate Emission from Surface Contamination Sites. EPA/OHEA/EPA. 600/8-85/002. Cowherd, C., Jr., G.E. Muleski, P.J. Engelhart and D.A. Gillett, Ed. Washington DC: Midwest Research Inst.. 1985.

US EPA (United States Environmental Protection Agency). "Guidelines for Carcinogenic Risk Assessment." Federal Register 51:33992-34003. 1986.

US EPA (United States Environmental Protection Agency). Superfund Exposure Assessment Manual (SEAM). EPA 540/1-88/001. Office of Remedial Response. Washington, DC, 1988.

US EPA (United States Environmental Protection Agency). Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A). Interim Final. EPA/540/1-89/002. Office of Emergency and Remedial Response. Washington, DC, 1989.

US EPA (United States Environmental Protection Agency). Human Health Evaluation Manual, Supplemental Guidance: Standard Default Exposure Factors. OSWER Directive. 9285.6-03. Office of Solid Waste and Emergency Response. Washington, DC, 1991.

US EPA (United States Environmental Protection Agency). Dermal Exposure Assessment: Principles and Applications. Office of Research and Development. EPA/600/8-91/011B. Washington, DC, 1992.

US EPA (United States Environmental Protection Agency). Estimating Exposure to Dioxin-Like Compounds. Office of Research and Development. EPA/600/6-88/005B. Washington, DC, 1992.

US EPA (United States Environmental Protection Agency). Provisional Guidance for Quantitative Risk Assessment of Polynuclear Aromatic Hydrocarbons. Office of Solid Waste and Environmental Remediation. EPA/600/R-93/089, July 1993

US EPA (United States Environmental Protection Agency). Land Use Directive in the CERCLA Remedy Selection Process. OSWER Directive 9355.7-04. Office of Solid Waste and Emergency Response. Washington, DC, May 1995.

US EPA (United States Environmental Protection Agency) Region 3. Technical Guidance Manual: Risk Assessment, Assessing Dermal Exposure From Soil. EPA/903-K-95-003. Office of Superfund Programs, Hazardous Waste Management Division. Washington, DC, 1995.

US EPA (United States Environmental Protection Agency) Region 4. Technical Services Supplemental Guidance to RAGS: Region 4 Bulletins. Waste Management Division, Atlanta, GA. 1995.

US EPA (United States Environmental Protection Agency). Supplement B to Compilation of Air Pollutant Emission Factors, Volume I: Stationary Point and Area Sources. AP-42, Fifth Edition, Supplement B. Office of Air Quality Planning and Standards, Office of Air and Radiation. Research Triangle Park, NC, 1996

US EPA (United States Environmental Protection Agency). Exposure Factors Handbook,. EPA/600/P-95/002F. Office of Research and Development, Washington, DC, August 1997.

US EPA (United States Environmental Protection Agency). Health Effects Assessment Summary Tables (HEAST). Office of Health and Environmental Assessment, Environmental Criteria and Assessment Office (ECAO). Cincinnati OH, 1997.

US EPA (United States Environmental Protection Agency). IRIS (Integrated Risk Information System). A Continuously Updated Electronic Database Maintained by the US Environmental Protection Agency. Bethesda, Maryland: National Library of Medicine, 1999.

US EPA (United States Environmental Protection Agency) Region 3. Updated Risk-Based Concentration Tables. Office of RCRA Technical & Program Support Branch. Philadelphia, PA, April, 1999.

Wallcave, L., H. Garcia, R. Fedlman, W. Linjinsky, and P. Shubik. 1971. Skin tumorigenesis in mice by petroleum asphalts and coal-tar pitches of known polynuclear aromatic hydrocarbon content. *Toxicol. Appl. Pharmacol.* 18, 41-52.

Weissenfels, W.D., Klewer, H.J. and Langhoff, J., J. *Appl. Microbiol. Biotechnol.* 1992, 36: 689.

White, J.C., Kelsey, J.W., Hatzinger, P.B and Alexander, M. Factors Affecting Sequestration and Bioavailability of Phenanthrene in Soils. *Environ. Toxicol. Chem.* 1997, 16(10): 2040.

Youngblood, W. W., and M. Blumer. 1975. Polycyclic aromatic hydrocarbons in the environment: homologous series in soils and recent marine sediments. *Geochim. Cosmochim. Acta* 39:1303-1315.

*Figure 1*

**Site Conceptual Model and Selection of Exposure Pathways**  
**Kerr McGee, Hattiesburg, MS**

| Scenario Timeframe | Medium   | Exposure Medium     | Exposure Point  | Receptor Population | Receptor Age | Exposure Route         | On-Site/Off-Site | Type of Analysis | Rationale for Selection or Exclusion of Exposure Pathway                             |
|--------------------|----------|---------------------|-----------------|---------------------|--------------|------------------------|------------------|------------------|--|
| Current            | Soil     | Surface Soil (0-1') | Exposure Unit 1 | Visitor             | Adolescent   | Dermal Oral Inhalation | On-Site          | None             | EU1 includes only surface water and sediment. Soils in this area are included in EU2 |
|                    |          |                     | Exposure Unit 2 | Visitor             | Adolescent   | Dermal Oral            | On-Site          | None             | EU1 includes only surface water and sediment. Soils in this area are included in EU2 |
|                    |          |                     | Exposure Unit 3 | Visitor             | Adolescent   | Dermal Oral            | On-Site          | Quantitative     | EU1 includes only surface water and sediment. Soils in this area are included in EU2 |
|                    |          |                     | Exposure Unit 4 | Visitor             | Adolescent   | Dermal Oral            | On-Site          | Quantitative     | Area potentially attractive for occasional recreational use                          |
|                    |          |                     | Exposure Unit 5 | Visitor             | Adolescent   | Dermal Oral            | On-Site          | Quantitative     | Area potentially attractive for occasional recreational use                          |
|                    |          |                     | Exposure Unit 1 | Maintenance Worker  | Adult        | Dermal Oral            | On-Site          | Quantitative     | Area potentially attractive for occasional recreational use                          |
|                    |          |                     | Exposure Unit 2 | Maintenance Worker  | Adult        | Dermal Oral            | On-Site          | Quantitative     | VOC's not present at levels of concern   |
|                    |          |                     | Exposure Unit 3 | Maintenance Worker  | Adult        | Dermal Oral            | On-Site          | Quantitative     | VOC's not present at levels of concern   |
|                    |          |                     | Exposure Unit 4 | Maintenance Worker  | Adult        | Dermal Oral            | On-Site          | Quantitative     | VOC's not present at levels of concern   |
|                    |          |                     | Exposure Unit 5 | Maintenance Worker  | Adult        | Dermal Oral            | On-Site          | Quantitative     | VOC's not present at levels of concern   |
|                    | Sediment | Exposure Unit 1     | Visitor         | Adolescent          | Dermal Oral  | Inhalation             | On-Site          | Quantitative     | Visitors may potentially wade in Gordon's Creek                                      |
|                    |          | Exposure Unit 4     | Visitor         | Adolescent          | Dermal Oral  | Inhalation             | On-Site          | Quantitative     | VOC's not present at levels of concern   |



Figure 1

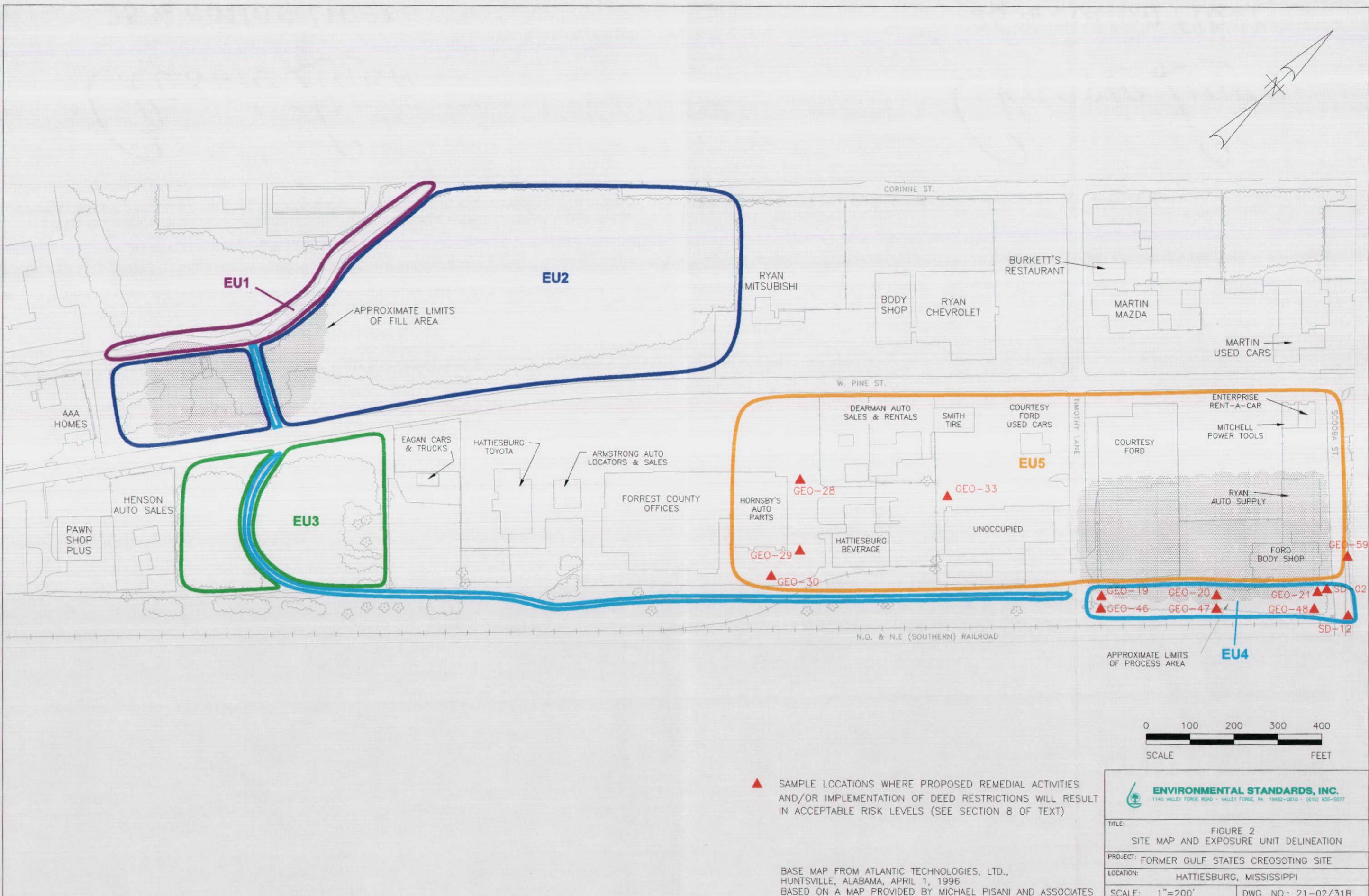
**Site Conceptual Model and Selection of Exposure Pathways**  
**Kerr McGee, Hattiesburg, MS**

| Scenario | Medium | Exposure Medium | Exposure Point  | Receptor Population | Receptor Age | Exposure Route               | On-Site/Off-Site | Type of Analysis                             | Rationale for Selection or Exclusion of Exposure Pathway  |
|----------|--------|-----------------|-----------------|---------------------|--------------|------------------------------|------------------|--|---|
|          |        |                 | Exposure Unit 1 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | None   | Sediment exposures addressed in EU1 under a future scenario   |
|          |        |                 | Exposure Unit 4 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | None   | Sediment exposures addressed in EU1 under a future scenario   |
|          |        |                 | Exposure Unit 6 | Resident            | Child/Adult  | Dermal<br>Oral<br>Inhalation | Off-Site         | Quantitative<br>Quantitative<br>Quantitative | Sediment exposures addressed in EU1 under a future scenario   |
|          |        |                 |                 |                     |              |                              |                  |  | Infrequent maintenance of drainage ditch<br>Infrequent maintenance of drainage ditch<br>VOCs not present at levels of concern                       |
|          |        |                 |                 |                     |              |                              |                  |  | Playing/working in drainage ditch<br>Playing/working in drainage ditch<br>VOCs not present at levels of concern                                     |
|          |        |                 |                 |                     |              |                              |                  |  |   |
|          |        |                 | Exposure Unit 1 | Visitor             | Adolescent   | Dermal<br>Oral<br>Inhalation | On-Site          | Quantitative<br>Quantitative<br>None         | Visitor may potentially wade in Gordon's Creek<br>Visitor may potentially wade in Gordon's Creek<br>VOCs not present at levels of concern           |
|          |        |                 | Exposure Unit 4 | Visitor             | Adolescent   | Dermal<br>Oral<br>Inhalation | On-Site          | Quantitative<br>Quantitative<br>None         | Visitor may potentially walk through drainage ditch<br>Visitor may potentially walk through drainage ditch<br>VOCs not present at levels of concern |
|          |        |                 |                 |                     |              |                              |                  |  |   |
|          |        |                 | Exposure Unit 1 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | None   | Surface Water exposures addressed in EU1 under a future scenario  |
|          |        |                 | Exposure Unit 4 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | None   | Surface Water exposures addressed in EU1 under a future scenario  |
|          |        |                 | Exposure Unit 6 | Resident            | Child/Adult  | Dermal<br>Oral<br>Inhalation | Off-Site         | Quantitative<br>Quantitative<br>None         | Surface Water exposures addressed in EU1 under a future scenario  |
|          |        |                 |                 |                     |              |                              |                  |  | Infrequent maintenance of drainage ditch<br>Infrequent maintenance of drainage ditch<br>VOCs not present at levels of concern                       |
|          |        |                 |                 |                     |              |                              |                  |  |   |
|          |        |                 | Exposure Unit 1 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | None   | EU1 includes only surface water and sediment. Soils in this area are included in EU2  |
|          |        |                 | Exposure Unit 2 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | None   | EU1 includes only surface water and sediment. Soils in this area are included in EU2  |
|          |        |                 | Exposure Unit 3 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | None   | EU1 includes only surface water and sediment. Soils in this area are included in EU2  |
|          |        |                 | Exposure Unit 4 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | Quantitative<br>Quantitative<br>None         | May potentially become a maintained area<br>COPCs not present at levels of concern  |
|          |        |                 | Exposure Unit 5 | Maintenance Worker  | Adult        | Dermal<br>Oral<br>Inhalation | On-Site          | None   | May potentially become a maintained area<br>COPCs eliminated during screening process   |
|          |        |                 |                 |                     |              |                              |                  |  | Surface Soil exposures addressed in EU4 under a current scenario<br>Surface Soil exposures addressed in EU4 under a current scenario                |
|          |        |                 |                 |                     |              |                              |                  |  | May potentially become a maintained area<br>VOCs not present at levels of concern   |
|          |        |                 |                 |                     |              |                              |                  |  |   |



**Figure 1**  
**Site Conceptual Model and Selection of Exposure Pathways**  
**Kerr McGee, Hattiesburg, MS**

| Scenario Timeframe | Medium          | Exposure Medium                     | Exposure Point      | Receptor Population | Receptor Age           | Exposure Route         | On-Site/Off-Site                       | Type of Analysis  | Rationale for Selection or Exclusion of Exposure Pathway  |  |
|--------------------|-----------------|-------------------------------------|---------------------|---------------------|------------------------|------------------------|--|---|---|--|
|                    |                 |                                     |                     |                     |                        |                        |  |   | EU1 includes only surface water and sediment. Soils in this area are included in EU2  |  |
|                    | Subsurface Soil | Subsurface Soil (0' to water table) | Exposure Unit 1     | Construction Worker | Adult                  | Dermal Oral Inhalation | On-Site                                | None  | EU1 includes only surface water and sediment. Soils in this area are included in EU2  |  |
|                    |                 |                                     | Exposure Unit 2     | Construction Worker | Adult                  | Dermal Oral Inhalation | On-Site                                | None  | EU1 includes only surface water and sediment. Soils in this area are included in EU2  |  |
|                    |                 |                                     | Exposure Unit 3     | Construction Worker | Adult                  | Dermal Oral Inhalation | On-Site                                | Quantitative Quantitative Quantitative  | Potentially constructable area in the future Potentially constructable area in the future Non-VOC entrained fugitive dust generation during potential construction activities                               |  |
|                    |                 |                                     | Exposure Unit 4     | Construction Worker | Adult                  | Dermal Oral Inhalation | On-Site                                | None  | COPCs eliminated during screening process   |  |
|                    |                 |                                     |                     |                     |                        |                        |  | None  | COPCs eliminated during screening process   |  |
|                    |                 |                                     |                     |                     |                        |                        |  | None  | COPCs eliminated during screening process   |  |
|                    |                 |                                     | Exposure Unit 5     | Construction Worker | Adult                  | Dermal Oral Inhalation | On-Site                                | Quantitative Quantitative Quantitative  | In frequent construction activities may occur in the future In frequent construction activities may occur in the future Non-VOC entrained fugitive dust generation during potential construction activities |  |
|                    |                 |                                     |                     |                     |                        |                        |  | None  |   |  |
| Sediment           | Sediment        | Exposure Unit 1                     | Maintenance Worker  | Adult               | Dermal Oral Inhalation | On-Site                | Quantitative Quantitative Quantitative | Potentially constructable area in the future Potentially constructable area in the future Non-VOC entrained fugitive dust generation during potential construction activities                         |   |  |
|                    |                 | Exposure Unit 4                     | Maintenance Worker  | Adult               | Dermal Oral Inhalation | On-Site                | None                                   | In frequent maintenance of Gordon's Creek   |   |  |
|                    |                 |                                     |                     |                     |                        |                        |  | VOCs not present at levels of concern   |   |  |
|                    |                 |                                     |                     |                     |                        |                        |  |   |   |  |
|                    |                 | Exposure Unit 1                     | Construction Worker | Adult               | Dermal Oral Inhalation | On-Site                | Quantitative Quantitative Quantitative | Sediment exposures addressed in EU4 under a current scenario Sediment exposures addressed in EU4 under a current scenario Sediment exposures addressed in EU4 under a current scenario                |   |  |
|                    |                 | Exposure Unit 4                     | Construction Worker | Adult               | Dermal Oral Inhalation | On-Site                | None                                   | In frequent construction activities may occur in the future   |   |  |
|                    |                 |                                     |                     |                     |                        |                        |  | VOCs not present at levels of concern   |   |  |
|                    |                 |                                     |                     |                     |                        |                        |  |   |   |  |
| Surface Water      | Surface Water   | Exposure Unit 1                     | Maintenance Worker  | Adult               | Dermal Oral Inhalation | On-Site                | Quantitative Quantitative Quantitative | In frequent construction activities may occur in the future In frequent construction activities may occur in the future VOCs not present at levels of concern   |   |  |
|                    |                 | Exposure Unit 4                     | Maintenance Worker  | Adult               | Dermal Oral Inhalation | On-Site                | None                                   | In frequent maintenance of Gordon's Creek   |   |  |
|                    |                 |                                     |                     |                     |                        |                        |  | In frequent maintenance of Gordon's Creek   |   |  |
|                    |                 |                                     |                     |                     |                        |                        |  |   |   |  |
|                    |                 | Exposure Unit 1                     | Construction Worker | Adult               | Dermal Oral Inhalation | On-Site                | Quantitative Quantitative Quantitative | Surface Water exposures addressed in EU4 under a current scenario Surface Water exposures addressed in EU4 under a current scenario Surface Water exposures addressed in EU4 under a current scenario |   |  |
|                    |                 | Exposure Unit 4                     | Construction Worker | Adult               | Dermal Oral Inhalation | On-Site                | None                                   | In frequent construction activities may occur in the future   |   |  |
|                    |                 |                                     |                     |                     |                        |                        |  | VOCs not present at levels of concern   |   |  |
|                    |                 |                                     |                     |                     |                        |                        |  |   |   |  |



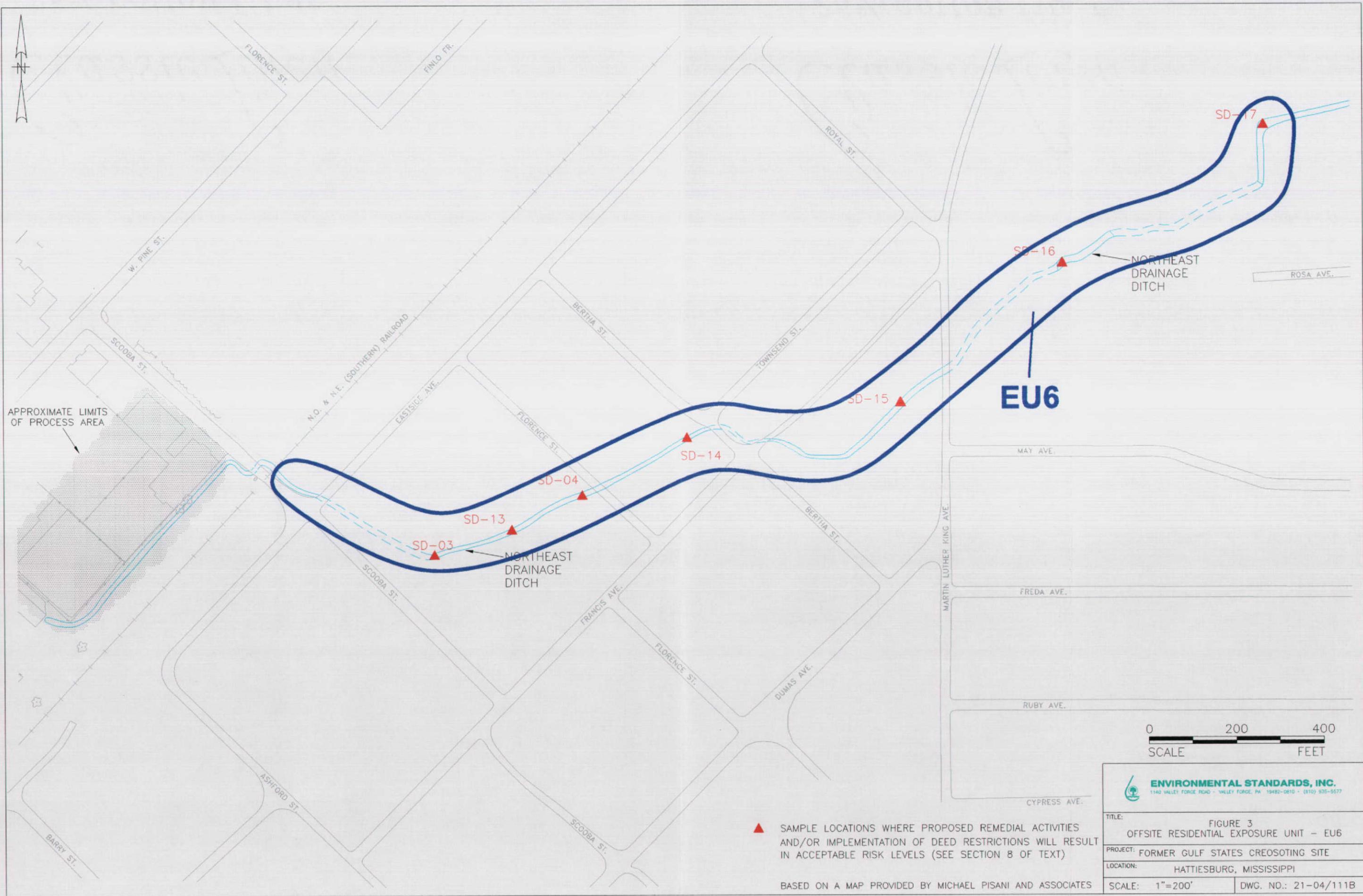
▲ SAMPLE LOCATIONS WHERE PROPOSED REMEDIAL ACTIVITIES  
AND/OR IMPLEMENTATION OF DEED RESTRICTIONS WILL RESULT  
IN ACCEPTABLE RISK LEVELS (SEE SECTION 8 OF TEXT)

BASE MAP FROM ATLANTIC TECHNOLOGIES, LTD.,  
HUNTSVILLE, ALABAMA, APRIL 1, 1996  
BASED ON A MAP PROVIDED BY MICHAEL PISANI AND ASSOCIATES

|           |  |                     |
|-----------|--|---------------------|
| TITLE:    | FIGURE 2<br>SITE MAP AND EXPOSURE UNIT DELINEATION |                     |
| PROJECT:  | FORMER GULF STATES CREOSOTING SITE                 |                     |
| LOCATION: | HATTIESBURG, MISSISSIPPI                           |                     |
| SCALE:    | 1"=200'  | DWG. NO.: 21-02/31B |



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|  |   |
|--|---|
| <b>ENVIRONMENTAL STANDARDS, INC.</b><br><small>1140 VALLEY FORCE ROAD - VALLEY FORCE, PA 15482-0810 - (812) 935-5577</small> |   |
| <b>TITLE:</b>  | FIGURE 3<br>OFFSITE RESIDENTIAL EXPOSURE UNIT - EU6 |
| <b>PROJECT:</b>  | FORMER GULF STATES CREOSOTING SITE                  |
| <b>LOCATION:</b>   | HATTIESBURG, MISSISSIPPI                            |
| <b>SCALE:</b>  | 1"=200' DWG. NO.: 21-04/111B                        |

**Table 25****Oral Exposure to EU1 Sediment by an Adolescent Visitor (Aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u><math>Cd * IngR * EF * ED * CF * ME</math></u> |                       |  |  |
|--|-----------|---|-----------------------|--|--|
|  |           | <u><math>BW * AT</math></u>                       |                       |  |  |
| Cd - Concentration in sediment =                     | mg/kg     | see below   |                       |  |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 100   | USEPA 1994, Region I  |  |  |
| EF - Exposure frequency =                            | days/year | 12  | USEPA 1991, HHEM      |  |  |
| ED - Exposure duration =                             | years     | 10  | reasonable assumption |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06  |                       |  |  |
| ME - Matrix effect =                                 |           | 1   | reasonable assumption |  |  |
| BW - Body weight =                                   | kg        | 45  | USEPA 1997, EFH       |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650  | reasonable assumption |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550   | USEPA 1991, HHEM      |  |  |

| Constituent            | Concentration<br>in Sediment<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average                         |                              |  |
|------------------------|---------------------------------------|--------------------------------------|----------------------------------|-----------------|---------------------------------|------------------------------|--|
|                        |                                       |                                      |                                  |                 | Lifetime<br>Intake<br>mg/kg-day | Daily<br>Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) |
| <b>Semivolatiles</b>   |                                       |                                      |                                  |                 |                                 |                              |  |
| Benzo(a)anthracene     | 5.90E-01                              | 4.31E-08                             | NA                               | NA              | 6.16E-09                        | 7.30E-01                     | 4.50E-09                                     |
| Benzo(a)pyrene         | 3.90E-01                              | 2.85E-08                             | NA                               | NA              | 4.07E-09                        | 7.30E+00                     | 2.97E-08                                     |
| Benzo(b)fluoranthene   | 5.80E-01                              | 4.24E-08                             | NA                               | NA              | 6.05E-09                        | 7.30E-01                     | 4.42E-09                                     |
| Benzo(k)fluoranthene   | 1.90E-01                              | 1.39E-08                             | NA                               | NA              | 1.98E-09                        | 7.30E-02                     | 1.45E-10                                     |
| Chrysene               | 5.30E-01                              | 3.87E-08                             | NA                               | NA              | 5.53E-09                        | 7.30E-03                     | 4.04E-11                                     |
| Dibenz(a,h)anthracene  | 6.20E-02                              | 4.53E-09                             | NA                               | NA              | 6.47E-10                        | 7.30E+00                     | 4.72E-09                                     |
| Indeno(1,2,3-cd)pyrene | 2.20E-01                              | 1.61E-08                             | NA                               | NA              | 2.30E-09                        | 7.30E-01                     | 1.68E-09                                     |

Total Cancer Risk = 4.52E-08



**Table 2****Dermal Exposure to EU1 Surface Water by an Adolescent Visitor (aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                                     | $C_w * SA * K_p * ABS * ET * EF * ED * CF$<br>BW * AT |  |                              |                                 |                             |             |          |
|---|-------------------------------------|---|--|------------------------------|---------------------------------|-----------------------------|-------------|----------|
| Cw - Concentration in surface water =                                   | mg/L                                | see below   |  |                              |                                 |                             |             |          |
| SA - Surface area available for exposure =                              | cm <sup>2</sup>                     | 3945  | calculated                             |                              |                                 |                             |             |          |
| SA <sub>t</sub> - Total skin surface area =                             | cm <sup>2</sup>                     | 12768.3   | USEPA 1997, EFH                        |                              |                                 |                             |             |          |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = |                                     | 30.9%   | USEPA 1997, EFH                        |                              |                                 |                             |             |          |
| K <sub>p</sub> - Dermal permeability constant =                         | cm/hr                               | see below   |  |                              |                                 |                             |             |          |
| ABS <sub>p</sub> - Absorption - cPAHs =                                 |                                     | 0.03  | USEPA 1995, Region III                 |                              |                                 |                             |             |          |
| ET - Exposure time =  | hrs/day                             | 1   | USEPA 1992, Dermal Exposure Assessment |                              |                                 |                             |             |          |
| EF - Exposure frequency =   | days/year                           | 12  | reasonable assumption                  |                              |                                 |                             |             |          |
| ED - Exposure duration =  | years                               | 10  | USEPA 1995, Region IV                  |                              |                                 |                             |             |          |
| CF - Conversion factor =  | L/cm <sup>3</sup>                   | 1.00E-03  |  |                              |                                 |                             |             |          |
| BW - Body weight =  | kg                                  | 45  | USEPA 1995, Region IV                  |                              |                                 |                             |             |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                    | days                                | 3650  | USEPA 1991, HHEM                       |                              |                                 |                             |             |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =                       | days                                | 25550   | USEPA 1991, HHEM                       |                              |                                 |                             |             |          |
| Constituent   | Concentration in Surface Water mg/L | K <sub>p</sub> cm/hr                                  | Average Daily Intake mg/kg-day         | Dermal Chronic RfD mg/kg-day | Lifetime Daily Intake mg/kg-day | Average Cancer Slope Factor |             |          |
|   |                                     |   |  |                              |                                 | 1/(mg/kg-day)               | Cancer Risk |          |
| <b>Semivolatiles</b>  |                                     |   |  |                              |                                 |                             |             |          |
| Benzo(a)anthracene  | 1.00E-03                            | 8.10E-01  | 7.00E-08                               | NA                           | NA                              | 1.00E-08                    | 1.46E+00    | 1.46E-08 |
| Benzo(a)pyrene  | 5.00E-04                            | 1.20E+00  | 5.19E-08                               | NA                           | NA                              | 7.41E-09                    | 1.46E+01    | 1.08E-07 |
| Benzo(b)fluoranthene  | 5.00E-04                            | 1.20E+00  | 5.19E-08                               | NA                           | NA                              | 7.41E-09                    | 1.46E+00    | 1.08E-08 |
| Benzo(k)fluoranthene  | 5.00E-04                            | 4.48E+01  | 1.94E-06                               | NA                           | NA                              | 2.77E-07                    | 1.46E-01    | 4.04E-08 |
| Chrysene  | 5.00E-04                            | 8.10E-01  | 3.50E-08                               | NA                           | NA                              | 5.00E-09                    | 1.46E-02    | 7.30E-11 |
| Dibenz(a,h)anthracene   | 5.00E-04                            | 2.70E+00  | 1.17E-07                               | NA                           | NA                              | 1.67E-08                    | 1.46E+01    | 2.43E-07 |
| Indeno(1,2,3-cd)pyrene  | 5.00E-04                            | 1.90E+00  | 8.22E-08                               | NA                           | NA                              | 1.17E-08                    | 1.46E+00    | 1.71E-08 |

NA - Not Available

Total Cancer Risk = 4.35E-07



*Statistical Summary and Selection of COPCs in EU1 Sediment  
Kerr McGee, Hattiesburg, MS*

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Detected | Qualifier | Mean mg/kg | Logarithmic Mean | Maximum mg/kg | Detected | Maximum mg/kg | Detected | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|----------|-----------|------------|------------------|---------------|----------|---------------|----------|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>   |            |               |                |                 |               |               |                       |          |               |          |           |            |                  |               |          |               |          |                                   |                          |
| 2-methylphthalene      | 91-57-6    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 7.40E-02              | J        | 2.92E-01      | 1.94E-01 |           | 5.10E-01   |                  | SD-07         |          | 3.08E-01      |          |                                   |                          |
| Acenaphthene           | 83-32-9    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.80E-01              | J        | 3.15E-01      | 2.85E-01 |           | 4.50E-01   |                  | SD-07         |          | 1.91E-01      |          |                                   |                          |
| Acenaphthylene         | 208-96-8   | 2             | 1              | 50              | 4.00E-02      | 4.00E-02      | 1.80E-02              | J        | 4.90E-02      | 3.95E-02 |           | 7.80E-02   | J                | SD-07         |          | 4.10E-02      |          |                                   |                          |
| Anthracene             | 120-12-7   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 2.60E-01              | J        | 3.60E-01      | 3.46E-01 |           | 4.60E-01   |                  | SD-07         |          | 1.41E-01      |          |                                   |                          |
| Benz(a)anthracene      | 56-55-3    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.80E-01              | J        | 3.85E-01      | 3.26E-01 |           | 5.90E-01   |                  | SD-07         |          | 2.90E-01      |          |                                   |                          |
| Benz(a)pyrene          | 50-32-8    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.20E-01              | J        | 2.55E-01      | 2.16E-01 |           | 3.90E-01   | J                | SD-07         |          | 1.91E-01      |          |                                   |                          |
| Benz(b)fluoranthene    | 205-99-2   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.70E-01              | J        | 3.75E-01      | 3.14E-01 |           | 5.80E-01   |                  | SD-07         |          | 2.90E-01      |          |                                   |                          |
| Benz(ghi)perylene      | 191-24-2   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 6.50E-02              | J        | 1.23E-01      | 1.08E-01 |           | 1.80E-01   | J                | SD-07         |          | 8.13E-02      |          |                                   |                          |
| Benz(k)fluoranthene    | 207-08-9   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 6.40E-02              | J        | 1.27E-01      | 1.10E-01 |           | 1.90E-01   | J                | SD-07         |          | 8.91E-02      |          |                                   |                          |
| Carbazole              | 86-74-8    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.60E-01              | J        | 3.65E-01      | 3.02E-01 |           | 5.70E-01   |                  | SD-07         |          | 2.90E-01      |          |                                   |                          |
| Chrysene               | 218-01-9   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.80E-01              | J        | 3.55E-01      | 3.09E-01 |           | 5.30E-01   |                  | SD-07         |          | 2.47E-01      |          |                                   |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 2             | 1              | 50              | 4.00E-02      | 4.00E-02      | 6.20E-02              | J        | 4.10E-02      | 3.52E-02 |           | 6.20E-02   | J                | SD-07         |          | 2.97E-02      |          |                                   |                          |
| Dibenzofuran           | 132-64-9   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.50E-01              | J        | 2.80E-01      | 2.48E-01 |           | 4.10E-01   |                  | SD-07         |          | 1.84E-01      |          |                                   |                          |
| Fluoranthene           | 206-44-0   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 6.80E-01              | J        | 1.19E+00      | 1.08E+00 |           | 1.70E+00   |                  | SD-07         |          | 7.21E-01      |          |                                   |                          |
| Fluorene               | 86-73-7    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 2.30E-01              | J        | 4.25E-01      | 3.78E-01 |           | 6.20E-01   |                  | SD-07         |          | 2.76E-01      |          |                                   |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 6.90E-02              | J        | 1.45E-01      | 1.23E-01 |           | 2.20E-01   | J                | SD-07         |          | 1.97E-01      |          |                                   |                          |
| Naphthalene            | 91-20-3    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.80E-01              | J        | 6.40E-01      | 4.45E-01 |           | 1.10E+00   |                  | SD-07         |          | 6.51E-01      |          |                                   |                          |
| Phenanthrene           | 85-01-8    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 7.20E-01              | J        | 1.21E+00      | 1.11E+00 |           | 1.70E+00   |                  | SD-07         |          | 6.93E-01      |          |                                   |                          |
| Pyrene                 | 129-00-0   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 4.80E-01              | J        | 9.40E-01      | 8.20E-01 |           | 1.40E+00   |                  | SD-07         |          | 6.51E-01      |          |                                   |                          |

**Table 1**  
**Statistical Summary and Selection of COPCs in EU1 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Unrestricted<br>Self TRG<br>mg/kg | Is Maximum<br>Detected ><br>TRG? | Is the 95%<br>UCL ><br>TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|--|----------------------------------|-----------------------------|
| <b>Semi-volatiles</b>  |                  |                                 |                                   |  |  |                                  |                             |
| 2-methylnaphthalene    | 1.67E+00         | 1.60E+22                        | Unknown                           | 5.10E-01                                 | 3.13E+03                                 | no                               |                             |
| Acenaphthene           | 1.17E+00         | 3.23E+04                        | Unknown                           | 4.50E-01                                 | 4.69E+03                                 | no                               |                             |
| Acenaphthylene         | 2.32E-01         | 8.34E+09                        | Unknown                           | 7.80E-02                                 | 4.69E+03                                 | no                               |                             |
| Anthracene             | 9.91E-01         | 2.23E+01                        | Unknown                           | 4.60E-01                                 | 2.35E+04                                 | no                               |                             |
| Benz(a)anthracene      | 1.68E+00         | 1.25E+08                        | Unknown                           | 5.90E-01                                 | 8.75E-01                                 | yes*                             |                             |
| Benz(a)pyrene          | 1.11E+00         | 6.25E+07                        | Unknown                           | 3.90E-01                                 | 8.75E-02                                 | yes                              |                             |
| Benz(b)fluoranthene    | 1.67E+00         | 4.79E+08                        | Unknown                           | 5.80E-01                                 | 8.75E-01                                 | yes*                             |                             |
| Benz(g)heptaphene      | 4.86E-01         | 2.08E+05                        | Unknown                           | 1.80E-01                                 | 2.35E+03                                 | no                               |                             |
| Benz(k)fluoranthene    | 5.25E-01         | 1.71E+06                        | Unknown                           | 1.90E-01                                 | 8.75E+00                                 | yes*                             |                             |
| Carbazole              | 1.66E+00         | 2.15E+09                        | Unknown                           | 5.70E-01                                 | 3.19E+01                                 | no                               |                             |
| Chrysene               | 1.46E+00         | 3.73E+06                        | Unknown                           | 5.30E-01                                 | 8.75E+01                                 | yes*                             |                             |
| Dibenz(a,h)anthracene  | 1.74E-01         | 2.15E+06                        | Unknown                           | 6.20E-02                                 | 8.75E-02                                 | yes*                             |                             |
| Dibenzofuran           | 1.10E+00         | 3.27E+05                        | Unknown                           | 4.10E-01                                 | 3.13E+02                                 | no                               |                             |
| Fluoranthene           | 4.41E+00         | 1.22E+05                        | Unknown                           | 1.70E+00                                 | 3.13E+03                                 | no                               |                             |
| Fluorene               | 1.66E+00         | 3.35E+05                        | Unknown                           | 6.20E-01                                 | 3.13E+03                                 | no                               |                             |
| Indeno(1,2,3-cd)pyrene | 6.21E-01         | 1.88E+07                        | Unknown                           | 2.20E-01                                 | 8.75E-01                                 | yes*                             |                             |
| Naphthalene            | 3.54E+00         | 6.10E+19                        | Unknown                           | 1.10E+06                                 | 6.45E+02                                 | no                               |                             |
| Phenanthrene           | 4.30E+00         | 2.92E+04                        | Unknown                           | 1.70E+00                                 | 2.35E+03                                 | no                               |                             |
| Pyrene                 | 3.84E+00         | 7.40E+06                        | Unknown                           | 1.40E+00                                 | 2.35E+03                                 | no                               |                             |

\* Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a C

**Table 2**  
**Statistical Summary and Selection of COPCs in EUI Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/L | Maximum mg/L | Mean mg/L | Logarithmic Mean mg/L | Maximum Detected mg/L | Maximum Detected Qualifier | Location of Maximum Concentration | Standard Deviation mg/L |
|------------------------|------------|---------------|----------------|-----------------|--------------|--------------|-----------|-----------------------|-----------------------|----------------------------|-----------------------------------|-------------------------|
| <b>Semi-volatiles</b>  |            |               |                |                 |              |              |           |                       |                       |                            |                                   |                         |
| Benzo(a)anthracene     | 56-55-3    | 2             | 1              | 50              | 1.00E-03     | 1.00E-03     | J         | 7.50E-04              | 7.07E-04              | 1.00E-03                   | J                                 | SW-08                   |
| Benzo(a)pyrene         | 50-32-8    | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00                   | NA                                | SW-08                   |
| Benzo(b)fluoranthene   | 205-99-2   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00                   | NA                                | SW-08                   |
| Benzo(k)fluoranthene   | 207-08-9   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00                   | NA                                | SW-08                   |
| Chrysene               | 218-01-9   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00                   | NA                                | SW-08                   |
| Dibenz(a,h)anthracene  | 53-70-3    | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00                   | NA                                | SW-08                   |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00                   | NA                                | SW-08                   |
| Fluoranthene           | 206-44-0   | 2             | 1              | 50              | 1.00E-03     | 1.00E-03     | 7.50E-03  | 4.00E-03              | 1.94E-03              | 7.50E-03                   | SW-08                             | 4.95E-03                |
| Pyrene                 | 129-00-0   | 2             | 1              | 50              | 1.00E-03     | 1.00E-03     | J         | 7.50E-04              | 7.07E-04              | 1.00E-03                   | J                                 | SW-08                   |

NA - Not applicable: constituent not detected in media.

**Table 2**  
**Statistical Summary and Selection of COPCs in EU1 Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/L | Logarithmic<br>95% UCL<br>mg/L | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/L | Human Health<br>Consumption of Water<br>& Organisms AWQC<br>mg/L | Is Maximum<br>Detected ><br>AWQC? |
|------------------------|-----------------|--------------------------------|-----------------------------------|---|--|-----------------------------------|
| <b>Semivolatiles</b>   |                 |                                |                                   |   |  |                                   |
| Benzo(a)anthracene     | 2.31E-03        | 4.37E-01                       | Unknown                           | 1.00E-03                                | 4.40E-06   | YES - COPC                        |
| Benzo(b)pyrene         | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Benzo(b)fluoranthene   | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Benzo(k)fluoranthene   | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Chrysene               | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Dibenz(a,h)anthracene  | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Indeno(1,2,3-cd)pyrene | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Fluoranthene           | 2.61E-02        | 2.90E+42                       | Unknown                           | 7.50E-03                                | 3.00E-01   | no                                |
| Pyrene                 | 2.33E-03        | 4.37E-01                       | Unknown                           | 1.00E-03                                | 9.60E-01   | no                                |

NA - Not Available

\* Retained as a COPC, as per MDEQ Comments (8/2/2000): constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.



**Table 3**  
**Statistical Summary and Selection of COPCs in EU2 Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | CAS Number | Total Samples | Hit %       | Minimum Detection Limit mg/kg | Maximum Detection Limit mg/kg | Minimum Detected mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Maximum Quantifier mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|-------------|-------------------------------|-------------------------------|------------------------|------------|------------------------|------------------------|--------------------------|-----------------------------------|--------------------------|
|                        |            | Hits          | Frequency % | mg/kg                         | mg/kg                         | Qualifier              | mg/kg      | mg/kg                  | mg/kg                  | mg/kg                    |                                   |                          |
| Semivaluates           |            |               |             |                               |                               |                        |            |                        |                        |                          |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 14            | 2           | 14.29                         | 3.30E-02                      | 7.00E-02               | J          | 3.06E-02               | 2.15E-02               | 1.60E-01                 | J                                 | SS-10                    |
| Acenaphthene           | 83-32-9    | 14            | 1           | 7.14                          | 3.30E-02                      | 4.90E-02               | J          | 1.88E-02               | 1.78E-02               | 4.90E-02                 | J                                 | GEO-13                   |
| Acenaphthylene         | 208-96-8   | 14            | 6           | 42.86                         | 3.30E-02                      | 3.30E-02               | J          | 1.59E-01               | 4.29E-02               | 1.30E-00                 | J                                 | GEO-13                   |
| Anthracene             | 120-12-7   | 14            | 7           | 50                            | 3.30E-02                      | 3.30E-02               | J          | 1.89E-01               | 5.00E-02               | 1.60E+00                 | J                                 | GEO-13                   |
| Benz(a)anthracene      | 56-55-3    | 14            | 12          | 85.71                         | 3.30E-02                      | 3.30E-02               | J          | 8.98E-01               | 2.28E-01               | 6.70E+00                 | J                                 | GEO-13                   |
| Benz(a)pyrene          | 50-32-8    | 14            | 11          | 78.57                         | 6.70E-02                      | 8.40E-02               | J          | 8.31E-01               | 2.82E-01               | 5.20E+00                 | J                                 | GEO-13                   |
| Benz(b)fluoranthene    | 205-99-2   | 14            | 12          | 85.71                         | 6.70E-02                      | 6.70E-02               | J          | 1.84E+00               | 5.95E-01               | 9.20E+00                 | J                                 | GEO-13                   |
| Benz(ghi)perylene      | 191-24-2   | 14            | 10          | 71.43                         | 6.70E-02                      | 6.70E-02               | J          | 5.17E-01               | 2.20E-01               | 2.30E+00                 | J                                 | GEO-13                   |
| Benz(k)fluoranthene    | 207-08-9   | 14            | 9           | 64.29                         | 1.30E-01                      | 1.30E-01               | J          | 7.01E-01               | 2.88E-01               | 3.60E+00                 | J                                 | GEO-13                   |
| Carbazole              | 86-74-8    | 14            | 4           | 28.57                         | 3.30E-02                      | 3.30E-02               | J          | 6.28E-02               | 2.94E-02               | 3.50E-01                 | J                                 | GEO-13                   |
| Cycloene               | 218-01-9   | 14            | 12          | 85.71                         | 3.30E-02                      | 3.30E-02               | J          | 1.19E+00               | 3.11E-01               | 8.00E+00                 | J                                 | GEO-13                   |
| Dibenz(a,h)anthracene  | 53-70-3    | 14            | 7           | 50                            | 6.70E-02                      | 7.20E-02               | J          | 1.85E-01               | 8.87E-02               | 9.10E-01                 | J                                 | GEO-13                   |
| Dibenzofuran           | 132-64-9   | 14            | 2           | 14.29                         | 3.30E-02                      | 7.20E-02               | J          | 2.63E-02               | 2.08E-02               | 9.80E-02                 | J                                 | SS-10                    |
| Di-n-butylphthalate    | 84-74-2    | 14            | 9           | 64.29                         | 3.30E-02                      | 7.20E-02               | J          | 4.30E-02               | 3.68E-02               | 1.10E-01                 | J                                 | SS-10                    |
| Fluoranthene           | 206-44-0   | 14            | 12          | 85.71                         | 3.30E-02                      | 6.60E-02               | J          | 1.40E+00               | 3.90E-01               | 1.20E+01                 | J                                 | GEO-13                   |
| Fluorene               | 86-73-7    | 14            | 2           | 14.29                         | 3.30E-02                      | 4.50E-02               | J          | 4.38E-02               | 2.21E-02               | 3.70E-01                 | J                                 | GEO-13                   |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 14            | 10          | 71.43                         | 6.70E-02                      | 9.60E-02               | J          | 6.59E-01               | 2.37E-01               | 3.70E+00                 | J                                 | GEO-13                   |
| Naphthalene            | 91-20-3    | 14            | 2           | 14.29                         | 3.30E-02                      | 8.80E-02               | J          | 3.26E-02               | 2.20E-02               | 1.70E-01                 | J                                 | SS-10                    |
| Phenanthrene           | 85-01-8    | 14            | 8           | 57.14                         | 3.30E-02                      | 3.70E-02               | J          | 1.28E-01               | 5.30E-02               | 7.40E-01                 | J                                 | GEO-13                   |
| Pyrene                 | 129-00-0   | 14            | 12          | 85.71                         | 6.70E-02                      | 9.80E-02               | J          | 1.70E+00               | 4.60E-01               | 1.40E+01                 | J                                 | GEO-13                   |



**Table 3**  
**Statistical Summary and Selection of COPCs in EU2 Soil (0'-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>Unrestricted<br>Soil TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? |                          | Is the 95%<br>UCL > TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|---|---|--------------------------|--------------------------|
|                        |                  |                                 |                                   |  |   | Is the<br>Maximum<br>Detected ><br>TRG? | Is the 95%<br>UCL > TRG? |                          |
| Semivolatiles          |                  |                                 |                                   |  |   |   |                          |                          |
| 2-methylnaphthalene    | 4.95E-02         | 4.29E-02                        | Unknown                           | 4.29E-02                                 | 3.13E+63                                    | no                                      |                          |                          |
| Acenaphthene           | 2.29E-02         | 2.17E-02                        | Unknown                           | 2.17E-02                                 | 4.69E+03                                    | no                                      |                          |                          |
| Acenaphthylene         | 3.26E-01         | 4.99E-01                        | Unknown                           | 4.99E-01                                 | 4.69E+03                                    | no                                      |                          |                          |
| Anthracene             | 3.91E-01         | 6.29E-01                        | Unknown                           | 6.29E-01                                 | 2.35E+04                                    | no                                      |                          |                          |
| Benz(a)anthracene      | 1.74E+00         | 9.91E+00                        | Lognormal                         | 6.70E+00                                 | 8.75E-01                                    | YES                                     | YES - COPC               |                          |
| Benz(a)pyrene          | 1.50E+00         | 5.08E+06                        | Lognormal                         | 5.08E+00                                 | 8.75E-02                                    | YES                                     | YES - COPC               |                          |
| Benz(b)fluoranthene    | 3.08E+00         | 2.53E+01                        | Lognormal                         | 9.20E+00                                 | 8.75E-01                                    | YES                                     | YES - COPC               |                          |
| Benzogghiptycene       | 8.46E-01         | 2.74E+00                        | Lognormal                         | 2.30E+00                                 | 2.35E+03                                    | no                                      |                          |                          |
| Benz(k)fluoranthene    | 1.19E+00         | 2.93E+00                        | Lognormal                         | 2.93E+00                                 | 8.75E+00                                    | no                                      |                          |                          |
| Carbazole              | 1.12E-01         | 1.24E-01                        | Unknown                           | 1.24E-01                                 | 3.19E+01                                    | no                                      |                          |                          |
| Chrysene               | 2.22E+00         | 1.68E+01                        | Lognormal                         | 8.00E+00                                 | 8.75E+01                                    | no                                      |                          |                          |
| Dibenz(a,h)anthracene  | 3.11E-01         | 4.93E-01                        | Unknown                           | 4.93E-01                                 | 8.75E-01                                    | YES                                     | YES - COPC               |                          |
| Dibenzofuran           | 3.83E-02         | 3.57E-02                        | Unknown                           | 3.57E-02                                 | 3.13E+02                                    | no                                      |                          |                          |
| Di-n-butylphthalate    | 5.48E-02         | 6.30E-02                        | Normal/Lognormal                  | 6.30E-02                                 | 2.28E+03                                    | no                                      |                          |                          |
| Fluoranthene           | 2.89E+00         | 1.66E+01                        | Lognormal                         | 1.20E+01                                 | 3.13E+03                                    | no                                      |                          |                          |
| Fluorene               | 8.84E-02         | 5.84E-02                        | Unknown                           | 5.84E-02                                 | 3.13E+03                                    | no                                      |                          |                          |
| Indeno(1,2,3-cd)pyrene | 1.15E+00         | 4.29E+00                        | Lognormal                         | 3.70E+00                                 | 8.75E-01                                    | YES                                     | YES - COPC               |                          |
| Naphthalene            | 5.34E-02         | 4.71E-02                        | Unknown                           | 4.71E-02                                 | 6.45E+02                                    | no                                      |                          |                          |
| Phenanthrene           | 2.26E-01         | 3.96E-01                        | Lognormal                         | 3.96E-01                                 | 2.35E+03                                    | no                                      |                          |                          |
| Pyrene                 | 3.43E+00         | 1.25E+01                        | Lognormal                         | 1.25E+01                                 | 2.35E+03                                    | no                                      |                          |                          |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 4

Statistical Summary and Selection of COPCs in EU2 Soil (0-6' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit | Minimum Detected mg/kg | Mean mg/kg | Logarithmic Mean | Maximum Detected mg/kg | Detected Qualifier | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------|------------------------|------------|------------------|------------------------|--------------------|-----------------------------|-----------------------------------|--------------------------|
| Semi-volatiles             |            |               |                |                 |               |               |                 |                        |            |                  |                        |                    |                             |                                   |                          |
| 2-methylnaphthalene        | 91-57-6    | 20            | 2              | 10              | 3.30E-02      | 7.00E-02      | J               | 2.71E-02               | 2.07E-02   | 1.60E-01         | J                      | SS-10              | 3.34E-02                    |                                   |                          |
| Acenaphthene               | 83-32-9    | 21            | 1              | 4.76            | 3.30E-02      | 4.90E-02      | J               | 2.51E-02               | 2.01E-02   | 4.90E-02         | J                      | GEO-13             | 2.95E-02                    |                                   |                          |
| Acenaphthylene             | 208-96-8   | 21            | 6              | 28.57           | 3.30E-02      | 3.00E-01      | J               | 1.19E-01               | 3.60E-02   | 1.30E+00         | J                      | GEO-13             | 2.91E-01                    |                                   |                          |
| Anthracene                 | 120-12-7   | 21            | 8              | 38.1            | 3.30E-02      | 3.90E-02      | J               | 1.37E-01               | 3.94E-02   | 1.60E+00         | J                      | GEO-13             | 3.54E-01                    |                                   |                          |
| Benz(a)anthracene          | 56-55-3    | 21            | 14             | 66.67           | 3.30E-02      | 3.80E-02      | J               | 6.10E-01               | 1.12E-01   | 6.70E+00         | J                      | GEO-13             | 1.49E+00                    |                                   |                          |
| Benzo(a)pyrene             | 50-32-8    | 21            | 12             | 57.14           | 3.70E-02      | 6.70E-02      | J               | 5.65E-01               | 1.25E-01   | 5.20E+00         | J                      | GEO-13             | 1.21E+00                    |                                   |                          |
| Benzo(b)fluoranthene       | 205-99-2   | 21            | 16             | 76.19           | 3.70E-02      | 6.70E-02      | J               | 1.29E+00               | 3.16E-01   | 9.20E+00         | J                      | GEO-13             | 2.26E+00                    |                                   |                          |
| Benzo(g)perylene           | 191-24-2   | 21            | 11             | 52.38           | 3.70E-02      | 6.70E-02      | J               | 3.54E-01               | 1.04E-01   | 2.30E+00         | J                      | GEO-13             | 6.08E-01                    |                                   |                          |
| Benzo(k)fluoranthene       | 207-08-9   | 21            | 13             | 61.9            | 3.70E-02      | 1.30E-01      | J               | 5.21E-01               | 1.84E-01   | 3.60E+00         | J                      | GEO-13             | 8.79E-01                    |                                   |                          |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 20            | 1              | 5               | 6.70E-02      | 7.80E-02      | J               | 5.15E-02               | 3.91E-02   | 3.70E-01         | J                      | GEO-13             | 7.50E-02                    |                                   |                          |
| Carbazole                  | 86-74-8    | 20            | 4              | 20              | 3.30E-02      | 3.90E-02      | J               | 4.96E-02               | 2.57E-02   | 3.50E-01         | J                      | GEO-13             | 8.92E-02                    |                                   |                          |
| Chrysene                   | 218-01-9   | 21            | 13             | 61.9            | 3.30E-02      | 7.40E-02      | J               | 8.03E-01               | 1.32E-01   | 8.00E+00         | J                      | GEO-13             | 1.83E+00                    |                                   |                          |
| Dibenz(a,h)anthracene      | 53-70-3    | 21            | 8              | 38.1            | 3.70E-02      | 6.70E-02      | J               | 1.29E-01               | 5.30E-02   | 9.10E-01         | J                      | GEO-13             | 2.29E-01                    |                                   |                          |
| Dibenzofuran               | 132-64-9   | 20            | 2              | 10              | 3.30E-02      | 3.90E-02      | J               | 2.41E-02               | 2.02E-02   | 9.80E-02         | J                      | SS-10              | 2.13E-02                    |                                   |                          |
| Di-n-butylphthalate        | 84-74-2    | 20            | 9              | 45              | 3.30E-02      | 7.80E-02      | J               | 4.15E-02               | 3.71E-02   | 1.10E-01         | J                      | SS-10              | 2.08E-02                    |                                   |                          |
| Fluoranthene               | 206-44-0   | 21            | 14             | 66.67           | 3.30E-02      | 5.00E-02      | J               | 9.54E-01               | 1.42E-01   | 1.20E+01         | J                      | GEO-13             | 2.63E+00                    |                                   |                          |
| Floorene                   | 86-73-7    | 21            | 4              | 19.05           | 3.30E-02      | 3.80E-02      | J               | 5.08E-02               | 2.45E-02   | 3.70E-01         | J                      | GEO-13             | 9.99E-02                    |                                   |                          |
| Indeno[1,2,3-cd]pyrene     | 193-39-5   | 21            | 11             | 52.38           | 3.70E-02      | 6.70E-02      | J               | 4.50E-01               | 1.11E-01   | 3.70E+00         | J                      | GEO-13             | 8.86E-01                    |                                   |                          |
| Naphthalene                | 91-20-3    | 21            | 2              | 9.52            | 3.30E-02      | 3.00E-01      | J               | 3.43E-02               | 2.31E-02   | 1.70E-01         | J                      | SS-10              | 4.47E-02                    |                                   |                          |
| Phenanthrene               | 85-01-8    | 21            | 9              | 42.86           | 3.30E-02      | 3.90E-02      | J               | 1.01E-01               | 4.22E-02   | 7.40E-01         | J                      | GEO-13             | 1.77E-01                    |                                   |                          |
| Phenol                     | 108-95-2   | 20            | 2              | 10              | 3.30E-02      | 7.80E-02      | J               | 3.51E-02               | 2.52E-02   | 1.90E-01         | J                      | GEO-13             | 4.24E-02                    |                                   |                          |
| Pyrene                     | 129-00-0   | 21            | 14             | 66.67           | 3.70E-02      | 6.70E-02      | J               | 1.16E+00               | 1.92E-01   | 1.40E+01         | J                      | GEO-13             | 3.05E+00                    |                                   |                          |

Table 4

Statistical Summary and Selection of COPCs in EU2 Soil (0-6' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent                | Logarithmic Distribution |                  |                | Exposure Point Concentration<br>mg/kg | Tier 1 TRG<br>mg/kg | Is the Maximum Detected > TRG? |  |
|----------------------------|--------------------------|------------------|----------------|---------------------------------------|---------------------|--------------------------------|--|
|                            | 95% UCL<br>mg/kg         | 95% UCL<br>mg/kg | 99% Confidence |                                       |                     | Is the 95% UCL > TRG?          |  |
| <b>Semivolatiles</b>       |                          |                  |                |                                       |                     |                                |  |
| 2-methylnaphthalene        | 4.00E-02                 | 3.22E-02         | Unknown        | 3.22E-02                              | 8.18E+04            | no                             |  |
| Acenaphthene               | 3.62E-02                 | 2.90E-02         | Unknown        | 2.90E-02                              | 1.23E+05            | no                             |  |
| Acenaphthylene             | 2.28E-01                 | 1.83E-01         | Unknown        | 1.83E-01                              | 1.23E+05            | no                             |  |
| Anthracene                 | 2.70E-01                 | 2.09E-01         | Unknown        | 2.09E-01                              | 6.13E+05            | no                             |  |
| Benzo(a)anthracene         | 1.17E+00                 | 2.80E+00         | Lognormal      | 2.80E+00                              | 7.84E+00            | no                             |  |
| Benzo(a)pyrene             | 1.02E+00                 | 2.64E+00         | Lognormal      | 2.64E+00                              | 7.84E-01            | YES                            |  |
| Benzo(b)fluoranthene       | 2.14E+00                 | 1.09E+01         | Lognormal      | 9.20E+00                              | 7.84E+00            | YES                            |  |
| Benzo(g,h,i)perylene       | 5.83E-01                 | 1.41E+00         | Lognormal      | 1.41E+00                              | 6.13E+04            | no                             |  |
| Benzo(k)fluoranthene       | 8.52E-01                 | 1.84E+00         | Lognormal      | 1.84E+00                              | 7.84E+01            | no                             |  |
| Bis(2-ethylhexyl)phthalate | 8.05E-02                 | 5.77E-02         | Unknown        | 5.77E-02                              | 4.09E+02            | no                             |  |
| Carbazole                  | 8.41E-02                 | 6.51E-02         | Unknown        | 6.51E-02                              | 2.86E+02            | no                             |  |
| Chrysene                   | 1.49E+00                 | 5.33E+00         | Lognormal      | 5.33E+00                              | 7.84E+02            | no                             |  |
| Dibenz(a,h)anthracene      | 2.16E-01                 | 2.39E-01         | Unknown        | 2.39E-01                              | 7.84E-01            | YES                            |  |
| Dibenzofuran               | 3.23E-02                 | 2.86E-02         | Unknown        | 2.86E-02                              | 8.18E+03            | no                             |  |
| Di-n-butylphthalate        | 4.95E-02                 | 5.24E-02         | Lognormal      | 5.24E-02                              | 2.28E+03            | no                             |  |
| Fluoranthene               | 1.94E+00                 | 5.34E+00         | Lognormal      | 5.34E+00                              | 8.17E+04            | no                             |  |
| Fluorene                   | 8.84E-02                 | 6.16E-02         | Unknown        | 6.16E-02                              | 8.17E+04            | no                             |  |
| Indeno(1,2,3-cd)pyrene     | 7.83E-01                 | 1.97E+00         | Lognormal      | 1.97E+00                              | 7.84E+00            | no                             |  |
| Naphthalene                | 5.11E-02                 | 4.37E-02         | Unknown        | 4.37E-02                              | 8.24E+02            | no                             |  |
| Phenanthrene               | 1.67E-01                 | 1.88E-01         | Unknown        | 1.88E-01                              | 6.13E+04            | no                             |  |
| Phenol                     | 5.15E-02                 | 4.60E-02         | Unknown        | 4.60E-02                              | 1.23E+05            | no                             |  |
| Pyrene                     | 2.31E+00                 | 7.47E+00         | Lognormal      | 7.47E+00                              | 6.13E+04            | no                             |  |

Table 3

*Statistical Summary and Selection of COPCs in EU2 Soil ( $\theta\text{-}10'$  bgs)*  
*Kerr-McGee, Hattiesburg, MS*

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Maximum mg/kg | Detected | Maximum mg/kg | Detected | Maximum mg/kg | Qualifier | Logarithmic Mean mg/kg | Mean mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|---------------|----------|---------------|----------|---------------|-----------|------------------------|------------|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>       |            |               |                |                 |               |               |                       |          |               |               |          |               |          |               |           |                        |            |                                   |                          |
| 2,4-dimethylphenol         | 105-67-9   | 23            | 1              | 4.35            | 6.70E-02      | 3.30E-01      | 1.10E+00              | J        | 8.68E-02      | 4.33E-02      | J        | 1.10E+00      | J        | 1.10E+00      | J         | SB-05                  | 2.23E-01   |                                   |                          |
| 2-methylnaphthalene        | 91-57-6    | 23            | 4              | 17.39           | 3.30E-02      | 3.90E-02      | 7.00E-02              | J        | 2.71E-02      | 2.07E-02      | J        | 1.60E-01      | J        | SS-10         | 6.21E-01  |                        |            |                                   |                          |
| Acenaphthene               | 83-32-9    | 26            | 3              | 11.54           | 2.80E-02      | 3.10E-01      | 4.90E-02              | J        | 2.51E-02      | 2.01E-02      | J        | 4.90E-02      | J        | 4.90E-02      | J         | GEO-13                 | 4.29E+01   |                                   |                          |
| Acenaphthylene             | 208-96-8   | 26            | 8              | 30.77           | 2.80E-02      | 3.10E-01      | 3.70E-02              | J        | 1.19E-01      | 3.60E-02      | J        | 1.30E+00      | J        | 1.30E+00      | J         | GEO-13                 | 1.97E+06   |                                   |                          |
| Anthracene                 | 120-12-7   | 26            | 10             | 38.46           | 7.90E-04      | 3.90E-02      | 4.10E-02              | J        | 1.37E-01      | 3.94E-02      | J        | 1.60E+00      | J        | 1.60E+00      | J         | GEO-13                 | 2.86E+01   |                                   |                          |
| Benz(a)anthracene          | 56-55-3    | 26            | 16             | 61.54           | 8.60E-04      | 3.80E-02      | 4.10E-02              | J        | 6.10E-01      | 1.12E-01      | J        | 6.70E+00      | J        | 6.70E+00      | J         | GEO-13                 | 1.48E+01   |                                   |                          |
| Benzo(a)pyrene             | 50-32-8    | 26            | 16             | 61.54           | 3.70E-02      | 6.70E-02      | 1.69E-03              | J        | 5.65E-01      | 1.25E-01      | J        | 5.20E+00      | J        | 5.20E+00      | J         | GEO-13                 | 5.75E+00   |                                   |                          |
| Benzo(b)fluoranthene       | 205-99-2   | 26            | 20             | 76.92           | 3.70E-02      | 6.70E-02      | 7.60E-04              | J        | 1.29E+00      | 3.16E-01      | J        | 9.20E+00      | J        | 9.20E+00      | J         | GEO-13                 | 8.58E+00   |                                   |                          |
| Benzo(g,h)perylene         | 191-24-2   | 26            | 14             | 53.85           | 1.70E-02      | 6.70E-02      | 1.80E-03              | J        | 3.54E-01      | 1.04E-01      | J        | 2.30E+00      | J        | 2.30E+00      | J         | GEO-13                 | 1.65E+00   |                                   |                          |
| Benzo(k)fluoranthene       | 267-08-9   | 26            | 17             | 65.38           | 3.70E-02      | 1.30E-01      | 5.10E-04              | J        | 5.21E-01      | 1.84E-01      | J        | 3.60E+00      | J        | 3.60E+00      | J         | GEO-13                 | 2.70E+00   |                                   |                          |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 23            | 1              | 4.35            | 6.70E-02      | 5.00E-01      | 3.70E-01              | J        | 5.15E-02      | 3.91E-02      | J        | 3.70E-01      | J        | 3.70E-01      | J         | GEO-13                 | 8.41E-02   |                                   |                          |
| Carbazole                  | 86-74-8    | 23            | 6              | 26.09           | 3.30E-02      | 3.90E-02      | 4.30E-02              | J        | 4.96E-02      | 2.57E-02      | J        | 3.50E-01      | J        | 3.50E-01      | J         | GEO-13                 | 1.16E+01   |                                   |                          |
| Chrysene                   | 218-01-9   | 26            | 15             | 57.69           | 2.50E-03      | 7.40E-02      | 5.10E-02              | J        | 8.03E-01      | 1.32E-01      | J        | 8.00E+00      | J        | 8.00E+00      | J         | GEO-13                 | 1.31E+01   |                                   |                          |
| Dibenz(a,h)anthracene      | 53-70-3    | 26            | 10             | 38.46           | 5.30E-04      | 6.70E-02      | 1.88E-02              | J        | 1.29E-01      | 5.30E-02      | J        | 9.10E-01      | J        | 9.10E-01      | J         | GEO-13                 | 7.82E+01   |                                   |                          |
| Dibenzofuran               | 132-64-9   | 23            | 4              | 17.39           | 3.30E-02      | 3.90E-02      | 7.20E-02              | J        | 2.41E-02      | 2.02E-02      | J        | 9.80E-02      | J        | 9.80E-02      | J         | GEO-13                 | 4.47E+01   |                                   |                          |
| Di-n-hutylphthalate        | 84-74-2    | 23            | 9              | 39.13           | 3.30E-02      | 2.50E-01      | 3.60E-02              | J        | 4.15E-02      | 3.71E-02      | J        | 1.10E-01      | J        | 1.10E-01      | J         | SS-10                  | 2.80E-02   |                                   |                          |
| Fluoranthene               | 206-44-0   | 26            | 16             | 61.54           | 2.00E-03      | 3.80E-02      | 5.00E-02              | J        | 9.54E-01      | 1.42E-01      | J        | 1.20E+01      | J        | 1.20E+01      | J         | GEO-13                 | 5.91E+01   |                                   |                          |
| Fluorene                   | 86-73-7    | 26            | 6              | 23.08           | 2.60E-03      | 3.80E-02      | 2.90E-02              | J        | 5.08E-02      | 2.45E-02      | J        | 3.70E-01      | J        | 3.70E-01      | J         | GEO-13                 | 5.48E+01   |                                   |                          |
| Indeno[1,2,3-cd]pyrene     | 193-39-5   | 26            | 14             | 53.85           | 1.10E-02      | 6.70E-02      | 1.40E-03              | J        | 4.50E-01      | 1.11E-01      | J        | 3.70E+00      | J        | 3.70E+00      | J         | GEO-13                 | 2.30E+00   |                                   |                          |
| Naphthalene                | 91-20-3    | 26            | 5              | 19.23           | 2.80E-02      | 3.10E-01      | 8.80E-02              | J        | 3.43E-02      | 2.31E-02      | J        | 1.70E-01      | J        | 1.70E-01      | J         | SS-10                  | 8.54E+01   |                                   |                          |
| Phenanthrene               | 85-01-8    | 26            | 11             | 42.31           | 2.10E-03      | 3.90E-02      | 3.70E-02              | J        | 1.01E-01      | 4.22E-02      | J        | 7.40E-01      | J        | 7.40E-01      | J         | GEO-13                 | 1.20E+02   |                                   |                          |
| Phenol                     | 108-93-2   | 23            | 2              | 8.7             | 3.30E-02      | 2.50E-01      | 1.10E-01              | J        | 3.51E-02      | 2.52E-02      | J        | 1.90E-01      | J        | 1.90E-01      | J         | GEO-13                 | 4.49E+02   |                                   |                          |
| Pyrene                     | 129-00-0   | 26            | 16             | 61.54           | 4.50E-03      | 6.70E-02      | 6.80E-02              | J        | 1.16E+00      | 1.92E-01      | J        | 1.40E+01      | J        | 1.40E+01      | J         | GEO-13                 | 5.37E+01   |                                   |                          |
| <b>Volatiles</b>           |            |               |                |                 |               |               |                       |          |               |               |          |               |          |               |           |                        |            |                                   |                          |
| Acetone                    | 67-64-1    | 3             | 1              | 33.33           | 7.00E-03      | 3.50E-02      | 6.30E-02              | J        | 2.80E-02      | 1.57E-02      | J        | 6.30E-02      | J        | 6.30E-02      | J         | SB-05                  | 3.11E-02   |                                   |                          |
| Ethylbenzene               | 100-41-4   | 3             | 2              | 66.67           | 1.00E-03      | 1.00E-03      | 6.80E-02              | J        | 9.62E-02      | 1.96E-02      | J        | 2.20E-01      | J        | 2.20E-01      | J         | SB-05                  | 1.12E+01   |                                   |                          |
| Toluene                    | 108-88-3   | 3             | 2              | 66.67           | 1.00E-03      | 1.00E-03      | 1.40E-02              | J        | 2.07E-02      | 6.93E-03      | J        | 4.75E-02      | J        | 4.75E-02      | J         | SB-05                  | 2.42E+02   |                                   |                          |
| Xylene (total)             | 1330-20-7  | 3             | 2              | 66.67           | 1.00E-03      | 1.00E-03      | 4.90E-01              | J        | 5.64E-01      | 6.65E-02      | J        | 1.20E+00      | J        | 1.20E+00      | J         | SB-05                  | 6.03E+01   |                                   |                          |

Table 5

Statistical Summary and Selection of COPCs in EU2 Soil (0-10' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent                | Logarithmic      |                  |                                   | Exposure Point<br>Concentration<br>mg/kg | Tier 1<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? |            |
|----------------------------|------------------|------------------|-----------------------------------|--|------------------------|---|------------|
|                            | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence |  |                        | UCL > TRG?                              |            |
| <b>Semi-volatiles</b>      |                  |                  |                                   |  |                        |   |            |
| 2,4-dimethylphenol         | 1.66E+01         | 8.51E-02         | Unknown                           | 8.51E-02                                 | 4.08E+04               | no                                      |            |
| 2-methylnaphthalene        | 4.10E+01         | 3.97E+01         | Unknown                           | 1.60E-01                                 | 8.18E+04               | no                                      |            |
| Acenaphthene               | 2.58E+01         | 8.98E+00         | Unknown                           | 4.90E-02                                 | 1.23E+05               | no                                      |            |
| Acenaphthylene             | 1.33E+00         | 1.26E+00         | Unknown                           | 1.26E+00                                 | 1.23E+05               | no                                      |            |
| Anthracene                 | 1.77E+01         | 3.27E+01         | Unknown                           | 1.60E+00                                 | 6.13E+05               | no                                      |            |
| Benz(a)anthracene          | 9.66E+00         | 9.96E+01         | Lognormal                         | 6.70E+00                                 | 7.84E+00               | no                                      | YES*       |
| Benzo(a)pyrene             | 4.02E+00         | 2.17E+01         | Lognormal                         | 5.20E+00                                 | 7.84E+01               | YES                                     | YES - COPC |
| Benzo(b)fluoranthene       | 6.36E+00         | 1.42E+02         | Lognormal                         | 9.20E+00                                 | 7.84E+00               | YES                                     | YES - COPC |
| Benzo(g,h)perylene         | 1.30E+00         | 5.31E+00         | Lognormal                         | 2.30E+00                                 | 6.13E+04               | no                                      |            |
| Benzo(k)fluoranthene       | 2.08E+00         | 1.71E+01         | Lognormal                         | 3.60E+00                                 | 7.84E+01               | no                                      | YES*       |
| Bis(2-ethylhexyl)phthalate | 9.44E-02         | 7.81E-02         | Unknown                           | 7.81E-02                                 | 4.09E+02               | no                                      |            |
| Carbazole                  | 7.55E+00         | 5.44E+00         | Unknown                           | 3.50E-01                                 | 2.86E+02               | no                                      |            |
| Chrysene                   | 8.74E+00         | 7.90E+01         | Lognormal                         | 8.00E+00                                 | 7.84E+02               | no                                      | YES*       |
| Dibenz(a,h)anthracene      | 5.85E-01         | 1.69E+00         | Lognormal                         | 9.10E-01                                 | 7.84E+01               | YES                                     | YES**      |
| Dibenzofuran               | 2.99E+01         | 2.24E+01         | Unknown                           | 9.80E-02                                 | 8.18E+03               | no                                      |            |
| Di-n-butylphthalate        | 5.59E-02         | 5.95E-02         | Lognormal                         | 5.95E-02                                 | 2.28E+03               | no                                      |            |
| Fluoranthene               | 3.71E+01         | 4.36E+02         | Lognormal                         | 1.20E+01                                 | 8.17E+04               | no                                      |            |
| Fluorene                   | 3.32E+01         | 2.15E+01         | Unknown                           | 3.70E-01                                 | 8.17E+04               | no                                      |            |
| Indeno(1,2,3-cd)pyrene     | 1.78E+00         | 9.60E+00         | Lognormal                         | 3.70E+00                                 | 7.84E+00               | no                                      | YES*       |
| Naphthalene                | 5.71E+01         | 2.83E+01         | Unknown                           | 1.70E-01                                 | 8.24E+02               | no                                      |            |
| Phenanthrene               | 7.34E+01         | 1.11E+02         | Unknown                           | 7.40E-01                                 | 6.13E+04               | no                                      |            |
| Phenol                     | 5.65E-02         | 5.50E-02         | Unknown                           | 5.50E-02                                 | 1.23E+05               | no                                      |            |
| Pyrene                     | 3.39E+01         | 2.55E+02         | Lognormal                         | 1.40E+01                                 | 6.13E+04               | no                                      |            |
| <b>Volatiles</b>           |                  |                  |                                   |  |                        |   |            |
| Acetone                    | 8.04E-02         | 1.17E+07         | Normal/Lognormal                  | 6.30E-02                                 | 1.04E+05               | no                                      |            |
| Ethylbenzene               | 2.86E-01         | 2.68E+42         | Normal/Lognormal                  | 2.20E-01                                 | 3.95E+02               | no                                      |            |
| Toluene                    | 6.15E-02         | 2.26E+21         | Normal/Lognormal                  | 4.75E-02                                 | 3.80E+01               | no                                      |            |
| Xylene (total)             | 1.58E+00         | 3.97E+75         | Normal/Lognormal                  | 1.20E+00                                 | 3.18E+02               | no                                      |            |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

\*\*Logarithmic 95% UCL is less than benchmark but retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

**Table 6**  
**Statistical Summary and Selection of COPCs in EU3 Soil (0'-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | CAS Number | Total Samples | Hit % | Minimum Detection Limit mg/kg | Maximum Detection Limit mg/kg | Mean Qualifier mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Maximum Detected mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|-------|-------------------------------|-------------------------------|----------------------|------------------------|------------------------|------------------------|-----------------------------------|--------------------------|
| Semivolatiles          |            |               |       |                               |                               |                      |                        |                        |                        |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 3             | 1     | 33.33                         | 3.30E-02                      | 2.30E-01             | J                      | 8.77E-02               | 2.30E-01               | J                                 | SS-16                    |
| Acenaphthylene         | 208-96-8   | 3             | 2     | 66.67                         | 3.30E-02                      | 1.20E-01             | J                      | 1.02E-01               | 1.70E-01               | J                                 | SS-16                    |
| Anthracene             | 120-12-7   | 3             | 2     | 66.67                         | 3.30E-02                      | 1.20E-01             | J                      | 1.02E-01               | 6.96E-02               | 6.96E-02                          | 7.83E-02                 |
| Benz(a)anthracene      | 56-55-3    | 3             | 3     | 100                           | 0.00E+00                      | 0.00E+00             | J                      | 5.60E-02               | 3.62E-01               | 2.46E-01                          | 5.40E-01                 |
| Benz(a)pyrene          | 50-32-8    | 3             | 2     | 66.67                         | 6.70E-02                      | 6.70E-02             | J                      | 5.60E-01               | 4.35E-01               | 2.37E-01                          | 7.10E-01                 |
| Benz(b)fluoranthene    | 205-99-2   | 3             | 3     | 100                           | 0.00E+00                      | 0.00E+00             | J                      | 1.90E-01               | 9.30E-01               | 6.83E-01                          | 1.40E+00                 |
| Benz(gi)perylene       | 191-24-2   | 3             | 3     | 100                           | 0.00E+00                      | 0.00E+00             | J                      | 8.00E-02               | 6.53E-01               | 4.03E-01                          | 1.20E+00                 |
| Benzo(k)fluoranthene   | 207-08-9   | 3             | 2     | 66.67                         | 1.30E-01                      | 4.70E-01             | J                      | 3.42E-01               | 2.46E-01               | 4.90E-01                          | SS-16                    |
| Carbazole              | 86-74-8    | 3             | 2     | 66.67                         | 3.30E-02                      | 4.60E-02             | J                      | 5.75E-02               | 4.37E-02               | 1.10E-01                          | J                        |
| Chrysene               | 218-01-9   | 3             | 3     | 100                           | 0.60E+00                      | 0.00E+00             | J                      | 1.10E-01               | 5.93E-01               | 4.25E-01                          | 8.70E-01                 |
| Dibenz(a,h)anthracene  | 53-70-3    | 3             | 2     | 66.67                         | 6.70E-02                      | 1.40E-01             | J                      | 1.11E-01               | 9.09E-02               | 1.60E-01                          | J                        |
| Dibenzofuran           | 132-64-9   | 3             | 2     | 66.67                         | 3.30E-02                      | 3.30E-02             | J                      | 4.85E-02               | 3.81E-02               | 9.30E-02                          | J                        |
| Di-n-butyl phthalate   | 84-74-2    | 3             | 3     | 100                           | 0.00E+00                      | 0.00E+00             | J                      | 8.30E-02               | 7.58E-02               | 1.10E-01                          | J                        |
| Fluoranthene           | 206-44-0   | 3             | 3     | 100                           | 0.00E+00                      | 0.00E+00             | J                      | 1.20E-01               | 5.27E-01               | 3.99E-01                          | 7.80E-01                 |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 3             | 3     | 100                           | 0.00E+00                      | 0.00E+00             | J                      | 8.60E-02               | 3.85E-01               | 2.89E-01                          | 6.00E-01                 |
| Naphthalene            | 91-20-3    | 3             | 1     | 33.33                         | 3.30E-02                      | 3.30E-02             | J                      | 1.60E-01               | 6.43E-02               | 3.52E-02                          | 1.60E-01                 |
| Phenanthrene           | 85-01-8    | 3             | 2     | 66.67                         | 3.30E-02                      | 3.30E-02             | J                      | 1.30E-01               | 1.32E-01               | 8.12E-02                          | 2.50E-01                 |
| Pyrene                 | 129-00-0   | 3             | 3     | 100                           | 0.00E+00                      | 0.00E+00             | J                      | 1.20E-01               | 6.90E-01               | 4.85E-01                          | 1.00E+00                 |

**Table 6**  
**Statistical Summary and Selection of COPCs in EU3 Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | Logarithmic      |                  | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier 1<br>Unrestricted<br>Soil TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? | Is the<br>Logarithmic<br>95% UCL ><br>TRG? |
|------------------------|------------------|------------------|--------------------------------|--|---|---|--|
|                        | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg |                                |  |   |   |  |
| <b>Semivolatile</b>    |                  |                  |                                |  |   |   |  |
| 2-methylnaphthalene    | 2.95E-01         | 2.43E+08         | Unknown                        | 2.30E-01                                 | 3.13E+03                                    | no                                      |  |
| Acenaphthylene         | 2.34E-01         | 3.45E+05         | Normal/Lognormal               | 1.70E-01                                 | 4.69E+03                                    | no                                      |  |
| Anthracene             | 2.34E-01         | 3.45E+05         | Normal/Lognormal               | 1.70E-01                                 | 2.35E+04                                    | no                                      |  |
| Benz(a)anthracene      | 8.11E-01         | 2.15E+06         | Normal/Lognormal               | 5.40E-01                                 | 8.75E-01                                    | yes*                                    |  |
| Benz(a)pyrene          | 1.03E+00         | 3.82E+11         | Normal/Lognormal               | 7.10E-01                                 | 8.75E-02                                    | YES - COPC                              |  |
| Benz(b)fluoranthene    | 2.02E+00         | 1.13E+05         | Normal/Lognormal               | 1.40E+00                                 | 8.75E-01                                    | YES - COPC                              |  |
| Benz(ghi)perylene      | 1.60E+00         | 1.70E+08         | Normal/Lognormal               | 1.20E+00                                 | 2.35E+03                                    | no                                      |  |
| Benz(k)fluoranthene    | 7.46E-01         | 1.06E+05         | Normal/Lognormal               | 4.90E-01                                 | 3.75E+00                                    | yes*                                    |  |
| Carbazole              | 1.38E-01         | 2.81E+02         | Normal/Lognormal               | 1.10E-01                                 | 3.19E+01                                    | no                                      |  |
| Chrysene               | 1.30E+00         | 2.63E+05         | Normal/Lognormal               | 8.70E-01                                 | 8.75E+01                                    | yes*                                    |  |
| Dibenz(a,h)anthracene  | 2.26E-01         | 1.35E+02         | Normal/Lognormal               | 1.60E-01                                 | 8.75E-02                                    | YES - COPC                              |  |
| Dibenzofuran           | 1.16E-01         | 5.59E+01         | Normal/Lognormal               | 9.30E-02                                 | 3.13E+02                                    | no                                      |  |
| Di-n-butyl phthalate   | 1.46E-01         | 1.52E+00         | Normal/Lognormal               | 1.10E-01                                 | 2.28E+03                                    | no                                      |  |
| Fluoranthene           | 1.13E+00         | 1.59E+04         | Normal/Lognormal               | 7.80E-01                                 | 3.13E+03                                    | no                                      |  |
| Indeno(1,2,3-cd)pyrene | 8.36E-01         | 1.56E+04         | Normal/Lognormal               | 6.00E-01                                 | 8.75E-01                                    | no                                      |  |
| Naphthalene            | 2.04E-01         | 6.64E+05         | Unknown                        | 1.60E-01                                 | 6.45E+02                                    | no                                      |  |
| Phenanthrene           | 3.29E-01         | 2.65E+07         | Normal/Lognormal               | 2.50E-01                                 | 2.35E+03                                    | no                                      |  |
| Pyrene                 | 1.52E+00         | 7.45E+05         | Normal/Lognormal               | 1.00E+00                                 | 2.35E+03                                    | no                                      |  |

\*Retained as a COPC, as per MDDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table

*Statistical Summary and Selection of COPCs in EU3 Soil (0-6' bgs)*  
*Kerr McGee, Harrisburg, MS*

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected mg/kg | Minimum Qualifier mg/kg | Maximum Qualifier mg/kg | Logarithmic Mean | Mean mg/kg | Detected mg/kg | Detected Qualifier mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------------|-------------------------|-------------------------|------------------|------------|----------------|--------------------------|-----------------------------------|--------------------------|
| Semi挥发物                |            |               |                |                 |               |               |                       |                |                         |                         |                  |            |                |                          |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 7             | 1              | 14.29           | 3.30E-02      | 4.00E-02      | 2.30E-01              | J              | 4.85E-02                | 2.62E-02                | 2.30E-01         |            | J              | SS-16                    | 8.00E-02                          |                          |
| Acenaphthylene         | 208-96-8   | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 1.20E-01              | J              | 5.47E-02                | 3.33E-02                | 1.70E-01         |            | J              | SS-16                    | 6.34E-02                          |                          |
| Anthracene             | 120-12-7   | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 1.20E-01              | J              | 5.47E-02                | 3.33E-02                | 1.70E-01         |            | J              | SS-16                    | 6.34E-02                          |                          |
| Benzo(a)anthracene     | 56-55-3    | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 5.60E-02              | J              | 1.66E-01                | 5.71E-02                | 5.40E-01         |            | J              | SS-15                    | 2.39E-01                          |                          |
| Benzo(a)pyrene         | 50-32-8    | 7             | 2              | 28.57           | 3.70E-02      | 6.70E-02      | 5.60E-01              | J              | 1.97E-01                | 5.62E-02                | 7.10E-01         |            | J              | SS-16                    | 3.02E-01                          |                          |
| Benzo(b)fluoranthene   | 205-99-2   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 1.90E-01              | J              | 4.10E-01                | 8.85E-02                | 1.40E+00         |            | J              | SS-16                    | 6.14E-01                          |                          |
| Benzo(ghi)perylene     | 191-24-2   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 8.00E-02              | J              | 2.91E-01                | 7.06E-02                | 1.20E+00         |            | J              | SS-16                    | 4.69E-01                          |                          |
| Benzo(k)fluoranthene   | 207-08-9   | 7             | 2              | 28.57           | 3.70E-02      | 1.30E-01      | 4.70E-01              | J              | 1.57E-01                | 5.72E-02                | 4.90E-01         |            | J              | SS-16                    | 2.21E-01                          |                          |
| Carbazole              | 86-74-8    | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 4.60E-02              | J              | 3.56E-02                | 2.72E-02                | 1.10E-01         |            | J              | SS-15                    | 3.44E-02                          |                          |
| Chrysene               | 218-01-9   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 1.10E-01              | J              | 2.65E-01                | 7.22E-02                | 8.70E-01         |            | J              | SS-16                    | 3.91E-01                          |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 7             | 2              | 28.57           | 3.70E-02      | 6.70E-02      | 1.40E-01              | J              | 5.86E-02                | 3.73E-02                | 1.60E-01         |            | J              | SS-16                    | 6.29E-02                          |                          |
| Dibenzofuran           | 132-64-9   | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 3.60E-02              | J              | 3.17E-02                | 2.57E-02                | 9.30E-02         |            | J              | SS-16                    | 2.78E-02                          |                          |
| Di-n-butylphthalate    | 84-74-2    | 7             | 3              | 42.86           | 7.50E-02      | 7.90E-02      | 4.00E-02              | J              | 5.74E-02                | 5.13E-02                | 1.10E-01         |            | J              | SS-16                    | 3.23E-02                          |                          |
| Fluoranthene           | 206-44-0   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 1.20E-01              | J              | 2.37E-01                | 7.03E-02                | 7.80E-01         |            | J              | SS-16                    | 3.40E-01                          |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 8.60E-02              | J              | 1.76E-01                | 6.13E-02                | 6.00E-01         |            | J              | SS-16                    | 2.49E-01                          |                          |
| Naphthalene            | 91-20-3    | 7             | 1              | 14.29           | 3.30E-02      | 4.00E-02      | 1.60E-01              | J              | 3.85E-02                | 2.48E-02                | 1.60E-01         |            | J              | SS-16                    | 5.36E-02                          |                          |
| Phenanthrene           | 85-01-8    | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 1.30E-01              | J              | 6.76E-02                | 3.55E-02                | 2.50E-01         |            | J              | SS-16                    | 9.05E-02                          |                          |
| Phenol                 | 108-95-2   | 7             | 2              | 28.57           | 3.30E-02      | 7.90E-02      | 9.60E-02              | J              | 6.47E-02                | 3.95E-02                | 2.30E-01         |            | J              | GFO-17                   | 7.81E-02                          |                          |
| Pyrene                 | 129-00-0   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 1.20E-01              | J              | 3.07E-01                | 7.64E-02                | 1.00E+00         |            | J              | SS-17                    | 4.58E-01                          |                          |

**Table 7**  
**Statistical Summary and Selection of COPCs in EU3 Soil (0-6' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|------------------------|---|
| Semivolatiles          |                  |                                 |                                   |  |                        |   |
| 2-methylnaphthalene    | 1.07E-01         | 1.70E-01                        | Unknown                           | 1.70E-01                                 | 8.18E+04               | no                                      |
| Acenaphthylene         | 1.01E-01         | 2.50E-01                        | Unknown                           | 1.70E-01                                 | 1.23E+05               | no                                      |
| Anthraene              | 1.01E-01         | 2.50E-01                        | Unknown                           | 1.70E-01                                 | 6.13E+05               | no                                      |
| Benz(a)anthracene      | 3.42E-01         | 5.64E+00                        | Unknown                           | 5.40E-01                                 | 7.84E+00               | no                                      |
| Benz(e)pyrene          | 4.19E-01         | 1.08E+01                        | Unknown                           | 7.10E-01                                 | 7.84E-01               | no                                      |
| Benzo(b)fluoranthene   | 8.61E-01         | 1.83E+02                        | Lognormal                         | 1.40E+00                                 | 7.84E+00               | no                                      |
| Benzo(g)perylene       | 6.35E-01         | 3.79E+01                        | Lognormal                         | 1.20E+00                                 | 6.13E+04               | no                                      |
| Benzo(k)fluoranthene   | 3.20E-01         | 4.72E+00                        | Lognormal                         | 4.90E-01                                 | 7.84E+01               | no                                      |
| Carbazole              | 6.08E-02         | 8.08E+02                        | Unknown                           | 8.08E-02                                 | 2.86E+02               | no                                      |
| Chrysene               | 5.52E-01         | 3.06E+01                        | Lognormal                         | 8.70E-01                                 | 7.84E+02               | no                                      |
| Dibenz(a,h)anthracene  | 1.05E-01         | 2.52E-01                        | Unknown                           | 1.60E-01                                 | 7.84E-01               | no                                      |
| Dibenzofuran           | 5.21E-02         | 6.21E-02                        | Unknown                           | 6.21E-02                                 | 8.18E+03               | no                                      |
| Di-n-butylphthalate    | 8.12E-02         | 9.38E-02                        | Unknown                           | 9.38E-02                                 | 2.28E+03               | no                                      |
| Fluoranthene           | 4.87E-01         | 2.06E+01                        | Lognormal                         | 7.80E-01                                 | 8.17E+04               | no                                      |
| Indeno(1,2,3-cd)pyrene | 3.59E-01         | 6.92E+00                        | Lognormal                         | 6.00E-01                                 | 7.84E+00               | no                                      |
| Naphthalene            | 7.79E-02         | 1.03E-01                        | Unknown                           | 1.03E-01                                 | 8.24E+02               | no                                      |
| Phenanthrene           | 1.34E-01         | 4.34E-01                        | Unknown                           | 2.50E-01                                 | 6.13E+04               | no                                      |
| Pheno                  | 1.22E-01         | 3.11E-01                        | Lognormal                         | 2.30E-01                                 | 1.23E+05               | no                                      |
| Pyrene                 | 6.43E-01         | 5.32E+01                        | Lognormal                         | 1.00E+00                                 | 6.13E+04               | no                                      |

Table 6

Statistical Summary and Selection of COPCs in EU3 Soil (0-20' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected mg/kg | Minimum mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected mg/kg | Detected Qualifier | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------------|---------------|------------|------------------------|---------------|----------------|--------------------|-----------------------------|-----------------------------------|--------------------------|
| Semivolatiles          |            |               |                |                 |               |               |                       |                |               |            |                        |               |                |                    |                             |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 7             | 1              | 14.29           | 3.30E-02      | 4.00E-02      | 2.30E-01              | J              | 4.85E-02      | 2.62E-02   | 2.30E-01               | J             | SS-16          | 8.00E-02           |                             |                                   |                          |
| Acenaphthylene         | 208-96-8   | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 1.20E-01              | J              | 5.47E-02      | 3.33E-02   | 1.70E-01               | J             | SS-16          | 6.34E-02           |                             |                                   |                          |
| Anthracene             | 120-12-7   | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 1.20E-01              | J              | 5.47E-02      | 3.33E-02   | 1.70E-01               | J             | SS-16          | 6.34E-02           |                             |                                   |                          |
| Benz(a)anthracene      | 56-55-3    | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 5.60E-02              | J              | 1.66E-01      | 5.71E-02   | 5.40E-01               | J             | SS-15          | 2.39E-01           |                             |                                   |                          |
| Benz(a)pyrene          | 50-32-8    | 7             | 2              | 28.57           | 3.70E-02      | 6.70E-02      | 5.60E-01              | J              | 1.97E-01      | 5.62E-02   | 7.10E-01               | J             | SS-16          | 3.02E-01           |                             |                                   |                          |
| Benz(b)fluoranthene    | 205-99-2   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 1.90E-01              | J              | 4.10E-01      | 8.85E-02   | 1.40E+00               | J             | SS-16          | 6.14E-01           |                             |                                   |                          |
| Benz(g,h)perylene      | 191-24-2   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 8.00E-02              | J              | 2.91E-01      | 7.06E-02   | 1.20E+00               | J             | SS-16          | 4.69E-01           |                             |                                   |                          |
| Benz(k)fluoranthene    | 207-08-9   | 7             | 2              | 28.57           | 3.70E-02      | 1.30E-01      | 4.70E-01              | J              | 1.57E-01      | 5.72E-02   | 4.90E-01               | J             | SS-16          | 2.21E-01           |                             |                                   |                          |
| Carbazole              | 86-74-8    | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 4.60E-02              | J              | 3.56E-02      | 2.72E-02   | 1.10E-01               | J             | SS-15          | 3.44E-02           |                             |                                   |                          |
| Chrysene               | 218-01-9   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 1.10E-01              | J              | 2.65E-01      | 7.22E-02   | 8.70E-01               | J             | SS-16          | 3.91E-01           |                             |                                   |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 7             | 2              | 28.57           | 3.70E-02      | 6.70E-02      | 1.40E-01              | J              | 5.86E-02      | 3.73E-02   | 1.60E-01               | J             | SS-16          | 6.29E-02           |                             |                                   |                          |
| Dibenzofuran           | 132-64-9   | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 3.60E-02              | J              | 3.17E-02      | 2.57E-02   | 9.30E-02               | J             | SS-16          | 2.78E-02           |                             |                                   |                          |
| Di-n-butylphthalate    | 84-74-2    | 7             | 3              | 42.86           | 7.50E-02      | 7.90E-02      | 4.00E-02              | J              | 5.74E-02      | 5.13E-02   | 1.10E-01               | J             | SS-16          | 3.23E-02           |                             |                                   |                          |
| Fluoranthene           | 206-44-6   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 1.20E-01              | J              | 2.37E-01      | 7.03E-02   | 7.80E-01               | J             | SS-16          | 3.40E-01           |                             |                                   |                          |
| Indeno(1,2,3-cd)pyrene | 191-39-5   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 8.60E-02              | J              | 1.76E-01      | 6.13E-02   | 6.00E-01               | J             | SS-16          | 2.49E-01           |                             |                                   |                          |
| Naphthalene            | 91-20-3    | 7             | 1              | 14.29           | 3.30E-02      | 4.00E-02      | 1.60E-01              | J              | 3.85E-02      | 2.48E-02   | 1.60E-01               | J             | SS-16          | 5.36E-02           |                             |                                   |                          |
| Phenanthrene           | 85-01-8    | 7             | 2              | 28.57           | 3.30E-02      | 4.00E-02      | 1.30E-01              | J              | 6.76E-02      | 3.55E-02   | 2.50E-01               | J             | SS-16          | 9.05E-02           |                             |                                   |                          |
| Phenol                 | 108-95-2   | 7             | 2              | 28.57           | 3.30E-02      | 7.90E-02      | 9.60E-02              | J              | 6.47E-02      | 3.95E-02   | 2.30E-01               | J             | GEO-17         | 7.81E-02           |                             |                                   |                          |
| Pyrene                 | 123-00-0   | 7             | 3              | 42.86           | 3.70E-02      | 4.00E-02      | 1.20E-01              | J              | 3.07E-01      | 7.64E-02   | 1.00E+00               | J             | SS-17          | 4.58E-01           |                             |                                   |                          |

**Table 8**  
**Statistical Summary and Selection of COPCs in EU3 Soil (0-20' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent<br>Semivolatiles | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>Restricted Soil<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? |
|------------------------------|------------------|---------------------------------|-----------------------------------|--|---|---|
| 2-methylnaphthalene          | 1.07E-01         | 1.70E-01                        | Unknown                           | 1.70E-01                                 | 8.18E+04                                  | no                                      |
| Acenaphthylene               | 1.01E-01         | 2.50E-01                        | Unknown                           | 1.70E-01                                 | 1.23E+05                                  | no                                      |
| Anthracene                   | 1.01E-01         | 2.50E-01                        | Unknown                           | 1.70E-01                                 | 6.13E+05                                  | no                                      |
| Benz(a)anthracene            | 3.42E-01         | 5.64E+00                        | Unknown                           | 5.40E-01                                 | 7.84E+00                                  | no                                      |
| Benzo(b)pyrene               | 4.19E-01         | 1.08E+01                        | Unknown                           | 7.10E-01                                 | 7.84E-01                                  | no                                      |
| Benzo(b)fluoranthene         | 8.61E-01         | 1.83E+02                        | Lognormal                         | 1.40E+00                                 | 7.84E+00                                  | no                                      |
| Benzo(g,h)perylene           | 6.35E-01         | 3.79E+01                        | Lognormal                         | 1.20E+00                                 | 6.13E+04                                  | no                                      |
| Benzo(k)fluoranthene         | 3.20E-01         | 4.72E+00                        | Lognormal                         | 4.90E-01                                 | 7.84E+01                                  | no                                      |
| Carbazole                    | 6.08E-02         | 8.08E-02                        | Unknown                           | 8.08E-02                                 | 2.86E+02                                  | no                                      |
| Chrysene                     | 5.52E-01         | 3.06E+01                        | Lognormal                         | 8.70E-01                                 | 7.84E+02                                  | no                                      |
| Dibenz(a,h)anthracene        | 1.05E-01         | 2.52E-01                        | Unknown                           | 1.60E-01                                 | 7.84E-01                                  | no                                      |
| Dibenzofuran                 | 5.21E-02         | 6.21E-02                        | Unknown                           | 6.21E-02                                 | 8.18E-03                                  | no                                      |
| Di-n-butylphthalate          | 8.12E-02         | 9.38E-02                        | Unknown                           | 9.38E-02                                 | 2.28E+03                                  | no                                      |
| Fluoranthene                 | 4.87E-01         | 2.06E+01                        | Lognormal                         | 7.80E-01                                 | 8.17E+04                                  | no                                      |
| Indeno(1,2,3-cd)pyrene       | 3.59E-01         | 6.92E+00                        | Lognormal                         | 6.00E-01                                 | 7.84E+00                                  | no                                      |
| Naphthalene                  | 7.79E-02         | 1.03E-01                        | Unknown                           | 1.03E-01                                 | 8.24E+02                                  | no                                      |
| Phenanthrene                 | 1.34E-01         | 4.34E-01                        | Unknown                           | 2.50E-01                                 | 6.13E+04                                  | no                                      |
| Phenol                       | 1.22E-01         | 3.11E-01                        | Lognormal                         | 2.30E-01                                 | 1.23E+05                                  | no                                      |
| Pyrene                       | 6.43E-01         | 5.32E+01                        | Lognormal                         | 1.00E+00                                 | 6.13E+04                                  | no                                      |

*Statistical Summary and Selection of COPCs in EU4 Sediment  
Kerr McGee, Hattiesburg, MS*

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected mg/kg | Minimum Detected mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Detected Qualifier | Maximum Qualifier mg/kg | Detected Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------------|------------------------|------------|------------------------|------------------------|--------------------|-------------------------|------------------------------|-----------------------------------|--------------------------|
| <b>Semivolatile</b>    |            |               |                |                 |               |               |                       |                |                        |            |                        |                        |                    |                         |                              |                                   |                          |
| 2,4-dimethylphenol     | 105-67-9   | 1             | 1              | 100             | 0.00E+00      | 1.50E+00      | J                     | 1.50E+00       | 1.50E+00               | 1.50E+00   | 1.50E+00               | J                      | SD-02              | NA                      | NA                           | NA                                |                          |
| 2-methylnaphthalene    | 91-57-6    | 1             | 1              | 100             | 0.00E+00      | 1.50E+03      | J                     | 1.50E+03       | 1.50E+03               | 1.50E+03   | 1.50E+03               | J                      | SD-02              | NA                      | NA                           | NA                                |                          |
| Acenaphthene           | 83-32-9    | 8             | 2              | 25              | 5.70E-01      | 1.00E+03      | 6.50E+00              | J              | 1.08E+02               | 6.96E+00   | 3.45E+02               | J                      | SD-12              | 1.99E+02                |                              |                                   |                          |
| Acenaphthylene         | 208-96-8   | 8             | 1              | 12.5            | 5.70E-01      | 1.85E+02      | 3.50E+01              | J              | 1.72E+01               | 3.57E+00   | 3.50E+01               | J                      | SD-02              | 3.26E+01                |                              |                                   |                          |
| Anthracene             | 120-12-7   | 8             | 4              | 50              | 3.80E-02      | 5.56E-01      | 1.80E+00              | Z              | 4.01E+02               | 1.78E+00   | 1.90E+03               | Z                      | SD-02              | 7.57E+02                |                              |                                   |                          |
| Benzof[a]anthracene    | 56-55-3    | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.29E-01              | Z              | 6.52E+01               | 4.98E+00   | 3.30E+02               | Z                      | SD-02              | 1.23E+02                |                              |                                   |                          |
| Benzof[a]pyrene        | 50-32-8    | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.80E-01              | Z              | 2.69E+01               | 4.43E+00   | 1.30E+02               | J                      | SD-02              | 4.81E+01                |                              |                                   |                          |
| Benzof[b]fluoranthene  | 205-99-2   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 4.25E-01              | Z              | 3.29E+01               | 6.00E+00   | 1.80E+02               | J                      | SD-02              | 6.31E+01                |                              |                                   |                          |
| Benzof[ghi]perylene    | 191-24-2   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 1.73E-01              | J              | 7.24E+00               | 2.05E+00   | 3.60E+01               | J                      | SD-02              | 1.25E+01                |                              |                                   |                          |
| Benzof[k]fluoranthene  | 207-08-9   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.13E-01              | Z              | 1.38E+01               | 2.88E+00   | 6.40E+01               | Z                      | SD-02              | 2.38E+01                |                              |                                   |                          |
| Carbazole              | 86-74-8    | 1             | 1              | 100             | 0.00E+00      | 0.00E+00      | 5.90E+02              | Z              | 5.90E+02               | 5.90E+02   | 5.90E+02               | Z                      | SD-02              | NA                      | NA                           | NA                                | NA                       |
| Chrysene               | 218-01-9   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.50E-01              | Z              | 5.44E+01               | 4.85E+00   | 2.90E+02               | Z                      | SD-02              | 1.05E+02                |                              |                                   |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 8             | 7              | 87.5            | 6.00E-02      | 6.00E-02      | 5.70E-02              | J              | 2.93E+00               | 5.99E-01   | 1.20E+01               | J                      | SD-02              | 4.76E+00                |                              |                                   |                          |
| Dibenzofuran           | 132-64-9   | 1             | 1              | 100             | 0.00E+00      | 0.00E+00      | 9.40E+02              | Z              | 9.40E+02               | 9.40E+02   | 9.40E+02               | Z                      | SD-02              | NA                      | NA                           | NA                                | NA                       |
| Fluoranthene           | 206-44-0   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.60E-01              | Z              | 3.27E+02               | 1.42E+01   | 1.60E+03               | Z                      | SD-02              | 6.16E+02                |                              |                                   |                          |
| Fluorene               | 86-73-7    | 8             | 3              | 37.5            | 5.30E-02      | 4.50E-01      | 7.40E+00              | Z              | 2.32E+02               | 1.93E+00   | 1.20E+03               | Z                      | SD-02              | 4.51E+02                |                              |                                   |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.23E-01              | J              | 1.08E+01               | 2.98E+00   | 4.70E+01               | J                      | SD-02              | 1.75E+01                |                              |                                   |                          |
| Naphthalene            | 91-20-3    | 8             | 2              | 25              | 5.70E-01      | 1.85E+02      | 8.20E+00              | J              | 3.89E+02               | 7.60E+00   | 3.00E+03               | J                      | SD-02              | 1.06E+03                |                              |                                   |                          |
| Phenanthrene           | 85-01-8    | 8             | 3              | 37.5            | 3.10E-02      | 1.05E+00      | 2.38E+01              | Z              | 6.49E+02               | 3.49E+00   | 3.20E+03               | Z                      | SD-02              | 1.24E+03                |                              |                                   |                          |
| Pyrene                 | 129-00-0   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 4.59E-01              | J              | 2.48E+02               | 1.67E+01   | 1.00E+03               | J                      | SD-02              | 4.44E+02                |                              |                                   |                          |

**Table 9**  
**Statistical Summary and Selection of COPCs in EU4 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>99% UCL |                  |            | Exposure Point<br>Concentration<br>mg/kg | Tier 1<br>Unrestricted<br>Soil TRG<br>mg/kg | Maximum<br>Detected ><br>TRG? | Is the 95%<br>UCL ><br>TRG? |
|------------------------|------------------|------------------------|------------------|------------|--|---|-------------------------------|-----------------------------|
|                        |                  | 95% UCL<br>mg/kg       | 99% UCL<br>mg/kg | Confidence |  |   |                               |                             |
| <b>Semi挥发物</b>         |                  |                        |                  |            |  |   |                               |                             |
| 2,4-dimethylphenol     | NA               | NA                     | Unknown          | 1.30E+00   | 1.56E+03                                 | no  |                               |                             |
| 2-methylnaphthalene    | NA               | NA                     | Unknown          | 1.50E+03   | 3.13E+03                                 | no  |                               |                             |
| Acenaphthene           | 2.41E+02         | 8.24E+05               | Lognormal        | 3.45E+02   | 4.69E+03                                 | no  |                               |                             |
| Acenaphthylene         | 3.90E+01         | 1.12E+03               | Lognormal        | 3.50E+01   | 4.69E+03                                 | no  |                               |                             |
| Anthracene             | 9.08E+02         | 1.74E+15               | Lognormal        | 1.90E+03   | 2.35E+04                                 | no  |                               |                             |
| Benz(a)anthracene      | 1.48E+02         | 3.91E+05               | Lognormal        | 3.30E+02   | 8.75E-01                                 | YES   |                               |                             |
| Benz(a)pyrene          | 5.91E+01         | 6.94E+03               | Lognormal        | 1.30E+02   | 8.75E-02                                 | YES   |                               |                             |
| Benz(b)fluoranthene    | 7.52E+01         | 4.80E+03               | Lognormal        | 1.80E+02   | 8.75E-01                                 | YES   |                               |                             |
| Benz(ghi)perylene      | 1.56E+01         | 2.97E+02               | Lognormal        | 3.60E+01   | 2.35E+03                                 | no  |                               |                             |
| Benzo(k)fluoranthene   | 2.98E+01         | 1.74E+03               | Lognormal        | 6.40E+01   | 8.75E+00                                 | YES   |                               |                             |
| Carbazole              | NA               | NA                     | Unknown          | 5.90E+02   | 3.19E+01                                 | YES   |                               |                             |
| Chrysene               | 1.25E+02         | 1.18E+05               | Lognormal        | 2.90E+02   | 8.75E+01                                 | YES   |                               |                             |
| Dibenz(a,h)anthracene  | 6.12E+00         | 9.77E+02               | Lognormal        | 1.20E+01   | 8.75E-02                                 | YES   |                               |                             |
| Dibenzofuran           | NA               | NA                     | Unknown          | 9.40E+02   | 3.13E+02                                 | YES   |                               |                             |
| Fluoranthene           | 7.40E+02         | 6.03E+07               | Lognormal        | 1.60E+03   | 3.13E+03                                 | no  |                               |                             |
| Fluorene               | 5.34E+02         | 1.26E+12               | Lognormal        | 1.20E+03   | 3.13E+03                                 | no  |                               |                             |
| Indeno(1,2,3-cd)pyrene | 2.26E+01         | 5.97E+02               | Lognormal        | 4.70E+01   | 8.75E-01                                 | YES   |                               |                             |
| Naphthalene            | 1.10E+03         | 8.93E+06               | Lognormal        | 3.00E+03   | 6.45E+02                                 | YES   |                               |                             |
| Phenanthrene           | 1.48E+03         | 1.30E+15               | Lognormal        | 3.20E+03   | 2.35E+03                                 | YES   |                               |                             |
| Pyrene                 | 5.46E+02         | 5.33E+06               | Lognormal        | 1.00E+03   | 2.35E+03                                 | no  |                               |                             |

*Statistical Summary and Selection of COPCs in EU4 Surface Water  
Kerr McGee, Harrisburg, MS*

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/L | Maximum mg/L | Detection Limit mg/L | Detected | Minimum mg/L | Maximum mg/L | Detected | Detected | Logarithmic Mean mg/L | Mean mg/L | Maximum mg/L | Detected | Detected | Maximum mg/L | Location of Qualifier | Maximum Concentration mg/L | Standard Deviation mg/L |
|----------------------------|------------|---------------|----------------|-----------------|--------------|--------------|----------------------|----------|--------------|--------------|----------|----------|-----------------------|-----------|--------------|----------|----------|--------------|-----------------------|----------------------------|-------------------------|
| <b>Semivaluates</b>        |            |               |                |                 |              |              |                      |          |              |              |          |          |                       |           |              |          |          |              |                       |                            |                         |
| Acenaphthene               | 83-32-9    | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 1.40E-02             | 1.40E-02 | 1.40E-02     | 1.40E-02     | J        | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Anthracene                 | 120-12-7   | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 1.30E-02             | 1.30E-02 | 1.30E-02     | 1.30E-02     | NA       | SW-02    | 0.00E+00              |           |              |          |          |              |                       |                            |                         |
| Benzo(a)anthracene         | 56-55-3    | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 5.00E-03             | 5.00E-03 | 5.00E-03     | 5.00E-03     | J        | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Benzo(a)pyrene             | 50-32-8    | 1             | 0              | 100             | 1.00E-03     | 1.00E-03     | 0.00E+00             | 0.00E+00 | 5.00E-04     | 5.00E-04     | NA       | SW-02    | 0.00E+00              |           |              |          |          |              |                       |                            |                         |
| Benzo(b)fluoranthene       | 205-99-2   | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 1.20E-02             | 1.20E-02 | 1.20E-02     | 1.20E-02     | NA       | SW-02    | 0.00E+00              |           |              |          |          |              |                       |                            |                         |
| Benzo(k)fluoranthene       | 207-08-9   | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 2.00E-03             | 2.00E-03 | 2.00E-03     | 2.00E-03     | J        | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 3.00E-03             | 3.00E-03 | 3.00E-03     | 3.00E-03     | J        | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Carbazole                  | 86-74-8    | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 1.00E-02             | 1.00E-02 | 1.00E-02     | 1.00E-02     | J        | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Chrysene                   | 218-01-9   | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 6.00E-03             | 6.00E-03 | 6.00E-03     | 6.00E-03     | J        | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Dibenz(a,h)anthracene      | 53-70-3    | 1             | 0              | 100             | 1.00E-03     | 1.00E-03     | 0.00E+00             | 0.00E+00 | 5.00E-04     | 5.00E-04     | NA       | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Dibenzofuran               | 132-64-9   | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 1.10E-02             | 1.10E-02 | 1.10E-02     | 1.10E-02     | NA       | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Fluoranthene               | 206-44-0   | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 3.90E-02             | 3.90E-02 | 3.90E-02     | 3.90E-02     | NA       | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Fluorene                   | 86-73-7    | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 1.20E-02             | 1.20E-02 | 1.20E-02     | 1.20E-02     | NA       | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Indeno(1,2,3-cd)pyrene     | 193-39-5   | 1             | 0              | 100             | 1.00E-03     | 1.00E-03     | 0.00E+00             | 0.00E+00 | 5.00E-04     | 5.00E-04     | NA       | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Phenanthrene               | 85-01-8    | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 1.70E-02             | 1.70E-02 | 1.70E-02     | 1.70E-02     | NA       | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |
| Pyrene                     | 129-00-0   | 1             | 1              | 100             | 0.00E+00     | 0.00E+00     | 2.10E-02             | 2.10E-02 | 2.10E-02     | 2.10E-02     | NA       | NA       | SW-02                 | 0.00E+00  |              |          |          |              |                       |                            |                         |

NA - Not applicable: constituent not detected in media.



**Table 10**  
**Statistical Summary and Selection of COPCs in EU4 Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                 | 95% UCL<br>mg/L | Logarithmic<br>95% UCL<br>mg/L | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/L | Human Health                                     |                      | Is Maximum<br>Detected ><br>AWQC? |
|-----------------------------|-----------------|--------------------------------|--------------------------------|---|--|----------------------|-----------------------------------|
|                             |                 |                                |                                |   | Consumption of Water<br>& Organisms AWQC<br>mg/L | Human Health<br>mg/L |                                   |
| <b>Semivolatile</b>         |                 |                                |                                |   |  |                      |                                   |
| Acenaphthene                | NA              | NA                             | Unknown                        | 1.40E-02                                | 1.20E+00   | no                   |                                   |
| Anthracene                  | NA              | NA                             | Unknown                        | 1.30E-02                                | 9.60E+00   | no                   |                                   |
| Benzo(a)anthracene          | NA              | NA                             | Unknown                        | 5.00E-03                                | 4.40E-06   | YES - COPC           |                                   |
| Benzo(a)pyrene              | NA              | NA                             | Unknown                        | 5.00E-04                                | 4.40E-06   | YES**                |                                   |
| Benzo(b)fluoranthene        | NA              | NA                             | Unknown                        | 1.20E-02                                | 4.40E-06   | YES - COPC           |                                   |
| Benzo(k)fluoranthene        | NA              | NA                             | Unknown                        | 2.00E-03                                | 4.40E-06   | YES - COPC           |                                   |
| Benz(2-ethylhexyl)phthalate | NA              | NA                             | Unknown                        | 3.00E-03                                | 1.80E-03   | YES - COPC           |                                   |
| Carbazole*                  | NA              | NA                             | Unknown                        | 1.00E-02                                | NA   | no                   |                                   |
| Chrysene                    | NA              | NA                             | Unknown                        | 6.00E-03                                | 4.40E-06   | YES - COPC           |                                   |
| Dibenz(a,h)anthracene       | NA              | NA                             | Unknown                        | 5.00E-04                                | 4.40E-06   | YES**                |                                   |
| Dibenzofuran*               | NA              | NA                             | Unknown                        | 1.10E-02                                | NA   | no                   |                                   |
| Fluoranthene                | NA              | NA                             | Unknown                        | 3.90E-02                                | 3.00E-01   | no                   |                                   |
| Fluorene                    | NA              | NA                             | Unknown                        | 1.20E-02                                | 1.30E+00   | no                   |                                   |
| Indeno(1,2,3-cd)pyrene      | NA              | NA                             | Unknown                        | 5.00E-04                                | 4.40E-06   | YES**                |                                   |
| Phenanthrene*               | NA              | NA                             | Unknown                        | 1.70E-02                                | NA   | no                   |                                   |
| Pyrene                      | NA              | NA                             | Unknown                        | 2.10E-02                                | 9.60E-01   | no                   |                                   |

NA - Not Available

\* Constituent will be retained as a COPC due to lack of published screening criteria.

\*\*Retained as a COPC, as per MDEQ Comments (8/2/2000): constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 11

Statistical Summary and Selection of COPCs in EU4 Soil (#-I' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Maximum mg/kg | Mean mg/kg | Logarithmic Mean | Detected | Maximum mg/kg | Mean mg/kg | Qualifier | Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|---------------|------------|------------------|----------|---------------|------------|-----------|---------------------|-----------------------------------|--------------------------|
| <b>SemiVolatile</b>    |            |               |                |                 |               |               |                       |          |               |               |            |                  |          |               |            |           |                     |                                   |                          |
| 2,4-dimethylphenol     | 105-67-9   | 3             | 1              | 33.33           | 4.10E-01      | 9.90E+00      | 2.50E-01              | J        | 1.80E+00      | 6.33E-01      | 2.50E-01   | 2.50E-01         | J        | GEO-19        | 2.73E+00   |           |                     |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 2.70E-01              | J        | 9.36E-01      | 3.42E+00      | 2.80E+02   | J                | GEO-21   | 1.61E+02      |            |           |                     |                                   |                          |
| 2-methylphenol         | 95-48-7    | 3             | 1              | 33.33           | 2.00E-01      | 5.00E+00      | 7.30E-02              | J        | 8.91E-01      | 2.63E-01      | 7.30E-02   | J                | GEO-19   | 1.39E+00      |            |           |                     |                                   |                          |
| 3- and 4-methylphenol  | 106-44-5   | 3             | 1              | 33.33           | 4.10E-01      | 9.90E+00      | 2.10E-01              | J        | 1.79E+00      | 5.97E-01      | 2.10E-01   | J                | GEO-19   | 2.74E+00      |            |           |                     |                                   |                          |
| Acenaphthene           | 83-32-9    | 6             | 2              | 33.33           | 2.00E-01      | 1.50E+00      | 1.00E+00              | J        | 1.58E+02      | 7.02E+00      | 1.90E+02   | J                | GEO-21   | 3.00E+02      |            |           |                     |                                   |                          |
| Acenaphthylene         | 208-96-8   | 6             | 3              | 50              | 2.80E+00      | 1.50E+03      | 1.40E+00              | J        | 1.37E+02      | 1.34E+01      | 4.70E+01   | J                | GEO-21   | 3.01E+02      |            |           |                     |                                   |                          |
| Anthracene             | 120-12-7   | 6             | 5              | 83.33           | 5.30E-02      | 5.30E+02      | 2.10E+00              | Z        | 6.35E+02      | 2.06E+01      | 3.00E+03   | Z                | GEO-48   | 1.20E+03      |            |           |                     |                                   |                          |
| Benz(a)anthracene      | 56-55-3    | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 2.10E+00              | Z        | 2.16E+02      | 3.61E+01      | 9.30E+02   | Z                | GEO-48   | 3.65E+02      |            |           |                     |                                   |                          |
| Benz(a)pyrene          | 50-32-8    | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 3.00E+00              | Z        | 1.35E+02      | 3.16E+01      | 5.00E+02   | Z                | GEO-48   | 1.99E+02      |            |           |                     |                                   |                          |
| Benz(b)fluoranthene    | 205-99-2   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 3.50E+00              | Z        | 1.86E+02      | 4.59E+01      | 5.30E+02   | Z                | GEO-48   | 2.43E+02      |            |           |                     |                                   |                          |
| Benz(g)phenylene       | 191-24-2   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 1.60E+00              | J        | 4.27E+01      | 1.45E+01      | 1.30E+02   | J                | GEO-48   | 5.43E+01      |            |           |                     |                                   |                          |
| Benz(k)fluoranthene    | 207-08-9   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 1.80E+00              | Z        | 8.28E+01      | 1.94E+01      | 2.90E+02   | Z                | GEO-48   | 1.18E+02      |            |           |                     |                                   |                          |
| Carbazole              | 86-74-8    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 6.00E-01              | J        | 7.88E+01      | 9.34E+00      | 2.30E+02   | Z                | GEO-21   | 1.31E+02      |            |           |                     |                                   |                          |
| Chrysene               | 218-01-9   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 2.70E+00              | Z        | 1.79E+02      | 3.72E+01      | 6.90E+02   | Z                | GEO-48   | 2.73E+02      |            |           |                     |                                   |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 4.80E-01              | J        | 1.83E+01      | 4.98E+00      | 6.40E+01   | J                | GEO-48   | 2.57E+01      |            |           |                     |                                   |                          |
| Dibenzofuran           | 132-64-9   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 3.40E-01              | J        | 6.37E+01      | 3.65E+00      | 1.90E+02   | J                | GEO-21   | 1.09E+02      |            |           |                     |                                   |                          |
| Fluoranthene           | 206-44-0   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 2.80E+00              | Z        | 9.04E+02      | 7.65E+01      | 4.60E+03   | Z                | GEO-48   | 1.83E+03      |            |           |                     |                                   |                          |
| Fluorene               | 86-73-7    | 6             | 4              | 66.67           | 2.90E-01      | 2.60E-01      | 1.40E+00              | Z        | 3.44E+02      | 4.84E+00      | 1.80E+03   | Z                | GEO-48   | 7.21E+02      |            |           |                     |                                   |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 2.00E+00              | Z        | 6.94E+01      | 1.94E+01      | 2.50E+02   | J                | GEO-48   | 9.91E+01      |            |           |                     |                                   |                          |
| Naphthalene            | 91-20-3    | 6             | 4              | 66.67           | 2.80E+00      | 1.20E+01      | 6.80E-01              | J        | 4.50E+02      | 1.30E+01      | 2.20E+03   | J                | GEO-48   | 8.79E+02      |            |           |                     |                                   |                          |
| N-nitrosodiphenylamine | 86-30-6    | 3             | 1              | 33.33           | 3.70E-02      | 5.00E+00      | 2.00E+01              | J        | 9.06E-01      | 2.10E-01      | 2.00E-01   | J                | GEO-20   | 1.38E+00      |            |           |                     |                                   |                          |
| Phenanthrene           | 85-01-8    | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 3.10E-01              | J        | 1.19E+03      | 2.33E+01      | 6.40E+03   | Z                | GEO-48   | 2.57E+03      |            |           |                     |                                   |                          |
| Pyrene                 | 129-00-0   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 5.10E+00              | Z        | 8.76E+02      | 9.14E+01      | 4.40E+03   | Z                | GEO-20   | 1.74E+03      |            |           |                     |                                   |                          |

**Table 11**  
**Statistical Summary and Selection of COPCs in EU4 Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>Unrestricted Soil<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? | Is the<br>95%<br>UCL > TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|---|---|-----------------------------|
| <b>Semivolatiles</b>   |                  |                                 |                                   |  |   |   |                             |
| 2,4-dimethylphenol     | 6.40E+00         | 1.83E+13                        | Normal/Lognormal                  | 2.50E-01                                 | 1.56E+03                                    | no                                      |                             |
| 2-methylnaphthalene    | 3.66E+02         | 4.06E+62                        | Lognormal                         | 2.80E+02                                 | 3.13E+03                                    | no                                      |                             |
| 2-methylphenol         | 3.24E+00         | 3.99E+15                        | Normal/Lognormal                  | 7.30E-02                                 | 3.91E+03                                    | no                                      |                             |
| 3- and 4-methylphenol  | 6.40E+00         | 9.25E+13                        | Lognormal                         | 2.10E-01                                 | 3.91E+02                                    | no                                      |                             |
| Acenaphthene           | 4.05E+02         | 5.72E+11                        | Lognormal                         | 1.90E+02                                 | 4.69E+03                                    | no                                      |                             |
| Acenaphthylene         | 3.84E+02         | 4.19E+06                        | Lognormal                         | 4.70E+01                                 | 4.69E+03                                    | no                                      |                             |
| Anthracene             | 1.62E+03         | 8.99E+17                        | Lognormal                         | 3.00E+03                                 | 2.35E+04                                    | no                                      |                             |
| Benz(a)anthracene      | 5.17E+02         | 1.56E+07                        | Lognormal                         | 9.30E+02                                 | 8.75E-01                                    | YES                                     | YES - COPC                  |
| Benz(a)pyrene          | 2.98E+02         | 9.02E+05                        | Normal/Lognormal                  | 5.00E+02                                 | 8.75E-02                                    | YES                                     | YES - COPC                  |
| Benz(b)fluoranthene    | 3.85E+02         | 1.50E+06                        | Normal/Lognormal                  | 5.30E+02                                 | 8.75E-01                                    | YES                                     | YES - COPC                  |
| Benz(ghi)perylene      | 8.73E+01         | 2.25E+04                        | Normal/Lognormal                  | 1.30E+02                                 | 2.35E+03                                    | no                                      |                             |
| Benz(k)fluoranthene    | 1.80E+02         | 6.89E+05                        | Normal/Lognormal                  | 2.90E+02                                 | 8.75E+00                                    | YES                                     | YES - COPC                  |
| Carbazole              | 3.00E+02         | 1.23E+39                        | Normal/Lognormal                  | 2.30E+02                                 | 3.19E+01                                    | YES                                     | YES - COPC                  |
| Chrysene               | 4.04E+02         | 3.19E+06                        | Normal/Lognormal                  | 6.90E+02                                 | 8.75E+01                                    | YES                                     | YES - COPC                  |
| Dibenz(a,h)anthracene  | 3.95E+01         | 3.75E+04                        | Normal/Lognormal                  | 6.40E+01                                 | 8.75E-02                                    | YES                                     | YES - COPC                  |
| Dibenzofuran           | 2.48E+02         | 7.02E+50                        | Lognormal                         | 1.90E+02                                 | 3.13E+02                                    | no                                      |                             |
| Fluoranthene           | 2.41E+03         | 3.17E+09                        | Lognormal                         | 4.60E+03                                 | 3.13E-03                                    | YES                                     | YES - COPC                  |
| Fluorene               | 9.37E+02         | 1.95E+16                        | Lognormal                         | 1.80E+03                                 | 3.13E+03                                    | no                                      |                             |
| Indeno(1,2,3-cd)pyrene | 1.51E+02         | 1.07E+05                        | Normal/Lognormal                  | 2.50E+02                                 | 8.75E-01                                    | YES                                     | YES - COPC                  |
| Naphthalene            | 1.17E+03         | 7.77E+12                        | Lognormal                         | 2.20E+03                                 | 6.45E+02                                    | YES                                     | YES - COPC                  |
| N-nitrosodiphenylamine | 3.24E+00         | 6.32E+24                        | Normal/Lognormal                  | 2.00E+01                                 | 1.30E+02                                    | no                                      |                             |
| Phenanthrene           | 3.31E+03         | 9.46E+14                        | Lognormal                         | 6.40E+03                                 | 2.35E+03                                    | YES                                     | YES - COPC                  |
| Pyrene                 | 2.31E+03         | 6.40E+08                        | Lognormal                         | 4.40E+03                                 | 2.35E+03                                    | YES                                     | YES - COPC                  |

Table 2

Statistical Summary and Selection of COPCs in EU4 Soil (0-6' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Maximum mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Detected | Maximum mg/kg | Detected | Maximum mg/kg | Qualifier | Concentration mg/kg | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|---------------|------------|------------------------|----------|---------------|----------|---------------|-----------|---------------------|--------------------------|
| <b>Semivolatiles</b>   |            |               |                |                 |               |               |                       |          |               |               |            |                        |          |               |          |               |           |                     |                          |
| 2,4-dimethylphenol     | 105-67-9   | 9             | 3              | 33.33           | 7.30E-02      | 9.90E+00      | 7.90E-02              | J        | 1.63E+00      | 2.30E-01      | 8.90E+00   |                        | J        | GEO-21        | 3.16E+00 |               |           |                     |                          |
| 2-methylnaphthalene    | 91-57-6    | 9             | 8              | 88.89           | 4.00E-02      | 4.00E-02      | 6.20E-02              | J        | 2.11E+02      | 1.91E+00      | 1.50E+03   |                        | J        | GEO-21        | 4.93E+02 |               |           |                     |                          |
| 2-methylphenol         | 95-48-7    | 9             | 1              | 11.11           | 3.70E-02      | 5.00E+00      | 7.30E-02              | J        | 4.61E-01      | 8.75E-02      | 7.30E-02   |                        | J        | GEO-19        | 8.70E-01 |               |           |                     |                          |
| 3- and 4-methylphenol  | 106-44-5   | 9             | 1              | 11.11           | 7.30E-02      | 9.90E+00      | 2.10E-01              | J        | 9.17E-01      | 1.82E-01      | 2.10E-01   |                        | J        | GEO-19        | 1.72E+00 |               |           |                     |                          |
| Acenaphthene           | 83-32-9    | 18            | 9              | 50              | 4.09E-02      | 1.50E+03      | 9.70E-02              | J        | 1.30E+02      | 2.05E+00      | 1.20E+03   |                        | J        | GEO-21        | 3.21E+02 |               |           |                     |                          |
| Acenaphthylene         | 208-96-8   | 18            | 7              | 38.89           | 3.80E-02      | 1.50E+03      | 8.30E-02              | J        | 4.90E+01      | 1.14E+00      | 5.00E+01   |                        | J        | GEO-21        | 1.76E+02 |               |           |                     |                          |
| Anthracene             | 120-12-7   | 18            | 14             | 77.78           | 2.60E-03      | 5.30E-02      | 1.20E-01              | Z        | 3.15E+02      | 1.93E+00      | 3.00E+03   |                        | Z        | GEO-48        | 8.06E+02 |               |           |                     |                          |
| Benz(a)anthracene      | 56-55-3    | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 4.90E-03              | J        | 9.72E+01      | 2.57E+00      | 9.30E+02   |                        | Z        | GEO-48        | 2.34E+02 |               |           |                     |                          |
| Benz(a)pyrene          | 50-32-8    | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 1.10E-02              | J        | 5.71E+01      | 2.04E+00      | 5.00E+02   |                        | Z        | GEO-48        | 1.29E+02 |               |           |                     |                          |
| Benz(b)fluoranthene    | 205-99-2   | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 1.10E-02              | J        | 7.89E+01      | 2.53E+00      | 5.30E+02   |                        | Z        | GEO-48        | 1.65E+02 |               |           |                     |                          |
| Benz(ghi)perylene      | 191-24-2   | 18            | 17             | 94.44           | 3.80E-02      | 3.80E-02      | 8.90E-03              | J        | 1.86E+01      | 8.49E-01      | 1.30E+02   |                        | J        | GEO-48        | 3.79E+01 |               |           |                     |                          |
| Benz(k)fluoranthene    | 207-08-9   | 18            | 16             | 88.89           | 3.80E-02      | 4.00E-02      | 5.60E-03              | J        | 3.27E+01      | 1.03E+00      | 2.90E+02   |                        | Z        | GEO-48        | 7.57E+01 |               |           |                     |                          |
| Carbazole              | 86-74-8    | 9             | 8              | 88.89           | 4.00E-02      | 4.00E-02      | 2.10E-01              | J        | 9.80E+01      | 2.78E+00      | 6.20E+02   |                        | Z        | GEO-21        | 2.10E+02 |               |           |                     |                          |
| Chrysene               | 218-01-9   | 18            | 17             | 94.44           | 5.10E-03      | 9.90E-03      | 9.90E-03              | J        | 8.58E+01      | 2.42E+00      | 6.90E+02   |                        | Z        | GEO-48        | 1.88E+02 |               |           |                     |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 18            | 14             | 77.78           | 2.60E-03      | 4.00E-02      | 3.80E-03              | J        | 7.52E+00      | 2.96E-01      | 6.40E+01   |                        | J        | GEO-48        | 1.68E+01 |               |           |                     |                          |
| Dibenzofuran           | 132-64-9   | 9             | 8              | 88.89           | 4.00E-02      | 4.00E-02      | 7.80E-02              | J        | 1.54E+02      | 2.43E+00      | 1.10E+03   |                        | J        | GEO-21        | 3.61E+02 |               |           |                     |                          |
| Fluoranthene           | 206-44-0   | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 1.00E-02              | J        | 4.29E+02      | 7.56E+00      | 4.60E+03   |                        | Z        | GEO-48        | 1.15E+03 |               |           |                     |                          |
| Fluorene               | 86-73-7    | 18            | 12             | 66.67           | 1.30E-02      | 2.60E-01      | 1.40E-01              | J        | 2.09E+02      | 1.40E+00      | 1.80E+03   |                        | Z        | GEO-48        | 5.31E+02 |               |           |                     |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 18            | 17             | 94.44           | 6.00E-03      | 6.00E-03      | 9.00E-03              | J        | 2.83E+01      | 1.05E+00      | 2.50E+02   |                        | J        | GEO-48        | 6.42E+01 |               |           |                     |                          |
| Naphthalene            | 91-20-3    | 18            | 12             | 66.67           | 4.00E-02      | 1.20E+01      | 7.60E-02              | J        | 3.73E+02      | 2.99E+00      | 3.50E+03   |                        | J        | GEO-21        | 9.37E+02 |               |           |                     |                          |
| N-nitrosodiphenylamine | 86-30-6    | 9             | 1              | 11.11           | 3.70E-02      | 5.00E+00      | 2.00E-01              | J        | 4.66E-01      | 8.11E-02      | 2.00E-01   |                        | Z        | GEO-20        | 8.69E-01 |               |           |                     |                          |
| Phenanthrene           | 85-01-8    | 18            | 17             | 94.44           | 4.00E-02      | 5.30E-03      | 1.60E-01              | J        | 6.51E+02      | 4.31E+00      | 6.40E+03   |                        | Z        | GEO-48        | 1.72E+03 |               |           |                     |                          |
| Pyrene                 | 129-00-0   | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 1.60E-02              | J        | 3.77E+02      | 7.72E+00      | 4.40E+03   |                        | Z        | GEO-20        | 1.06E+03 |               |           |                     |                          |

**Table 12**  
**Statistical Summary and Selection of COPCs in EU4 Soil (0-6' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg | 99% Confidence | Logarithmic Distribution<br>Concentration<br>mg/kg | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>TRG | Is the<br>Maximum<br>Detected ><br>TRG? |                             |
|------------------------|------------------|------------------|----------------|--|--|---------------|---|-----------------------------|
|                        |                  |                  |                |  |  |               | Is the 95%<br>UCL ><br>TRG?             | Is the 95%<br>UCL ><br>TRG? |
| <b>Semivolatiles</b>   |                  |                  |                |  |  |               |   |                             |
| 2,4-dimethylphenol     | 3.59E+00         | 1.33E+02         | Lognormal      | 8.90E+00   | 4.08E+04                                 | no            |   |                             |
| 2-methylnaphthalene    | 5.17E+02         | 1.40E+11         | Lognormal      | 1.50E+03   | 8.18E+04                                 | no            |   |                             |
| 2-methylphenol         | 1.00E+00         | 1.71E+01         | Lognormal      | 7.30E-02   | 1.02E+05                                 | no            |   |                             |
| 3- and 4-methylphenol  | 1.98E+00         | 3.49E+01         | Lognormal      | 2.10E-01   | 1.02E+04                                 | no            |   |                             |
| Acenaphthene           | 2.61E+02         | 7.82E+05         | Lognormal      | 1.20E+03   | 1.23E+05                                 | no            |   |                             |
| Acenaphthylene         | 1.21E+02         | 8.90E+03         | Lognormal      | 5.00E+01   | 1.23E+05                                 | no            |   |                             |
| Anthracene             | 6.46E+02         | 3.29E+08         | Lognormal      | 3.00E+03   | 6.13E+05                                 | no            |   |                             |
| Benzo(a)anthracene     | 1.93E+02         | 1.41E+06         | Lognormal      | 9.30E+02   | 7.84E+00                                 | YES           | YES - COPC                              |                             |
| Benzo(a)pyrene         | 1.10E+02         | 1.87E+05         | Lognormal      | 5.00E+02   | 7.84E-01                                 | YES           | YES - COPC                              |                             |
| Benzo(b)fluoranthene   | 1.47E+02         | 5.68E+05         | Lognormal      | 5.30E+02   | 7.84E+00                                 | YES           | YES - COPC                              |                             |
| Benzo(g,h,i)perylene   | 3.42E+01         | 2.74E+04         | Lognormal      | 1.30E+02   | 6.13E+04                                 | no            |   |                             |
| Benzo(k)fluoranthene   | 6.37E+01         | 2.16E+05         | Lognormal      | 2.90E+02   | 7.84E+01                                 | YES           | YES - COPC                              |                             |
| Carbazole              | 2.28E+02         | 1.00E+08         | Lognormal      | 6.20E+02   | 2.86E+02                                 | YES           | YES - COPC                              |                             |
| Chrysene               | 1.63E+02         | 3.02E+06         | Lognormal      | 6.90E+02   | 7.84E+02                                 | no            |   | YES*                        |
| Dibenz(a,h)anthracene  | 1.44E+01         | 1.67E+04         | Lognormal      | 6.40E+01   | 7.84E-01                                 | YES           | YES - COPC                              |                             |
| Dibenzofuran           | 3.78E+02         | 4.19E+09         | Lognormal      | 1.10E+03   | 8.18E+03                                 | no            |   |                             |
| Fluoranthene           | 8.99E+02         | 1.07E+07         | Lognormal      | 4.60E+03   | 8.17E+04                                 | no            |   |                             |
| Fluorene               | 4.27E+02         | 4.71E+07         | Lognormal      | 1.80E+03   | 8.17E+04                                 | no            |   |                             |
| Indeno(1,2,3-cd)pyrene | 5.47E+01         | 1.10E+05         | Lognormal      | 2.50E+02   | 7.84E+00                                 | YES           | YES - COPC                              |                             |
| Naphthalene            | 7.57E+02         | 2.43E+07         | Lognormal      | 3.50E+03   | 8.24E+02                                 | YES           | YES - COPC                              |                             |
| N-nitrosodiphenylamine | 1.00E+00         | 2.74E+01         | Lognormal      | 2.00E-01   | 1.17E+03                                 | no            |   |                             |
| Phenanthrene           | 1.35E+03         | 6.44E+08         | Lognormal      | 6.40E+03   | 6.13E+04                                 | no            |   |                             |
| Pyrene                 | 8.10E+02         | 5.56E+06         | Lognormal      | 4.40E+03   | 6.13E+04                                 | no            |   |                             |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table I<sub>3</sub>

Statistical Summary and Selection of COPCs in EU4 Soil (0-20' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Minimum Detection Limit mg/kg | Maximum Detection Limit mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Maximum Detected mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-------------------------------|-------------------------------|------------|------------------------|------------------------|------------------------|-----------------------------------|--------------------------|
| Semivariates           |            |               |                |                 |               |               |                               |                               |            |                        |                        |                        |                                   |                          |
| 2,4-dimethylphenol     | 105-67-9   | 11            | 4              | 36.36           | 7.30E-02      | 9.90E+00      | 7.90E-02                      | 9.90E+00                      | 1.95E-01   | 8.90E+00               | J                      | GEO-21                 | 2.89E+00                          |                          |
| 2-methylnaphthalene    | 91-57-6    | 11            | 10             | 90.91           | 4.00E-02      | 6.20E-02      | 4.00E-02                      | 6.20E-02                      | 1.50E+03   | 1.50E+03               | J                      | GEO-21                 | 4.47E+02                          |                          |
| 2-methylphenol         | 95-48-7    | 11            | 2              | 18.18           | 3.70E-02      | 5.00E+00      | 5.10E-02                      | 7.23E-02                      | 3.83E-01   | 7.30E-02               | J                      | GEO-19                 | 7.97E-01                          |                          |
| 3- and 4-methylphenol  | 106-44-5   | 11            | 1              | 9.09            | 7.30E-02      | 9.90E+00      | 2.10E-01                      | 7.20E-01                      | 7.58E-01   | 1.37E-01               | J                      | GEO-19                 | 1.58E+00                          |                          |
| Acenaphthene           | 83-32-9    | 21            | 12             | 57.14           | 4.00E-02      | 1.50E+03      | 9.70E-02                      | 1.14E+02                      | 1.14E+02   | 2.39E+00               | 1.20E+03               | J                      | GEO-21                            | 2.98E+02                 |
| Acenaphthylene         | 208-96-8   | 21            | 8              | 38.1            | 3.80E-02      | 1.50E+03      | 8.30E-02                      | 4.22E+01                      | 9.69E-01   | 5.00E+01               | J                      | GEO-21                 | 1.63E+02                          |                          |
| Anthracene             | 120-12-7   | 21            | 17             | 80.95           | 2.60E-03      | 5.30E-02      | 1.20E-01                      | 2.71E+02                      | 2.71E+02   | 2.00E+00               | 3.00E+03               | Z                      | GEO-48                            | 7.51E+02                 |
| Benz[a]anthracene      | 56-55-3    | 21            | 21             | 100             | 0.00E+00      | 0.00E+00      | 4.90E-03                      | 9.40E-03                      | 8.42E+01   | 2.53E+00               | 9.30E+02               | Z                      | GEO-48                            | 2.18E+02                 |
| Benz[a]pyrene          | 50-32-8    | 21            | 21             | 100             | 0.00E+00      | 0.00E+00      | 1.10E-02                      | 1.10E-02                      | 4.93E+01   | 1.83E+00               | 5.00E+02               | Z                      | GEO-48                            | 1.21E+02                 |
| Benz[0]fluoranthene    | 205-99-2   | 21            | 21             | 100             | 0.00E+00      | 0.00E+00      | 1.10E-02                      | 1.10E-02                      | 6.80E+01   | 2.27E+00               | 5.30E+02               | Z                      | GEO-48                            | 1.55E+02                 |
| Benzog[hi]perylene     | 191-24-2   | 21            | 19             | 90.48           | 3.80E-02      | 4.00E-02      | 8.90E-03                      | 1.61E+01                      | 7.03E-01   | 1.30E+02               | J                      | GEO-48                 | 3.56E+01                          |                          |
| Benzo[K]fluoranthene   | 207-08-9   | 21            | 19             | 90.48           | 3.80E-02      | 4.00E-02      | 5.60E-03                      | 1.28E+01                      | 9.39E-01   | 2.90E+02               | Z                      | GEO-48                 | 7.07E+01                          |                          |
| Carbazole              | 86-74-8    | 11            | 10             | 90.91           | 4.00E-02      | 4.00E-02      | 2.10E-01                      | 8.10E+01                      | 2.49E+00   | 6.20E+02               | J                      | GEO-21                 | 1.91E+02                          |                          |
| Chrysene               | 218-01-9   | 21            | 20             | 95.24           | 5.10E-03      | 5.10E-03      | 9.90E-03                      | 7.42E+01                      | 2.29E+00   | 6.90E+02               | Z                      | GEO-48                 | 1.76E+02                          |                          |
| Dibenz[a,h]anthracene  | 53-70-3    | 21            | 16             | 76.19           | 2.60E-03      | 4.00E-02      | 3.80E-03                      | J                             | 6.48E+00   | 2.65E-01               | 6.40E+01               | J                      | GEO-48                            | 1.57E+01                 |
| Dibenzofuran           | 132-64-9   | 11            | 10             | 90.91           | 4.00E-02      | 4.00E-02      | 7.80E-02                      | J                             | 1.30E+02   | 2.74E+00               | 1.10E+03               | J                      | GEO-21                            | 3.27E+02                 |
| Fluoranthene           | 206-44-0   | 21            | 21             | 100             | 0.00E+00      | 0.00E+00      | 1.00E-02                      | J                             | 3.72E+02   | 7.99E+00               | 4.60E+03               | Z                      | GEO-48                            | 1.07E+03                 |
| Fluorene               | 86-73-7    | 21            | 15             | 71.43           | 1.30E-02      | 2.60E-01      | 1.40E-01                      | J                             | 1.82E+02   | 1.76E+00               | 1.80E+03               | Z                      | GEO-48                            | 4.94E+02                 |
| Indeno[1,2,3-cd]pyrene | 193-39-5   | 21            | 20             | 95.24           | 6.00E-03      | 6.00E-03      | 9.00E-03                      | J                             | 2.44E+01   | 9.11E-01               | 2.50E+02               | J                      | GEO-48                            | 6.00E+01                 |
| Naphthalene            | 91-20-3    | 21            | 15             | 71.43           | 4.00E-02      | 1.20E+01      | 7.60E-02                      | J                             | 3.27E+02   | 3.63E+00               | 3.50E+03               | Z                      | GEO-21                            | 8.72E+02                 |
| N-nitrosodiphenylamine | 86-30-6    | 11            | 1              | 9.09            | 3.70E-02      | 5.00E+00      | 2.00E-01                      | J                             | 3.84E-01   | 6.25E-02               | 2.00E+01               | Z                      | GEO-20                            | 7.98E-01                 |
| Phenanthrene           | 85-01-8    | 21            | 20             | 95.24           | 4.00E-02      | 4.00E-02      | 5.30E-03                      | J                             | 5.68E+02   | 5.46E+00               | 6.40E+03               | Z                      | GEO-48                            | 1.00E+03                 |
| Phenol                 | 108-95-2   | 11            | 1              | 9.09            | 7.30E-02      | 9.90E+00      | 1.00E-01                      | J                             | 7.48E-01   | 1.28E-01               | 1.00E-01               | J                      | GEO-20                            | 1.58E+00                 |
| Pyrene                 | 129-00-0   | 21            | 21             | 100             | 0.00E+00      | 0.00E+00      | 1.60E-02                      | J                             | 3.26E+02   | 7.75E+00               | 4.40E+03               | Z                      | GEO-20                            | 9.82E+02                 |

**Table I3**  
**Statistical Summary and Selection of COPCs in EU4 Soil (0-20' bgs)**  
**Kerr McGee, Hatiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg | Logarithmic<br>99%<br>Confidence | Distribution<br>99%<br>Concentration | Exposure Point<br>mg/kg | Tier 1                          |                                     | Is the<br>UCL > TRG?<br>no |
|------------------------|------------------|------------------|----------------------------------|--------------------------------------|-------------------------|---------------------------------|-------------------------------------|----------------------------|
|                        |                  |                  |                                  |                                      |                         | Restricted<br>Soil TRG<br>mg/kg | Maximum<br>Detected ><br>TRG?<br>no |                            |
| <b>Senivolatiles</b>   |                  |                  |                                  |                                      |                         |                                 |                                     |                            |
| 2,4-dimethylphenol     | 2.94E+00         | 2.48E+01         | Lognormal                        | 8.90E+00                             | 4.08E+04                |                                 |                                     |                            |
| 2-methylnaphthalene    | 4.23E+02         | 4.96E+08         | Lognormal                        | 1.50E+03                             | 8.18E+04                |                                 |                                     |                            |
| 2-methylphenol         | 8.19E-01         | 3.98E+00         | Unknown                          | 7.30E-02                             | 1.02E-05                |                                 |                                     |                            |
| 3- and 4-methylphenol  | 1.62E+00         | 9.14E+00         | Unknown                          | 2.10E-01                             | 1.02E-04                |                                 |                                     |                            |
| Acenaphthene           | 2.26E+02         | 1.38E+05         | Lognormal                        | 1.20E+03                             | 1.23E+05                |                                 |                                     |                            |
| Acenaphthylene         | 1.03E+02         | 2.70E+03         | Lognormal                        | 5.00E+01                             | 1.23E+05                |                                 |                                     |                            |
| Anthracene             | 5.54E+02         | 1.16E+07         | Lognormal                        | 3.00E+03                             | 6.13E+05                |                                 |                                     |                            |
| Benz(a)anthracene      | 1.66E+02         | 1.53E+05         | Lognormal                        | 9.30E+02                             | 7.84E+00                | YES - COPC                      |                                     |                            |
| Benz(a)pyrene          | 9.47E+01         | 2.66E+04         | Lognormal                        | 5.00E+02                             | 7.84E+01                | YES - COPC                      |                                     |                            |
| Benz(b)fluoranthene    | 1.26E+02         | 6.79E+04         | Lognormal                        | 5.30E+02                             | 7.84E+00                | YES - COPC                      |                                     |                            |
| Benz(g,h)perylene      | 2.95E+01         | 5.59E+03         | Lognormal                        | 1.30E+02                             | 6.13E+04                |                                 |                                     |                            |
| Benz(k)fluoranthene    | 5.48E+01         | 2.80E+04         | Lognormal                        | 2.90E+02                             | 7.84E+01                | YES - COPC                      |                                     |                            |
| Carbazole              | 1.86E+02         | 1.17E+06         | Lognormal                        | 6.20E+02                             | 2.86E+02                | YES - COPC                      |                                     |                            |
| Chrysene               | 1.40E+02         | 2.66E+05         | Lognormal                        | 6.90E+02                             | 7.84E+02                | *YES                            |                                     |                            |
| Dibenz(a,h)anthracene  | 1.24E+01         | 2.39E+03         | Lognormal                        | 6.40E+01                             | 7.84E+01                | YES - COPC                      |                                     |                            |
| Dibenzofuran           | 3.09E+02         | 3.01E+07         | Lognormal                        | 1.10E+03                             | 8.18E+03                |                                 |                                     |                            |
| Fluoranthene           | 7.73E+02         | 1.06E+06         | Lognormal                        | 4.60E+03                             | 8.17E+04                |                                 |                                     |                            |
| Fluorene               | 3.68E+02         | 4.04E+06         | Lognormal                        | 1.80E+03                             | 8.17E+04                |                                 |                                     |                            |
| Indeno(1,2,3-cd)pyrene | 4.70E+01         | 1.53E+04         | Lognormal                        | 2.50E+02                             | 7.84E+00                | YES - COPC                      |                                     |                            |
| Naphthalene            | 6.55E+02         | 2.85E+06         | Lognormal                        | 3.50E+03                             | 8.24E+02                | YES - COPC                      |                                     |                            |
| N-nitrosodiphenylamine | 8.20E-01         | 5.77E+00         | Unknown                          | 2.00E-01                             | 1.17E+03                | no                              |                                     |                            |
| Phenanthrene           | 1.17E+03         | 4.13E+07         | Lognormal                        | 6.40E+03                             | 6.13E+04                | no                              |                                     |                            |
| Phenol                 | 1.61E+00         | 8.43E+00         | Unknown                          | 1.00E-01                             | 1.23E+05                | no                              |                                     |                            |
| Pyrene                 | 6.96E+02         | 5.88E+05         | Lognormal                        | 4.40E+03                             | 6.13E+04                | no                              |                                     |                            |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000): constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

*Table I<sub>a</sub>*  
*Statistical Summary and Selection of COPCs in EU5 Soil (0-1' bgs)*  
*Kerr McGee, Hattiesburg, MS*

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum Detection Limit mg/kg | Minimum Detected mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Detected Qualifier | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|-------------------------------|------------------------|------------|------------------------|------------------------|--------------------|-----------------------------|-----------------------------------|--------------------------|
| Semivolatiles          |            |               |                |                 |               |                               |                        |            |                        |                        |                    |                             |                                   |                          |
| 2,4-dimethylphenol     | 105-67-9   | 6             | 1              | 16.67           | 7.60E-02      | 4.55E-01                      | 1.10E-01               | J          | 1.07E-01               | 8.03E-02               | 1.10E-01           | J                           | GEO-30                            | 8.38E-02                 |
| 2-methylnaphthalene    | 91-57-6    | 6             | 4              | 66.67           | 3.80E-02      | 3.90E-02                      | 5.10E-02               | J          | 1.68E-01               | 1.71E-01               | 9.20E-01           | J                           | GEO-30                            | 3.69E+00                 |
| 2-methylphenol         | 95-48-7    | 6             | 1              | 16.67           | 3.80E-02      | 2.09E-01                      | 4.20E-02               | J          | 4.94E-02               | 3.77E-02               | 4.20E-02           | J                           | GEO-30                            | 3.92E-02                 |
| 3- and 4-methylphenol  | 106-44-5   | 6             | 1              | 16.67           | 7.60E-02      | 4.13E-01                      | 1.40E-01               | J          | 1.08E-01               | 8.22E-02               | 1.40E-01           | J                           | GEO-30                            | 7.94E-02                 |
| Acenaphthene           | 83-32-9    | 8             | 2              | 25              | 3.80E-02      | 2.40E-00                      | 1.60E-01               | J          | 6.91E-00               | 1.92E-01               | 5.35E-01           | J                           | GEO-33                            | 1.88E+01                 |
| Acenaphthylene         | 208-96-8   | 8             | 5              | 62.5            | 3.80E-02      | 2.40E-00                      | 4.80E-02               | J          | 2.61E-00               | 4.28E-01               | 1.60E-01           | J                           | GEO-33                            | 5.47E+00                 |
| Anthracene             | 120-12-7   | 8             | 4              | 50              | 1.00E-02      | 4.40E-02                      | 1.30E-01               | J          | 1.07E-01               | 2.00E-01               | 7.97E-01           | J                           | GEO-33                            | 2.79E+01                 |
| Benz(a)anthracene      | 56-55-3    | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 2.60E-01               | Z          | 1.30E-01               | 1.39E+00               | 8.35E+01           | J                           | GEO-33                            | 2.87E+01                 |
| Benz(a)pyrene          | 50-32-8    | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 3.10E-01               | J          | 8.82E+00               | 1.38E+00               | 5.25E+01           | J                           | GEO-33                            | 1.79E+01                 |
| Benz(b)fluoranthene    | 205-99-2   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 4.40E-01               | Z          | 1.38E-01               | 2.15E+00               | 7.95E+01           | J                           | GEO-33                            | 2.71E+01                 |
| Benz(ghi)perylene      | 191-24-2   | 8             | 6              | 75              | 3.80E-02      | 3.80E-02                      | 2.30E-01               | J          | 4.30E+00               | 5.42E-01               | 2.55E+01           | J                           | GEO-33                            | 8.71E+00                 |
| Benz(k)fluoranthene    | 207-08-9   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 2.10E-01               | Z          | 5.07E+00               | 1.01E+00               | 2.85E+01           | J                           | GEO-33                            | 9.68E+00                 |
| Carbazole              | 86-74-8    | 6             | 3              | 50              | 3.80E-02      | 3.90E-02                      | 5.30E-01               | J          | 2.63E+00               | 2.10E-01               | 1.35E+01           | J                           | GEO-33                            | 5.36E+00                 |
| Chrysene               | 218-01-9   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 2.70E-01               | Z          | 1.35E-01               | 1.62E+00               | 8.25E+01           | J                           | GEO-33                            | 2.83E+01                 |
| Dibenz(a,h)anthracene  | 53-70-3    | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 6.30E-02               | J          | 1.33E+00               | 3.02E-01               | 7.45E+00           | J                           | GEO-33                            | 2.53E+00                 |
| Dibenzofuran           | 132-64-9   | 6             | 4              | 66.67           | 3.80E-02      | 3.90E-02                      | 3.90E-02               | J          | 4.94E+00               | 1.77E-01               | 2.90E+01           | J                           | GEO-33                            | 1.18E+01                 |
| Fluoranthene           | 206-44-0   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 1.30E-01               | J          | 5.01E+01               | 2.31E+00               | 3.55E+02           | J                           | GEO-33                            | 1.23E+02                 |
| Fluorene               | 86-73-7    | 8             | 4              | 50              | 3.80E-02      | 5.20E-02                      | 3.30E-01               | J          | 8.10E+00               | 1.90E-01               | 6.30E+01           | J                           | GEO-33                            | 2.22E+01                 |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 2.60E-01               | J          | 5.49E+00               | 1.02E+00               | 3.10E-01           | J                           | GEO-33                            | 1.05E+01                 |
| Naphthalene            | 91-20-3    | 8             | 5              | 62.5            | 3.80E-02      | 5.60E-01                      | 7.50E-02               | J          | 1.63E+00               | 2.97E-01               | 6.85E+00           | J                           | GEO-33                            | 2.65E+00                 |
| Phenanthrene           | 85-01-8    | 8             | 6              | 75              | 3.80E-02      | 3.90E-02                      | 1.50E-01               | J          | 3.25E+01               | 7.56E-01               | 2.45E+02           | J                           | GEO-33                            | 8.59E+01                 |
| Phenol                 | 108-95-2   | 6             | 4              | 66.67           | 7.70E-02      | 4.13E-01                      | 1.40E-01               | J          | 1.99E-01               | 1.64E-01               | 3.80E-01           | J                           | GEO-29                            | 1.13E+01                 |
| Pyrene                 | 129-00-0   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02                      | 2.50E-01               | J          | 3.78E+01               | 2.54E+00               | 2.60E+02           | J                           | GEO-33                            | 9.01E+01                 |

**Table 14**  
**Statistical Summary and Selection of COPCs in EU5 Soil (θ-I' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                 | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier 1                            |                               | Is the Maximum<br>Detected > TRG?<br>UCL > TRG? |
|-----------------------------|------------------|---------------------------------|-----------------------------------|--|-----------------------------------|-------------------------------|---|
|                             |                  |                                 |                                   |  | Unrestricted Soil<br>TRG<br>mg/kg | Is the 95%<br>Detected > TRG? |   |
| <b>Senivolatiles</b>        |                  |                                 |                                   |  |                                   |                               |   |
| 2,4-dimethylphenol          | 1.76E-01         | 4.42E-01                        | Normal/Lognormal                  | 1.10E-01                                 | 1.56E+03                          | no                            |   |
| 2-methylnaphthalene         | 4.71E+00         | 4.63E+04                        | Lognormal                         | 9.20E+00                                 | 3.13E+03                          | no                            |   |
| 2-methylphenol <sup>*</sup> | 8.17E-02         | 1.82E-01                        | Normal/Lognormal                  | 4.20E-02                                 | 3.91E+03                          | no                            |   |
| 3- and 4-methylphenol       | 1.74E-01         | 4.55E-01                        | Normal/Lognormal                  | 1.40E-01                                 | 3.91E+02                          | no                            |   |
| Acenaphthene                | 1.95E+01         | 3.47E+04                        | Lognormal                         | 5.35E+01                                 | 4.69E+03                          | no                            |   |
| Acenaphthylene              | 6.27E+00         | 1.25E+03                        | Lognormal                         | 1.60E+01                                 | 4.69E+03                          | no                            |   |
| Anthracene                  | 2.94E+01         | 1.81E+07                        | Lognormal                         | 7.97E+01                                 | 2.35E+04                          | no                            |   |
| Benz(a)anthracene           | 3.23E+01         | 9.31E+04                        | Lognormal                         | 8.35E+01                                 | 8.75E-01                          | YES - COPC                    |   |
| Benz(a)pyrene               | 2.08E+01         | 2.16E+04                        | Lognormal                         | 5.25E+01                                 | 8.75E-02                          | YES - COPC                    |   |
| Benz(b)fluoranthene         | 3.20E+01         | 8.21E+04                        | Lognormal                         | 7.95E+01                                 | 8.75E-01                          | YES - COPC                    |   |
| Benz(ghi)perylene           | 1.01E+01         | 2.82E+04                        | Lognormal                         | 2.55E+01                                 | 2.35E+03                          | no                            |   |
| Benz(k)fluoranthene         | 1.16E+01         | 3.72E+03                        | Lognormal                         | 2.85E+01                                 | 8.75E+00                          | YES - COPC                    |   |
| Carbazole                   | 7.04E+00         | 8.43E+06                        | Lognormal                         | 1.35E+01                                 | 3.19E+01                          | no                            |   |
| Chrysene                    | 3.25E+01         | 1.14E+05                        | Lognormal                         | 8.25E+01                                 | 8.75E+01                          | YES*                          |   |
| Dibenz(a,h)anthracene       | 3.02E+00         | 1.76E+02                        | Lognormal                         | 7.45E+00                                 | 8.75E-02                          | YES - COPC                    |   |
| Dibenzofuran                | 1.46E+01         | 4.28E+06                        | Lognormal                         | 2.90E+01                                 | 3.13E+02                          | no                            |   |
| Fluoranthene                | 1.33E+02         | 2.93E+07                        | Lognormal                         | 3.55E+02                                 | 3.13E+03                          | no                            |   |
| Fluorene                    | 2.30E+01         | 1.21E+05                        | Lognormal                         | 6.30E+01                                 | 3.13E+03                          | no                            |   |
| Indeno(1,2,3-cd)pyrene      | 1.25E+01         | 5.34E+03                        | Lognormal                         | 3.10E+01                                 | 8.75E-01                          | YES - COPC                    |   |
| Naphthalene                 | 3.41E+00         | 1.19E+03                        | Lognormal                         | 6.85E+00                                 | 6.45E+02                          | no                            |   |
| Phenanthrene                | 9.00E+01         | 2.29E+07                        | Lognormal                         | 2.45E+02                                 | 2.35E+03                          | no                            |   |
| Phenol                      | 2.92E-01         | 7.33E-01                        | Normal/Lognormal                  | 3.80E-01                                 | 4.69E+04                          | no                            |   |
| Pyrene                      | 9.82E+01         | 4.13E+06                        | Lognormal                         | 2.60E+02                                 | 2.35E+03                          | no                            |   |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

*Table I*  
*Statistical Summary and Selection of COPCs in EU5 Soil (0-6' bgs)*  
*Kerr McGee, Hattiesburg, MS*

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detected Limit mg/kg | Minimum Detected mg/kg | Mean mg/kg | Logarithmic Mean | Maximum mg/kg | Detected mg/kg | Maximum Qualifier mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|----------------------|------------------------|------------|------------------|---------------|----------------|-------------------------|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>       |            |               |                |                 |               |               |                      |                        |            |                  |               |                |                         |                                   |                          |
| 2,4-dimethylphenol         | 105-67-9   | 18            | 1              | 5.56            | 7.50E-02      | 4.55E-01      | 1.10E-01             | J                      | 6.17E-02   | 4.98E-02         | 1.10E-01      | J              | GEO-30                  | 5.60E-02                          |                          |
| 2-methylnaphthalene        | 91-57-6    | 18            | 5              | 27.78           | 3.80E-02      | 4.10E-02      | 5.10E-02             | J                      | 6.21E-01   | 4.99E-02         | 9.20E+00      | J              | GEO-33                  | 2.15E+00                          |                          |
| 2-methylphenol             | 95-48-7    | 18            | 1              | 5.56            | 3.80E-02      | 2.09E-01      | 4.20E-02             | J                      | 2.95E-02   | 2.44E-02         | 4.20E-02      | J              | GEO-30                  | 2.57E-02                          |                          |
| 3- and 4-methylphenol      | 106-44-5   | 18            | 1              | 5.56            | 7.50E-02      | 4.13E-01      | 1.40E-01             | J                      | 6.22E-02   | 5.02E-02         | 1.40E-01      | J              | GEO-30                  | 5.46E-02                          |                          |
| Acenaphthene               | 83-32-9    | 24            | 4              | 16.67           | 2.90E-02      | 2.40E+00      | 1.10E-01             | J                      | 2.53E+00   | 6.24E-02         | 5.35E+01      | J              | GEO-33                  | 1.09E+01                          |                          |
| Acenaphthylene             | 208-96-8   | 24            | 6              | 25              | 3.80E-02      | 2.40E+00      | 4.80E-02             | J                      | 9.59E-01   | 7.93E-02         | 1.60E+01      | J              | GEO-33                  | 3.26E+00                          |                          |
| Anthracene                 | 120-12-7   | 24            | 6              | 25              | 5.40E-04      | 6.00E-02      | 1.30E-01             | J                      | 3.94E+00   | 4.33E-02         | 7.97E+01      | J              | GEO-33                  | 1.63E+01                          |                          |
| Benzo(a)anthracene         | 56-55-3    | 24            | 13             | 54.17           | 3.80E-02      | 4.10E-02      | 6.80E-03             | Z                      | 4.81E+00   | 1.11E-01         | 8.35E+01      | Z              | GEO-33                  | 1.71E+01                          |                          |
| Benzo(a)pyrene             | 50-32-8    | 24            | 13             | 54.17           | 3.80E-02      | 4.10E-02      | 8.30E-03             | Z                      | 3.17E+00   | 1.15E-01         | 5.25E+01      | Z              | GEO-33                  | 1.07E+01                          |                          |
| Benzo(b)fluoranthene       | 205-99-2   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 9.00E-03             | Z                      | 4.99E+00   | 1.59E-01         | 7.95E+01      | J              | GEO-33                  | 1.63E+01                          |                          |
| Benzo(ghi)perylene         | 191-24-2   | 24            | 12             | 50              | 3.80E-02      | 4.10E-02      | 6.70E-03             | J                      | 1.55E+00   | 7.48E-02         | 2.55E+01      | J              | GEO-33                  | 5.22E+00                          |                          |
| Benzo(k)fluoranthene       | 207-08-9   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 4.70E-03             | Z                      | 1.84E+00   | 9.91E-02         | 2.85E+01      | Z              | GEO-33                  | 5.86E+00                          |                          |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 18            | 2              | 11.11           | 7.50E-02      | 4.13E-01      | 1.40E-01             | J                      | 6.84E-02   | 5.41E-02         | 1.50E-01      | J              | GEO-32                  | 5.80E-02                          |                          |
| Carbazole                  | 86-74-8    | 18            | 4              | 22.22           | 3.80E-02      | 4.10E-02      | 5.30E-01             | J                      | 9.78E-01   | 5.52E-02         | 1.35E+01      | J              | GEO-33                  | 3.17E+00                          |                          |
| Chrysene                   | 218-01-9   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 2.40E-03             | J                      | 4.91E+00   | 1.16E-01         | 8.25E+01      | J              | GEO-33                  | 1.69E+01                          |                          |
| Dibenz(a,h)anthracene      | 53-70-3    | 24            | 12             | 50              | 3.80E-02      | 4.10E-02      | 1.70E-03             | J                      | 4.89E-01   | 4.58E-02         | 7.45E+00      | J              | GEO-33                  | 1.53E+00                          |                          |
| Dibenzofuran               | 132-64-9   | 18            | 6              | 33.33           | 3.80E-02      | 4.10E-02      | 3.90E-02             | J                      | 1.82E+00   | 5.87E-02         | 2.90E+01      | J              | GEO-33                  | 6.81E+00                          |                          |
| Fluoranthene               | 206-44-0   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 1.30E-02             | Z                      | 1.85E+01   | 1.82E-01         | 3.55E+02      | J              | GEO-33                  | 7.23E+01                          |                          |
| Fluorene                   | 86-73-7    | 24            | 8              | 33.33           | 2.90E-03      | 5.20E-02      | 3.60E-03             | J                      | 2.98E+00   | 4.72E-02         | 6.30E+01      | J              | GEO-33                  | 1.28E+01                          |                          |
| Indeno(1,2,3-cd)pyrene     | 193-39-5   | 24            | 13             | 54.17           | 3.80E-02      | 4.10E-02      | 7.80E-03             | J                      | 1.97E+00   | 9.68E-02         | 3.10E+01      | J              | GEO-33                  | 6.37E+00                          |                          |
| Naphthalene                | 91-20-3    | 24            | 6              | 25              | 2.90E-02      | 5.60E-01      | 7.50E-02             | J                      | 5.93E-01   | 6.22E-02         | 6.85E+00      | J              | GEO-33                  | 1.63E+00                          |                          |
| Phenanthrene               | 85-01-8    | 24            | 13             | 54.17           | 3.80E-02      | 4.10E-02      | 6.80E-03             | J                      | 1.21E+01   | 1.17E-01         | 2.45E+02      | J              | GEO-33                  | 5.00E+01                          |                          |
| Phenol                     | 108-95-2   | 18            | 13             | 72.22           | 7.50E-02      | 4.13E-01      | 1.00E-01             | J                      | 1.48E-01   | 1.22E-01         | 3.80E-01      | J              | GEO-29                  | 8.68E-02                          |                          |
| Pyrene                     | 129-00-0   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 1.60E-02             | J                      | 1.39E+01   | 1.87E-01         | 2.60E+02      | J              | GEO-33                  | 5.30E+01                          |                          |

**Table 15**  
**Statistical Summary and Selection of COPCs in EU5 Soil (0-6' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|----------------------------|------------------|---------------------------------|-----------------------------------|--|--|-----------------------------------|--------------------------|
| Semivolatiles              |                  |                                 |                                   |  |  |                                   |                          |
| 2,4-dimethylphenol         | 8.47E-02         | 7.83E-02                        | Unknown                           | 7.85E-02                                 | 4.08E+04                               | no                                | no                       |
| 2-methylnaphthalene        | 1.50E+00         | 1.43E+00                        | Unknown                           | 1.43E+00                                 | 8.18E+04                               | no                                | no                       |
| 2-methylphenol             | 4.01E-02         | 3.65E-02                        | Unknown                           | 3.69E-02                                 | 1.02E+05                               | no                                | no                       |
| 3- and 4-methylphenol      | 8.46E-02         | 7.90E-02                        | Unknown                           | 7.99E-02                                 | 1.02E+04                               | no                                | no                       |
| Acenaphthene               | 6.35E+00         | 3.21E+00                        | Unknown                           | 3.21E+00                                 | 1.23E+05                               | no                                | no                       |
| Acenaphthylene             | 2.10E+00         | 2.70E+00                        | Unknown                           | 2.70E+00                                 | 1.23E+05                               | no                                | no                       |
| Anthracene                 | 9.62E+00         | 6.15E+01                        | Unknown                           | 6.15E+01                                 | 6.13E+05                               | no                                | yes - COPC               |
| Benz(a)anthracene          | 1.08E+01         | 7.77E+01                        | Unknown                           | 7.77E+01                                 | 7.84E+00                               | yes - COPC                        | yes - COPC               |
| Benz(a)pyrene              | 6.93E+00         | 4.10E+01                        | Unknown                           | 4.10E+01                                 | 7.84E+01                               | yes - COPC                        | yes - COPC               |
| Benz(b)fluoranthene        | 1.07E+01         | 1.30E+02                        | Unknown                           | 7.95E+01                                 | 7.84E+00                               | yes - COPC                        | yes - COPC               |
| Benz(ghi)perylene          | 3.38E+00         | 8.53E+00                        | Unknown                           | 8.53E+00                                 | 6.13E+04                               | no                                | no                       |
| Benz(k)fluoranthene        | 3.89E+00         | 1.97E+01                        | Unknown                           | 1.97E+01                                 | 7.84E+01                               | no                                | yes*                     |
| Bis(2-ethylhexyl)phthalate | 9.21E+02         | 9.19E-02                        | Unknown                           | 9.19E-02                                 | 4.09E+02                               | no                                | yes*                     |
| Carbazole                  | 2.28E+00         | 4.56E+00                        | Unknown                           | 4.56E+00                                 | 2.86E+02                               | no                                | yes*                     |
| Chrysene                   | 1.08E+01         | 1.27E+02                        | Unknown                           | 8.25E+01                                 | 7.84E+02                               | no                                | yes - COPC               |
| Dibenz(a,h)anthracene      | 1.02E+00         | 2.04E+00                        | Unknown                           | 2.04E+00                                 | 7.84E-01                               | yes - COPC                        | yes - COPC               |
| Dibenzofuran               | 4.61E+00         | 4.75E+00                        | Unknown                           | 4.75E+00                                 | 8.18E+03                               | no                                | no                       |
| Fluoranthene               | 4.38E+01         | 7.28E+02                        | Unknown                           | 3.55E+02                                 | 8.17E+04                               | no                                | no                       |
| Fluorene                   | 7.47E+00         | 7.47E+00                        | Unknown                           | 7.47E+00                                 | 8.17E+04                               | no                                | yes - COPC               |
| Indeno(1,2,3-cd)pyrene     | 4.20E+00         | 1.71E+01                        | Unknown                           | 1.71E+01                                 | 7.84E+00                               | yes - COPC                        | yes - COPC               |
| Naphthalene                | 1.17E+00         | 1.53E+00                        | Unknown                           | 1.53E+00                                 | 8.24E+02                               | no                                | no                       |
| Phenanthrene               | 2.96E+01         | 1.26E+02                        | Unknown                           | 1.26E+02                                 | 6.13E+04                               | no                                | no                       |
| Phenol                     | 1.84E+01         | 2.27E-01                        | Normal/Lognormal                  | 2.27E-01                                 | 1.23E+05                               | no                                | no                       |
| Pyrene                     | 3.24E+01         | 4.64E+02                        | Unknown                           | 2.60E+02                                 | 6.13E+04                               | no                                | no                       |

\* Logarithmic 95% UCL is less than benchmark but retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 1

Statistical Summary and Selection of COPCs in EU5 Soil (0-20' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent                | CAS Number | Total Number of Samples | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Minimum Detected mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|-------------------------|-----------------|---------------|---------------|-----------------------|------------------------|------------|------------------------|------------------------|-----------------------------|-----------------------------------|--------------------------|
|                            |            |                         |                 |               |               |                       |                        |            |                        |                        |                             |                                   |                          |
| <b>Semi-volatiles</b>      |            |                         |                 |               |               |                       |                        |            |                        |                        |                             |                                   |                          |
| 2,4-dimethylphenol         | 105-07-9   | 23                      | 1               | 4.35          | 6.70E-02      | 1.30E+00              | 1.10E-01               | J          | 8.81E-02               | 5.57E-02               | 1.10E-01                    | J                                 | GEO-30                   |
| 2-methylnaphthalene        | 91-57-6    | 23                      | 8               | 34.78         | 3.30E-02      | 4.10E-02              | 5.10E-02               | J          | 2.37E+01               | 4.40E+02               | 1.19E-01                    | J                                 | SB-05                    |
| 2-methylphenol             | 95-48-7    | 23                      | 1               | 4.35          | 3.80E-02      | 1.30E+00              | 4.20E-02               | J          | 6.29E-02               | 3.18E-02               | 4.20E-02                    | J                                 | GEO-30                   |
| 3- and 4-methylphenol      | 106-44-5   | 23                      | 1               | 4.35          | 7.50E-02      | 2.00E+00              | 1.40E-01               | J          | 1.10E-01               | 6.12E-02               | 1.40E-01                    | J                                 | GEO-30                   |
| Acenaphthene               | 83-32-9    | 30                      | 7               | 23.33         | 2.90E-02      | 2.40E+00              | 1.0E-01                | J          | 1.29E+01               | 1.15E-01               | 2.90E+02                    | J                                 | SB-05                    |
| Acenaphthylene             | 208-96-8   | 30                      | 9               | 30            | 3.30E-02      | 2.40E+00              | 4.80E-02               | J          | 1.15E+00               | 9.90E-02               | 1.60E-01                    | J                                 | GEO-33                   |
| A-ithracene                | 120-12-7   | 30                      | 9               | 30            | 5.40E-04      | 6.00E-02              | 1.30E-01               | J          | 7.18E+00               | 6.99E-02               | 9.80E+01                    | J                                 | SB-05                    |
| Benzof[a]anthracene        | 56-55-3    | 30                      | 17              | 56.67         | 3.30E-02      | 4.10E-02              | 6.80E-03               | 2          | 6.52E+00               | 1.56E-01               | 8.35E+01                    | J                                 | GEO-33                   |
| Benzof[a]pyrene            | 50-32-8    | 30                      | 17              | 56.67         | 3.80E-02      | 6.70E-02              | 8.30E-03               | 2          | 3.56E+00               | 1.55E-01               | 5.25E+01                    | J                                 | GEO-33                   |
| Benzof[b]fluoranthene      | 205-99-2   | 30                      | 18              | 60            | 3.80E-02      | 6.70E-02              | 9.00E-03               | 2          | 5.47E+00               | 2.09E-01               | 7.95E+01                    | J                                 | GEO-33                   |
| Benzof[g]perylene          | 191-24-2   | 30                      | 16              | 53.33         | 3.80E-02      | 6.70E-02              | 6.70E-03               | J          | 1.51E+00               | 9.66E-02               | 2.55E+01                    | J                                 | GEO-33                   |
| Benzof[k]fluoranthene      | 207-08-9   | 30                      | 18              | 60            | 3.80E-02      | 1.30E-01              | 4.70E-03               | 2          | 1.99E+00               | 1.32E-01               | 2.85E+01                    | J                                 | GEO-33                   |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 23                      | 2               | 8.7           | 6.70E-02      | 1.30E+00              | 1.40E-01               | J          | 9.40E-02               | 6.04E-02               | 1.50E-01                    | J                                 | GEO-32                   |
| Carbazole                  | 86-74-8    | 23                      | 7               | 30.43         | 3.30E-02      | 4.10E-02              | 5.30E-01               | J          | 4.29E+00               | 9.66E-02               | 6.90E-01                    | J                                 | SB-05                    |
| Chrysene                   | 218-01-9   | 30                      | 18              | 60            | 3.30E-02      | 4.10E-02              | 2.40E-03               | J          | 6.33E+00               | 1.59E-01               | 8.23E+01                    | J                                 | GEO-33                   |
| Dibenz(a,l)anthracene      | 53-70-3    | 30                      | 15              | 50            | 3.80E-02      | 3.30E-01              | 1.70E-03               | J          | 4.87E-01               | 5.48E-02               | 7.45E+00                    | J                                 | GEO-33                   |
| Dibenzofuran               | 132-64-9   | 23                      | 9               | 39.13         | 3.30E-02      | 4.10E-02              | 3.90E-02               | J          | 1.47E+01               | 1.23E-01               | 2.70E+02                    | J                                 | SB-05                    |
| Fluoranthene               | 206-44-0   | 30                      | 18              | 60            | 3.30E-02      | 4.10E-02              | 1.30E-02               | 2          | 3.13E+01               | 2.82E-01               | 4.30E+02                    | J                                 | SB-05                    |
| Fluorene                   | 86-73-7    | 30                      | 11              | 36.67         | 2.90E-03      | 5.20E-02              | 3.60E-03               | J          | 1.52E+01               | 8.67E-02               | 3.30E+02                    | J                                 | SB-05                    |
| Indeno[1,2,3-cd]pyrene     | 193-39-5   | 30                      | 17              | 56.67         | 3.80E-02      | 6.70E-02              | 7.80E-03               | J          | 1.92E+00               | 1.22E-01               | 3.10E+01                    | J                                 | GEO-33                   |
| Naphthalene                | 91-20-3    | 30                      | 9               | 30            | 2.90E-02      | 5.60E-01              | 7.50E-02               | J          | 3.80E+01               | 1.33E-01               | 9.10E+02                    | J                                 | SB-05                    |
| Phenanthrene               | 85-01-8    | 30                      | 17              | 56.67         | 3.30E-02      | 4.10E-02              | 6.80E-03               | J          | 3.77E+01               | 2.06E+02               | 7.10E+02                    | J                                 | SB-05                    |
| Phenol                     | 108-95-2   | 23                      | 13              | 56.52         | 3.30E-02      | 6.70E-01              | 1.00E-01               | J          | 1.36E-01               | 9.66E-02               | 3.80E+01                    | J                                 | GEO-29                   |
| Pyrene                     | 129-00-0   | 30                      | 18              | 60            | 3.80E-02      | 6.70E-02              | 1.60E-02               | J          | 2.04E+01               | 2.85E-01               | 2.60E+02                    | J                                 | GEO-33                   |
| <b>Volatiles</b>           |            |                         |                 |               |               |                       |                        |            |                        |                        |                             |                                   |                          |
| Acetone                    | 67-64-1    | 5                       | 5               | 100           | 0.00E+00      | 0.00E+00              | 9.00E-03               | J          | 4.40E-02               | 2.95E-02               | 1.00E-01                    | J                                 | SB-05                    |
| Benzene                    | 71-43-2    | 5                       | 2               | 40            | 1.00E-03      | 1.00E-03              | 5.00E-03               | J          | 2.70E-03               | 1.34E-03               | 7.00E-03                    | J                                 | SB-05                    |
| Ethylbenzene               | 100-41-4   | 5                       | 3               | 60            | 1.00E-03      | 1.00E-03              | 2.40E-02               | J          | 3.82E-02               | 8.02E-03               | 1.20E-01                    | J                                 | GEO-02                   |
| Styrene                    | 100-42-5   | 5                       | 1               | 20            | 1.00E-03      | 1.00E-03              | 1.00E-01               | J          | 2.04E-02               | 1.44E-03               | 1.00E-01                    | J                                 | SB-05                    |
| Toluene                    | 108-88-3   | 5                       | 3               | 60            | 1.00E-03      | 1.00E-03              | 1.30E-02               | J          | 3.38E-02               | 5.85E-03               | 1.40E-01                    | J                                 | SB-05                    |
| Xylene (total)             | 1330-20-7  | 5                       | 3               | 60            | 1.00E-03      | 1.00E-03              | 7.50E-02               | J          | 2.27E-01               | 2.10E-02               | 7.80E-01                    | J                                 | SB-05                    |

**Table 16**  
**Statistical Summary and Selection of COPCs in EU5 Soil (0-20' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG?<br>> TRG? |
|----------------------------|------------------|---------------------------------|-----------------------------------|--|--|---|
| <b>Semi-volatiles</b>      |                  |                                 |                                   |  |  |   |
| 2,4-dimethylphenol         | 1.36E-01         | 1.12E-01                        | Unknown                           | 1.10E-01                                 | 4.08E+04                               | no  |
| 2-methylnaphthalene        | 5.67E+01         | 6.89E+02                        | Unknown                           | 4.40E+02                                 | 8.18E+04                               | no  |
| 2-methylphenol             | 1.11E-01         | 7.64E-02                        | Unknown                           | 4.20E-02                                 | 1.02E+05                               | no  |
| 3- and 4-methylphenol      | 1.83E-01         | 1.37E-01                        | Unknown                           | 1.37E-01                                 | 1.02E+04                               | no  |
| Acenaphthene               | 2.95E+01         | 8.04E+01                        | Unknown                           | 8.04E+01                                 | 1.23E+05                               | no  |
| Acenaphthylene             | 2.19E+00         | 3.82E+00                        | Unknown                           | 3.82E+00                                 | 1.23E+05                               | no  |
| Anthracene                 | 1.42E+01         | 4.31E+02                        | Unknown                           | 9.80E+01                                 | 6.13E+05                               | no  |
| Benz(a)anthracene          | 1.25E+01         | 1.52E+02                        | Unknown                           | 8.35E+01                                 | 7.84E+00                               | YES-COPC                                    |
| Benz(a)pyrene              | 6.82E+00         | 4.42E+01                        | Unknown                           | 4.42E+01                                 | 7.84E-01                               | YES-COPC                                    |
| Benz(b)fluoranthene        | 1.04E+01         | 1.19E+02                        | Unknown                           | 7.95E+01                                 | 7.84E+00                               | YES-COPC                                    |
| Benzofluoroperylene        | 2.99E+00         | 7.40E+00                        | Unknown                           | 7.40E+00                                 | 6.13E+04                               | no  |
| Benzofluoranthene          | 3.74E+00         | 1.68E+01                        | Unknown                           | 1.68E+01                                 | 7.84E+01                               | no  |
| Bis(2-ethylhexyl)phthalate | 1.42E-01         | 1.24E-01                        | Unknown                           | 1.24E-01                                 | 2.86E+02                               | no  |
| Carbazole                  | 9.49E+00         | 6.44E+01                        | Unknown                           | 6.44E+01                                 | 7.84E+02                               | no  |
| Chrysene                   | 1.21E+01         | 2.00E+02                        | Unknown                           | 8.25E+01                                 | 7.84E-01                               | YES-COPC                                    |
| Dibenz(a,h)anthracene      | 9.29E-01         | 1.53E+00                        | Unknown                           | 1.53E+00                                 | 8.18E+03                               | YES*  |
| Dibenzo-furan              | 3.48E+01         | 4.46E+02                        | Unknown                           | 4.46E+02                                 | 8.17E+04                               | no  |
| Fluoranthene               | 6.22E+01         | 3.34E+03                        | Unknown                           | 4.39E+02                                 | 8.17E+04                               | no  |
| Fluorene                   | 3.42E+01         | 2.24E+02                        | Unknown                           | 2.24E+02                                 | 7.84E+00                               | yes   |
| Indeno(1,2,3-cd)pyrene     | 3.73E+00         | 1.32E+01                        | Unknown                           | 1.32E+01                                 | 8.24E+02                               | no  |
| Naphthalene                | 9.01E+01         | 2.69E+02                        | Unknown                           | 2.69E+02                                 | 1.17E+03                               | no  |
| Phenanthrene               | 7.98E+01         | 2.37E+03                        | Unknown                           | 7.10E+02                                 | 6.13E+04                               | no  |
| Phenol                     | 1.72E-01         | 2.53E-01                        | Normal/Lognormal                  | 2.53E-01                                 | 1.23E+05                               | no  |
| Pyrene                     | 4.03E+01         | 1.08E+03                        | Unknown                           | 2.60E+02                                 | 6.13E+04                               | no  |
| <b>Volatiles</b>           |                  |                                 |                                   |  |  |   |
| Acetone                    | 8.01E-02         | 9.07E-01                        | Normal/Lognormal                  | 1.00E-01                                 | 1.04E+05                               | no  |
| Benzene                    | 5.65E-03         | 2.77E-01                        | Normal/Lognormal                  | 7.00E-03                                 | 1.36E+00                               | no  |
| Ethylbenzene               | 8.54E-02         | 1.58E+06                        | Normal/Lognormal                  | 1.20E-01                                 | 3.95E+02                               | no  |
| Styrene                    | 6.28E-02         | 1.19E+04                        | Unknown                           | 1.00E-01                                 | 3.84E+02                               | no  |
| Toluene                    | 9.08E-02         | 1.16E+05                        | Lognormal                         | 1.40E-01                                 | 3.80E+01                               | no  |
| Xylene (total)             | 5.41E-01         | 2.82E+13                        | Normal/Lognormal                  | 7.80E-01                                 | 3.18E+02                               | no  |

\*Logarithmic 95% UCL is less than benchmark but retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

**Table 17**  
**Statistical Summary and Selection of COPCs in EU6 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                   | CAS Number | Total Number of Samples | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Maximum Qualifier mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|-------------------------------|------------|-------------------------|-----------------|---------------|---------------|-----------------------|------------|------------------------|------------------------|-------------------------|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>          |            |                         |                 |               |               |                       |            |                        |                        |                         |                                   |                          |
| 1,2,4-trichlorobenzene        | 120-82-1   | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 1,2-dichlorobenzene           | 95-50-1    | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 1,3-dichlorobenzene           | 541-73-1   | 3                       | 1               | 35.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 1,4-dichlorobenzene           | 106-46-7   | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 2,2'-oxybis (1-chloropropane) | 108-60-1   | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 2,4,5-trichlorophenol         | 95-95-4    | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 2,4,6-trichlorophenol         | 88-06-2    | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 2,4-dichlorophenol            | 120-83-2   | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 2,4-dimethylphenol            | 105-67-9   | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 2,4-dinitrophenol             | 51-28-5    | 3                       | 1               | 33.33         | 2.40E-01      | 2.50E-01              | 2.30E+00   | 8.48E-01               | 3.26E-01               | 2.30E+00                | SD-04                             | 1.26E+00                 |
| 2,4-dinitrotoluene            | 121-14-2   | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 1.62E-01               | 8.94E-02               | 4.00E-01                | SD-04                             | 2.07E-01                 |
| 2,6-dinitrotoluene            | 606-20-2   | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 2-chloronaphthalene           | 91-58-7    | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 2-chlorophenol                | 95-57-8    | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 2-methylnaphthalene           | 91-57-6    | 3                       | 3               | 100           | 0.00E+00      | 0.00E+00              | 9.10E-02   | J                      | 1.28E+01               | 1.15E+01                | SD-04                             | 2.18E+01                 |
| 2-methylphenol                | 95-48-7    | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 2-nitroaniline*               | 88-74-4    | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 2-nitrophenol*                | 88-75-5    | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 3- and 4-methylphenol         | 106-44-5   | 3                       | 3               | 100           | 0.00E+00      | 0.00E+00              | 9.30E-02   | J                      | 3.34E-01               | 2.02E-01                | SD-04                             | 4.03E-01                 |
| 3,3'-dichlorobenzidine        | 91-94-1    | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 3-nitroaniline*               | 99-09-2    | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 4,6-dinitro-2-methylphenol    | 534-52-1   | 3                       | 1               | 33.33         | 2.10E-01      | 2.10E-01              | 2.00E+00   | 7.37E-01               | 2.80E-01               | 2.00E+00                | SD-04                             | 1.09E+00                 |
| 4-bromophenylphenylether*     | 101-55-3   | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 4-chloro-3-methylphenol*      | 59-50-7    | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 4-chloronaniline              | 106-47-8   | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 4-chlorophenylphenylether*    | 7005-72-3  | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |
| 4-nitroaniline*               | 100-01-6   | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| 4-nitrophenoxy*               | 100-02-7   | 3                       | 1               | 33.33         | 2.10E-01      | 2.10E-01              | 2.00E+00   | 7.37E-01               | 2.80E-01               | 2.00E+00                | SD-04                             | 1.09E+00                 |
| Acenaphthene                  | 83-32-9    | 8                       | 4               | 50            | 1.80E+00      | 3.50E+00              | 1.00E-01   | J                      | 1.89E+01               | 1.96E+00                | J                                 | 4.90E+01                 |
| Acenaphthylene                | 208-96-8   | 8                       | 3               | 37.5          | 1.80E+00      | 3.50E+00              | 1.70E-01   | J                      | 2.92E+00               | 1.69E+00                | J                                 | 3.14E+00                 |
| Anthracene                    | 120-12-7   | 8                       | 7               | 87.5          | 4.57E-01      | 4.57E-01              | 8.80E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | Z                                 | SD-16                    |
| Benz(a)anthracene             | 56-55-3    | 8                       | 8               | 100           | 0.00E+00      | 0.00E+00              | 9.30E-01   | 1.85E+01               | 4.83E+00               | 1.00E+02                | J                                 | SD-04                    |
| Benz(a)pyrene                 | 50-32-8    | 8                       | 8               | 100           | 0.00E+00      | 0.00E+00              | 9.70E-01   | J                      | 1.32E+01               | 5.59E+00                | J                                 | SD-03                    |
| Benz(b)fluoranthene           | 205-99-2   | 8                       | 8               | 100           | 0.00E+00      | 0.00E+00              | 1.40E+00   | 1.90E+01               | 7.10E+00               | 7.80E+01                | SD-03                             | 1.80E+01                 |
| Benz(h)perylene               | 191-24-2   | 8                       | 8               | 100           | 0.00E+00      | 0.00E+00              | 4.20E-01   | 6.50E+00               | 2.58E+00               | 3.20E+01                | SD-03                             | 1.07E+01                 |
| Benz(k)fluoranthene           | 207-08-9   | 8                       | 8               | 100           | 0.00E+00      | 0.00E+00              | 5.00E-01   | 6.96E+00               | 3.14E+00               | 2.30E+01                | SD-03                             | 8.76E+00                 |
| Bis(2-chloroethoxy)methane*   | 111-91-1   | 3                       | 1               | 33.33         | 8.40E-02      | 8.50E-02              | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01                | SD-04                             | 4.37E-01                 |
| Bis(2-chloroethyl)ether       | 111-44-4   | 3                       | 1               | 33.33         | 4.20E-02      | 4.30E-02              | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01                | SD-04                             | 2.19E-01                 |

*Table I-7*  
*Statistical Summary and Selection of COPCs in EU6 Sediment*  
*Kerr McGee, Hattiesburg, MS*

| Constituent                  | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>Unrestricted Soil<br>TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? |                          |
|------------------------------|------------------|---------------------------------|-----------------------------------|--|---|-----------------------------------|--------------------------|
|                              |                  |                                 |                                   |  |   | Is the 95% UCL ><br>TRG?          | Is the 95% UCL ><br>TRG? |
| 1,2,4-trichlorobenzene       | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 5.27E+02                                    | no                                | no                       |
| 1,2-dichlorobenzene          | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 2.79E+02                                    | no                                | no                       |
| 1,3-dichlorobenzene          | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 2.35E+03                                    | no                                | no                       |
| 1,4-dichlorobenzene          | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 2.66E+01                                    | no                                | no                       |
| 2,2'-oxybis(1-chloropropane) | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 5.93E+00                                    | no                                | no                       |
| 2,4,5-trichlorophenol        | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | 7.82E+03                                    | no                                | no                       |
| 2,4,6-trichlorophenol        | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | 5.81E+01                                    | no                                | no                       |
| 2,4-dichlorophenol           | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | 2.35E+02                                    | no                                | no                       |
| 2,4-dimethylphenol           | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | 1.56E+03                                    | no                                | no                       |
| 2,4-dinitrophenol            | 2.97E+00         | 4.39E+11                        | Lognormal                         | 2.30E+00                                 | 1.56E+02                                    | no                                | no                       |
| 2,4-dinitrothiophene         | 5.10E-01         | 1.19E+06                        | Lognormal                         | 4.00E-01                                 | 1.56E+02                                    | no                                | no                       |
| 2,6-dinitrothiophene         | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 7.82E+01                                    | no                                | no                       |
| 2-chloronaphthalene          | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 6.26E+03                                    | no                                | no                       |
| 2-chlorophenol               | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 3.91E+02                                    | no                                | no                       |
| 2-methylnaphthalene          | 4.96E+01         | 3.37E+41                        | Normal/Lognormal                  | 3.80E+01                                 | 3.13E+03                                    | no                                | no                       |
| 2-methylphenol               | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 3.91E+03                                    | no                                | no                       |
| 2-nitroaniline*              | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | NA  | no                                | no                       |
| 2-nitrophenol*               | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | NA  | no                                | no                       |
| 3- and 4-methylphenol        | 1.01E+00         | 2.30E+05                        | Normal/Lognormal                  | 8.00E-01                                 | 3.91E+02                                    | no                                | no                       |
| 3,3'-dichlorobenzidine       | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | 1.42E+00                                    | no                                | no                       |
| 3-nitroaniline*              | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | NA  | no                                | no                       |
| 4,6-dinitro-2-methylphenol   | 2.58E+00         | 4.94E+11                        | Unknown                           | 2.00E+00                                 | 7.82E+00                                    | no                                | no                       |
| 4-bromophenylphenylether*    | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | NA  | no                                | no                       |
| 4-chloro-3-methylphenol*     | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | NA  | no                                | no                       |
| 4-chloroaniline              | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 3.13E+02                                    | no                                | no                       |
| 4-chlorophenylphenylether*   | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | NA  | no                                | no                       |
| 4-nitroaniline*              | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | NA  | no                                | no                       |
| 4-nitrophenol!               | 2.58E+00         | 4.94E+11                        | Unknown                           | 2.00E+00                                 | 6.26E+02                                    | no                                | no                       |
| Acenaphthene                 | 5.17E+01         | 1.82E+03                        | Lognormal                         | 1.40E+02                                 | 4.69E+03                                    | no                                | no                       |
| Acenaphthylene               | 5.02E+00         | 2.23E+01                        | Normal/Lognormal                  | 8.90E+00                                 | 4.69E+03                                    | no                                | no                       |
| Anthracene                   | 1.13E+01         | 1.03E+02                        | Lognormal                         | 2.39E+01                                 | 2.35E+04                                    | no                                | no                       |
| Benz(a)anthracene            | 4.14E+01         | 7.51E+02                        | Lognormal                         | 1.00E+02                                 | 8.75E-01                                    | YES                               | YES - COPC               |
| Benz(a)pyrene                | 2.53E+01         | 1.63E+02                        | Lognormal                         | 4.90E+01                                 | 8.75E-02                                    | YES                               | YES - COPC               |
| Benz(b)fluoranthene          | 3.79E+01         | 2.88E+02                        | Lognormal                         | 7.80E+01                                 | 8.75E-01                                    | YES                               | YES - COPC               |
| Benz(g)phenylenes            | 1.37E+01         | 6.89E+01                        | Lognormal                         | 3.20E+01                                 | 2.35E+03                                    | no                                | no                       |
| Benz(k)fluoranthene          | 1.28E+01         | 8.18E+01                        | Lognormal                         | 2.30E+01                                 | 8.75E+00                                    | YES                               | YES - COPC               |
| Bis(2-chloroethoxy)methane*  | 1.03E+00         | 1.77E+11                        | Unknown                           | 8.00E-01                                 | NA  | no                                | no                       |
| Bis(2-chloroethyl)ether      | 5.16E-01         | 7.97E+10                        | Lognormal                         | 4.00E-01                                 | 2.73E-01                                    | YES                               | YES - COPC               |

**Table I7**  
**Statistical Summary and Selection of COPCs in EU6 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | CAS Number | Total Number of Samples | Hit Hits | Frequency % | Minimum mg/kg | Maximum mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Maximum Detected mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |          |
|----------------------------|------------|-------------------------|----------|-------------|---------------|---------------|------------|------------------------|------------------------|------------------------|-----------------------------------|--------------------------|----------|
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 3                       | 1        | 33.33       | 1.50E-01      | 2.50E-01      | 8.80E-01   | 3.60E-01               | 2.02E+01               | 8.80E-01               | SD-04                             | 4.51E-01                 |          |
| Butylbenzylphthalate       | 85-68-7    | 3                       | 1        | 33.33       | 8.40E-02      | 8.50E-02      | 8.00E-01   | 2.95E-01               | 1.13E+01               | 8.00E-01               | SD-04                             | 4.37E-01                 |          |
| Carbazole                  | 86-74-8    | 3                       | 3        | 100         | 0.00E+00      | 0.00E+00      | 2.20E-01   | 3.37E+01               | 2.77E+00               | 1.00E+02               | J                                 | 5.74E-01                 |          |
| Chrysene                   | 218-01-9   | 8                       | 6        | 75          | 9.40E-01      | 1.20E+00      | 1.30E+00   | 1.82E+01               | 4.61E+00               | 7.60E+01               | J                                 | 2.75E-01                 |          |
| Dibenz(a,h)anthracene      | 53-70-3    | 8                       | 8        | 100         | 0.00E+00      | 0.00E+00      | 1.50E-01   | 2.03E+00               | 8.14E+01               | 9.60E+00               | SD-04                             | 3.23E+00                 |          |
| Dibenzofuran               | 132-64-9   | 3                       | 3        | 100         | 0.00E+00      | 0.00E+00      | 1.00E-01   | 5.02E+01               | 1.93E+00               | 1.50E+02               | J                                 | 8.64E-01                 |          |
| Diethylphthalate           | 84-66-2    | 3                       | 1        | 33.33       | 8.40E-02      | 8.50E-02      | 8.00E-01   | 2.95E-01               | 1.13E+01               | 8.00E-01               | SD-04                             | 4.37E-01                 |          |
| Dimethylphthalate          | 131-11-3   | 3                       | 1        | 33.33       | 8.40E-02      | 8.50E-02      | 8.00E-01   | 2.95E-01               | 1.13E+01               | 8.00E-01               | SD-04                             | 4.37E-01                 |          |
| Di-n-butylphthalate        | 84-74-2    | 3                       | 1        | 33.33       | 8.40E-02      | 8.50E-02      | 8.00E-01   | 2.95E-01               | 1.13E+01               | 8.00E-01               | SD-04                             | 4.37E-01                 |          |
| Di-n-octylphthalate        | 117-84-0   | 3                       | 1        | 33.33       | 8.40E-02      | 8.50E-02      | 8.00E-01   | 2.95E-01               | 1.13E+01               | 8.00E-01               | SD-04                             | 4.37E-01                 |          |
| Fluoranthene               | 206-44-0   | 8                       | 8        | 100         | 0.00E+00      | 0.00E+00      | 9.20E-01   | Z                      | 6.71E+01               | 8.91E+00               | 4.70E+02                          | SD-04                    | 1.63E+02 |
| Fluorene                   | 86-73-7    | 8                       | 6        | 75          | 3.20E-01      | 3.30E-01      | 1.80E-01   | J                      | 3.32E+01               | 1.07E+00               | 2.60E+02                          | SD-04                    | 9.16E+01 |
| Hexachlorobenzene          | 118-74-1   | 3                       | 1        | 33.33       | 4.20E-02      | 4.30E-02      | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01               | SD-04                             | 2.19E+01                 |          |
| Hexachlorobutadiene        | 87-68-3    | 3                       | 1        | 33.33       | 8.40E-02      | 8.50E-02      | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01               | SD-04                             | 4.37E+01                 |          |
| Hexachlorocyclopentadiene  | 77-47-4    | 3                       | 1        | 33.33       | 2.10E-01      | 2.10E-01      | 2.00E+00   | 7.37E-01               | 2.80E-01               | 2.00E+00               | SD-04                             | 1.09E+00                 |          |
| Hexachloroethane           | 67-72-1    | 3                       | 1        | 33.33       | 4.20E-02      | 4.30E-02      | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01               | SD-04                             | 2.19E+01                 |          |
| Indeno(1,2,3-cd)pyrene     | 193-39-5   | 8                       | 8        | 100         | 0.00E+00      | 0.00E+00      | 5.40E-01   | 8.32E+00               | 3.53E+00               | 3.90E+01               | SD-03                             | 1.30E+01                 |          |
| Isophorone                 | 78-59-1    | 3                       | 1        | 33.33       | 4.20E-02      | 4.30E-02      | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01               | SD-04                             | 2.19E+01                 |          |
| Naphthalene                | 91-20-3    | 8                       | 3        | 37.5        | 1.80E+00      | 3.50E+00      | 1.60E-01   | 2.91E+00               | 1.48E+00               | 1.40E+01               | SD-04                             | 4.52E+00                 |          |
| Nitrobenzene               | 98-95-3    | 3                       | 1        | 33.33       | 4.20E-02      | 4.30E-02      | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01               | SD-04                             | 2.19E+01                 |          |
| N-nitrosodi-n-propylamine  | 621-64-7   | 3                       | 1        | 33.33       | 4.20E-02      | 4.30E-02      | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01               | SD-04                             | 2.19E+01                 |          |
| N-nitrosodiphenylamine     | 86-30-6    | 3                       | 1        | 33.33       | 4.20E-02      | 4.30E-02      | 4.00E-01   | 1.48E-01               | 5.65E-02               | 4.00E-01               | SD-04                             | 2.19E+01                 |          |
| Pentachlorophenol          | 87-86-5    | 3                       | 1        | 33.33       | 2.10E-01      | 2.10E-01      | 2.00E+00   | 7.37E-01               | 2.80E-01               | 2.00E+00               | SD-04                             | 1.09E+00                 |          |
| Phenanthrene               | 85-01-8    | 8                       | 5        | 62.5        | 1.30E-01      | 1.20E+00      | 6.60E-01   | 1.10E+02               | 1.84E+00               | 8.70E+02               | SD-04                             | 3.07E+02                 |          |
| Phenol                     | 108-95-2   | 3                       | 1        | 33.33       | 8.40E-02      | 8.50E-02      | 8.00E-01   | 2.95E-01               | 1.13E-01               | 8.00E-01               | SD-04                             | 4.37E+01                 |          |
| Pyrene                     | 129-00-0   | 8                       | 8        | 100         | 0.00E+00      | 0.00E+00      | 1.60E+00   | 4.88E+01               | 1.16E+01               | 3.00E+02               | SD-04                             | 1.02E+02                 |          |

*Table 17*  
*Statistical Summary and Selection of COPCs in EU6 Sediment*  
*Kerr McGee, Hattiesburg, MS*

| Constituent                        | Tier I           |                  |                                | Tier I                                   |                                |   |
|------------------------------------|------------------|------------------|--------------------------------|--|--------------------------------|---|
|                                    | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Unrestricted Soil<br>TRG mg/kg | Is the Maximum<br>Detected > TRG?<br>no |
| Bis(2-ethylhexyl)phthalate         | 1.12E+00         | 2.81E+06         | Normal/Lognormal               | 8.80E-01                                 | 4.56E+01                       | no                                      |
| Butylbenzylphthalate               | 1.03E+00         | 1.77E+11         | Unknown                        | 8.00E-01                                 | 9.28E+02                       | no                                      |
| Carbazole                          | 1.30E+02         | 3.75E+43         | Normal/Lognormal               | 1.00E+02                                 | 3.19E+01                       | YES                                     |
| Chrysene                           | 3.66E+01         | 2.31E+03         | Lognormal                      | 7.60E+01                                 | 8.75E+01                       | YES*                                    |
| Dibenz(a,h)anthracene              | 4.20E+00         | 2.18E+01         | Lognormal                      | 9.60E+00                                 | 8.75E-02                       | YES                                     |
| Dibenzofuran                       | 1.96E+02         | 1.09E+63         | Lognormal                      | 1.50E+02                                 | 3.13E+02                       | no                                      |
| Diethylphthalate                   | 1.03E+00         | 1.77E+11         | Unknown                        | 8.00E-01                                 | 1.97E+03                       | no                                      |
| Dimethylphthalate                  | 1.03E+00         | 1.77E+11         | Unknown                        | 8.00E-01                                 | 7.82E+05                       | no                                      |
| Di-n-butylphthalate                | 1.03E+00         | 1.77E+11         | Unknown                        | 8.00E-01                                 | 2.28E+03                       | no                                      |
| Di-n-octylphthalate                | 1.03E+00         | 1.77E+11         | Unknown                        | 8.00E-01                                 | 1.56E+03                       | no                                      |
| Fluoranthene                       | 1.76E+02         | 8.33E+03         | Lognormal                      | 4.70E+02                                 | 3.13E+03                       | no                                      |
| Fluorene                           | 9.46E+01         | 1.95E+04         | Lognormal                      | 2.60E+02                                 | 3.13E+03                       | no                                      |
| Hexachlorobenzene                  | 5.16E-01         | 7.97E+10         | Lognormal                      | 4.00E-01                                 | 3.99E-01                       | YES                                     |
| Hexachlorobutadiene                | 1.03E+00         | 1.77E+11         | Unknown                        | 8.00E-01                                 | 8.82E-02                       | YES                                     |
| Hexachlorocyclopentadiene          | 2.58E+00         | 4.94E+11         | Unknown                        | 2.00E+00                                 | 9.51E-01                       | YES                                     |
| Hexachloroethane                   | 5.16E-01         | 7.97E+10         | Lognormal                      | 4.00E-01                                 | 4.56E+01                       | no                                      |
| Indeno(1,2,3- <i>cd</i> )pyrene    | 1.70E+01         | 8.48E+01         | Lognormal                      | 3.90E+01                                 | 8.75E-01                       | YES                                     |
| Isoaphrone                         | 5.16E-01         | 7.97E+10         | Lognormal                      | 4.00E-01                                 | 6.72E+02                       | no                                      |
| Naphthalene                        | 5.93E+00         | 1.92E+01         | Lognormal                      | 1.40E+01                                 | 6.45E+02                       | no                                      |
| Nitrobenzene                       | 5.16E-01         | 7.97E+10         | Lognormal                      | 4.00E-01                                 | 8.41E+00                       | no                                      |
| N-nitrosodi- <i>n</i> -propylamine | 5.16E-01         | 7.97E+10         | Lognormal                      | 4.00E-01                                 | 9.12E-02                       | YES                                     |
| N-nitrosodiphenylamine             | 5.16E-01         | 7.97E+10         | Lognormal                      | 4.00E-01                                 | 1.30E+02                       | no                                      |
| Perachlorophenol                   | 2.58E+00         | 4.94E+11         | Unknown                        | 2.00E+00                                 | 2.66E+00                       | no                                      |
| Phenanthrene                       | 3.16E+02         | 1.66E+06         | Lognormal                      | 8.70E+02                                 | 2.35E+03                       | no                                      |
| Phenol                             | 1.03E+00         | 1.77E+11         | Unknown                        | 8.00E-01                                 | 4.69E+04                       | no                                      |
| Pyrene                             | 1.17E+02         | 1.89E+03         | Lognormal                      | 3.00E+02                                 | 2.35E+03                       | no                                      |
| NA - Not Available                 |                  |                  |                                |  |                                |   |

\* Constituent will be retained as a COPC due to lack of Tier I TRG.

\*\* Retained as a COPC, as per MDEQ Comments (8/2/2000): constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

\*\*\* Logarithmic 95% UCL is less than benchmark but retained as a COPC, as per MDEQ Comments (8/2/2000): constituent is a member of carcinogenic PAH family one of which has been retained as a COPC.

**Table 54**  
**Dermal Exposure to EU1 Sediment by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                                       | <u>Cs*SA*AH*ABS*EF*ED*CF</u><br>BW*AT |                                       |                 |  |          |          |
|---|---------------------------------------|---------------------------------------|---------------------------------------|-----------------|--|----------|----------|
| Cs - Concentration in sediment =                            | mg/kg                                 | chem. spec.                           |                                       |                 |  |          |          |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                  | 3000                                  | calculated                            |                 |  |          |          |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                       | 20000                                 | USEPA 1997, EFH                       |                 |  |          |          |
| Fs - Fraction of skin surface area available for exposure = |                                       | 15.0%                                 | USEPA 1997, EFH                       |                 |  |          |          |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                    | 0.13                                  | USEPA 1997, EFH                       |                 |  |          |          |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                                       | 0.03                                  | USEPA 1995, Region III                |                 |  |          |          |
| EF - Exposure frequency =                                   | days/year                             | 8                                     | reasonable assumption                 |                 |  |          |          |
| ED - Exposure duration =                                    | years                                 | 1                                     | reasonable assumption                 |                 |  |          |          |
| CF - Conversion factor =                                    | kg/mg                                 | 1.00E-06                              |                                       |                 |  |          |          |
| BW - Body weight =  | kg                                    | 70                                    | USEPA 1995, Region IV                 |                 |  |          |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                                  | 365                                   | USEPA 1991, HHEM                      |                 |  |          |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                                  | 25550                                 | USEPA 1991, HHEM                      |                 |  |          |          |
| Constituent   | Concentration in<br>Sediment<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day  | Dermal<br>Subchronic RFD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day |          |          |
|   |                                       |                                       |                                       |                 | Cancer Slope<br>1/(mg/kg-day)                    |          |          |
| Semivolatiles   |                                       |                                       |                                       |                 | Cancer<br>Risk                                   |          |          |
| Benzo(a)anthracene  | 5.90E-01                              | 2.16E-09                              | NA                                    | NA              | 3.09E-11   | 1.46E+00 | 4.51E-11 |
| Benzo(a)pyrene  | 3.90E-01                              | 1.43E-09                              | NA                                    | NA              | 2.04E-11   | 1.46E+01 | 2.98E-10 |
| Benzo(b)fluoranthene  | 5.80E-01                              | 2.12E-09                              | NA                                    | NA              | 3.04E-11   | 1.46E+00 | 4.43E-11 |
| Benzo(k)fluoranthene  | 1.90E-01                              | 6.96E-10                              | NA                                    | NA              | 9.94E-12   | 1.46E-01 | 1.45E-12 |
| Chrysene  | 5.30E-01                              | 1.94E-09                              | NA                                    | NA              | 2.77E-11   | 1.46E-02 | 4.05E-13 |
| Dibenz(a,h)anthracene                                       | 6.20E-02                              | 2.27E-10                              | NA                                    | NA              | 3.24E-12   | 1.46E+01 | 4.74E-11 |
| Indeno(1,2,3-cd)pyrene                                      | 2.20E-01                              | 8.06E-10                              | NA                                    | NA              | 1.15E-11   | 1.46E+00 | 1.68E-11 |

NA - Not Available

Total Cancer Risk = 4.53E-10



**Table 18**  
**Statistical Summary and Selection of COPCs in EU6 Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | CAS Number | Total Samples | Hit Frequency | Minimum mg/L | Maximum mg/L | Detection Limit | Detected | Minimum mg/L | Mean mg/L | Logarithmic Mean | Maximum mg/L | Detected | Detected | Maximum mg/L | Location of Maximum Concentration | Standard Deviation mg/L |
|------------------------|------------|---------------|---------------|--------------|--------------|-----------------|----------|--------------|-----------|------------------|--------------|----------|----------|--------------|-----------------------------------|-------------------------|
| <b>Semi挥发性</b>         |            |               |               |              |              |                 |          |              |           |                  |              |          |          |              |                                   |                         |
| Acenaphthene           | 83-32-9    | 2             | 1             | 50           | 1.00E-03     | 9.00E-03        | J        | 4.75E-03     | 2.12E-03  | 9.00E-03         | J            |          |          | SW-03        | 6.01E-03                          |                         |
| Benz(a)anthracene      | 56-55-3    | 2             | 0             | 0            | 1.00E-03     | 1.00E-03        | 0.00E+00 | NA           | 5.00E-04  | 5.00E-04         | 0.00E+00     | NA       |          | SW-03        | 0.00E+00                          |                         |
| Benz(a)pyrene          | 50-32-8    | 2             | 0             | 0            | 1.00E-03     | 1.00E-03        | 0.00E+00 | NA           | 5.00E-04  | 5.00E-04         | 0.00E+00     | NA       |          | SW-03        | 0.00E+00                          |                         |
| Benz(b)fluoranthene    | 205-99-2   | 2             | 1             | 50           | 1.00E-03     | 1.00E-03        | 9.00E-03 | J            | 4.75E-03  | 2.12E-03         | 9.00E-03     | J        |          | SW-03        | 6.01E-03                          |                         |
| Benz(k)fluoranthene    | 207-08-9   | 2             | 0             | 0            | 1.00E-03     | 1.00E-03        | 0.00E+00 | NA           | 5.00E-04  | 5.00E-04         | 0.00E+00     | NA       |          | SW-03        | 0.00E+00                          |                         |
| Chrysene               | 218-01-9   | 2             | 0             | 0            | 1.00E-03     | 1.00E-03        | 0.00E+00 | NA           | 5.00E-04  | 5.00E-04         | 0.00E+00     | NA       |          | SW-03        | 0.00E+00                          |                         |
| Dibenz(a,h)anthracene  | 53-70-3    | 2             | 0             | 0            | 1.00E-03     | 1.00E-03        | 0.00E+00 | NA           | 5.00E-04  | 5.00E-04         | 0.00E+00     | NA       |          | SW-03        | 0.00E+00                          |                         |
| Fluoranthene           | 206-44-0   | 2             | 2             | 100          | 0.00E+00     | 0.00E+00        | 1.20E-02 |              | 1.25E-02  | 1.25E-02         | 1.30E-02     |          |          | SW-03        | 7.07E-04                          |                         |
| Fluorene               | 86-73-7    | 2             | 1             | 50           | 1.00E-03     | 1.00E-03        | 1.10E-02 |              | 5.75E-03  | 2.35E-03         | 1.10E-02     |          |          | SW-03        | 7.42E-03                          |                         |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 2             | 0             | 0            | 1.00E-03     | 1.00E-03        | 0.00E+00 | NA           | 5.00E-04  | 5.00E-04         | 0.00E+00     | NA       |          | SW-03        | 0.00E+00                          |                         |

NA - Not applicable: constituent not detected in media.

**Table 18**  
**Statistical Summary and Selection of COPCs in EU6 Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            |                 |                                |                                   |   | Human Health                                     |                                   |  |
|------------------------|-----------------|--------------------------------|-----------------------------------|---|--|-----------------------------------|--|
|                        | 95% UCL<br>mg/L | Logarithmic<br>95% UCL<br>mg/L | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/L | Consumption of Water<br>& Organisms AWQC<br>mg/L | Is Maximum<br>Detected ><br>AWQC? |  |
| <b>Semi-volatiles</b>  |                 |                                |                                   |   |  |                                   |  |
| Aceanaphthene          | 3.16E-02        | 6.39E+48                       | Unknown                           | 9.00E-03                                | 1.20E+00   | no                                |  |
| Benz(a)anthracene      | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |  |
| Benzo(a)pyrene         | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |  |
| Benzo(b)fluoranthene   | 3.16E-02        | 6.39E+48                       | Unknown                           | 9.00E-03                                | 4.40E-06   | YES -COPC                         |  |
| Benzo(k)fluoranthene   | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |  |
| Chrysene               | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |  |
| Dibenz(a,h)anthracene  | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |  |
| Fluoranthene           | 1.57E-02        | 1.53E-02                       | Unknown                           | 1.30E-02                                | 3.00E-01   | no                                |  |
| Fluorene               | 3.89E-02        | 1.87E+56                       | Unknown                           | 1.10E-02                                | 1.30E+00   | no                                |  |
| Indeno(1,2,3-cd)pyrene | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |  |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

**Table 19**  
**Summary of Human Health Exposure Parameters**  
**Kerr McGee, Hattiesburg, MS**

| Receptors:   |                      | Adolescent<br>Visitor | Maintenance<br>Worker | Construction<br>Worker | Off-Site<br>Resident<br>Child | Off-Site<br>Resident<br>Adult |
|--|----------------------|-----------------------|-----------------------|------------------------|-------------------------------|-------------------------------|
| Parameter  | Units                |                       |                       |                        |                               |                               |
| Surface area available for exposure - soil           | cm <sup>2</sup> /day | 3052                  | 1                     | 3000                   | 1                             | 5560                          |
| Surface area available for exposure - sed. & sw      | cm <sup>2</sup> /day | 3945                  | 1                     | 3000                   | 1                             | 2229                          |
| Total skin surface area                              | cm <sup>2</sup>      | 12768.3               | 2                     | 20000                  | 2                             | 7213                          |
| Skin surface area available for exposure - soil      | %                    | 23.9%                 | 2                     | 15%                    | 2                             | 23.9%                         |
| Skin surface area available for exposure - sed. & sw | %                    | 30.9%                 | 2                     | 15.0%                  | 2                             | 30.9%                         |
| Adherence factor - soil                              | mg/cm <sup>2</sup>   | 0.026                 | 2                     | 0.038                  | 2                             | 0.026                         |
| Adherence factor - sed.                              | mg/cm <sup>3</sup>   | 0.33                  | 2                     | 0.038                  | 2                             | 0.33                          |
| Dermal absorption factor - benzo(a)pyrene            |                      | 0.03                  | 3                     | 0.03                   | 3                             | 0.03                          |
| Dermal absorption factor - other PAHs                |                      | 0.1                   | 3                     | 0.1                    | 3                             | 0.1                           |
| Exposure time  | hours/day            | 1                     | 5                     | 1                      | 5                             | 1                             |
| Exposure frequency - soils                           | days/year            | 12                    | 5                     | 150                    | 5                             | 80                            |
| Exposure frequency - sed. & sw                       | days/year            | 12                    | 5                     | 2                      | 5                             | 8                             |
| Exposure duration                                    | years                | 10                    | 6                     | 25                     | 6                             | 1                             |
| Body weight  | kg                   | 45                    | 6                     | 70                     | 6                             | 70                            |
| Averaging time - noncarcinogenic                     | days                 | 3650                  | 7                     | 9125                   | 7                             | 365                           |
| Averaging time - carcinogenic                        | days                 | 25550                 | 7                     | 25550                  | 7                             | 25550                         |
| Ingestion rate - soil                                | mg/day               | 100                   | 2                     | 100                    | 2                             | 480                           |
| Ingestion rate - surface water                       | L/hour               | 0.01                  | 6                     | 0.01                   | 6                             | 0.01                          |
| Matrix effect - PAHs                                 |                      | 1                     | 5                     | 1                      | 5                             | 1                             |
| Inhalation rate                                      | m <sup>3</sup> /day  | NA                    | NA                    | 20                     | 6                             | NA                            |
| Retention factor - semivolatiles                     |                      | NA                    | NA                    | 0.75                   | 8                             | NA                            |

NA - Not Applicable

1 Calculated

2 USEPA 1997, Exposure Factors Handbook

3 USEPA 1995, Region III Technical Guidance Manual: Assessing Dermal Exposure to Soil

4 USEPA 1992, Dermal Exposure Assessment

5 Reasonable Maximum

6 USEPA 1995, Region IV

7 USEPA 1991, HHEM Supplemental Guidances

8 International Commission on Radiological Protection, 1968



**Table 20**

**Particulate Emission Rate for Vehicular Movement and Excavation**

**Kerr McGee, Hattiesburg, MS**

**Vehicular Movement**

$$E = k * (5.9) * (s/12)(S/30) * (W/3)^{0.7}((w/4)^{0.5}) * ((365-p)/365) = \quad 16.49 \quad \text{lbs/vehicle mile}$$

|     |       |   |                    |
|-----|-------|---|--------------------|
| E = | 16.49 | particulate emission rate (lbs/vehicle mile - 30 miles travelled total over 80 - 8 hr days) |                    |
| k = | 0.5   | particle size multiplier  | US EPA AP-42, 1996 |
| s = | 50    | percent silt content  | Site Specific      |
| S = | 15    | mean vehicle speed (mi/hr)  | US EPA SEAM, 1988  |
| W = | 12.5  | mean vehicle weight (ton)   | US EPA SEAM, 1988  |
| w = | 8     | mean number of wheels per vehicle   | US EPA SEAM, 1988  |
| p = | 110   | mean number of days with ≥0.01 inches of precipitation per year                             | US EPA SEAM, 1988  |

Emission Rate      lbs/sec =  $(E \text{ lbs/mi}) * (30 \text{ mi/job}) * (\text{job}/80 \text{ days}) * (1 \text{ day}/8 \text{ hrs}) * (1 \text{ hr}/3600 \text{ sec})$

2.15E-04      lbs/sec

9.74E-02      g/sec

0.00010      kg/sec

**Excavation**

$$E = (1.0 * s^{1.5})M^{1.4} = \quad 7.90E+00 \quad \text{lbs/hour}$$

|     |          |                                      |               |
|-----|----------|--------------------------------------|---------------|
| E = | 7.90E+00 | particulate emission factor (lbs/hr) |               |
| s = | 50       | percent silt content                 | Site Specific |
| M = | 15.1     | percent soil moisture content        | Site Specific |

Emission Rate = 2.20E-03      lbs/sec

0.996      g/sec

0.000996      kg/sec



**Table 21**  
**Summary of Windrose Data**  
**Kerr McGee, Hattiesburg, MS**

**GRAPHICAL EXPOSURE MODELING SYSTEM**  
**STAR STATION JACKSON/THOMPSON, MS 1974-1978**

| DIRECTION | FREQUENCY | WINDSPEED | DIRECTION                          | FREQUENCY | WINDSPEED |
|-----------|-----------|-----------|------------------------------------|-----------|-----------|
| N         | 3.33325   | 0.03      | S                                  | 0.05336   | 3.08      |
| NNE       | 1.89301   | 0.03      | SSW                                | 0.09995   | 3.29      |
| NE        | 3.56791   | 0.07      | SW                                 | 0.10061   | 3.65      |
| ENE       | 0.12132   | 4.04      | WSW                                | 0.14723   | 3.93      |
| E         | 0.04843   | 3.39      | W                                  | 0.05047   | 3.7       |
| ESE       | 0.04328   | 3.12      | WNW                                | 0.04341   | 3.51      |
| SE        | 0.03686   | 3         | NW                                 | 0.02908   | 3.25      |
| SSE       | 0.05274   | 2.99      | NNW                                | 0.0406    | 3.26      |
| STABILITY | FREQUENCY | WINDSPEED | AUXILIARY VARIABLES                |           |           |
| 1         | 259.2     | 0.13      | Afternoon mixing height (meters)   | 1409      |           |
| 2         | 0.053     | 0.24      | Nocturnal mixing height (meters)   | 472       |           |
| 3         | 11.3      | 1         | Ambient air temperature (Kelvin)   | 303.6     |           |
| 4         | 0.01264   | 2.17      | Precipitation frequency (fraction) | 289.8     |           |
| 5         | 0.08137   | 2.98      | Precipitation intensity (mm/hour)  | 73.66     |           |
| 6         | 0.1315    | 3.91      | Grand average windspeed (m/s)      | 4.69      |           |



**Table 22**  
**Summary of Toxicity Values**  
**Kerr McGee, Huntington, MS**

| Chemical                    | mg/kg-day | Source | mg/kg-day | Source | mg/kg-day | Source    | mg/kg-day | Source   | mg/kg-day | Source | mg/kg-day | Source | Inhalation |
|-----------------------------|-----------|--------|-----------|--------|-----------|-----------|-----------|----------|-----------|--------|-----------|--------|------------|
|                             |           |        |           |        |           |           |           |          |           |        |           |        | CSF        |
|                             | RfD       | RfD    | RfD       | RfD    | RfD       | RfD       | RfD       | RfD      | RfD       | RfD    | CSF       | CSF    |            |
| Semivolatiles               |           |        |           |        |           |           |           |          |           |        |           |        | Source     |
| 2-Methylnaphthalene         | 2.00E-02  | E      | 5.70E-05  | H      | 0.5       | Region IV | 1.00E-02  |          |           |        |           |        |            |
| 2-Nitroaniline              |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| 2-Nitrophenol               |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| 3-Nitroaniline              |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| 4-Bromophenylphenylether    | 5.80E-02  | O      |           |        |           | 0.5       | Region IV | 2.90E-02 |           |        |           |        |            |
| 4-Chloro-3-methylphenol     |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| 4-Chlorophenylphenylether   |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| 4-Nitroaniline              |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Acenaphthylene              |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Benz(a)anthracene           |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Benz(a)pyrene               |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Benz(b)fluoranthene         |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Benz(g,h,i)perylene         |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Benz(k)fluoranthene         |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Bis(2-chloroethoxy)methane  |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Bis(2-chloroethyl)ether     |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Bis(2-ethylhexyl) phthalate | 2.00E-02  | IRIS   |           |        |           | 0.5       | Region IV | 2.00E-02 | W         |        |           |        |            |
| Carbazole                   |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Chrysene                    |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Dibenz(a,h)anthracene       |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Dibenzofuran                | 4.00E-03  | E      |           |        |           | 0.5       | Region IV | 2.00E-03 |           |        |           |        |            |
| Fluoranthene                | 4.00E-02  | IRIS   |           |        |           | 0.5       | Region IV | 2.00E-02 | 4.00E-01  | H      |           |        |            |
| Fluorene                    | 4.00E-02  | IRIS   |           |        |           | 0.5       | Region IV | 2.00E-02 | 4.00E-01  | H      |           |        |            |
| Hexachlorobenzene           | 8.00E-04  | IRIS   |           |        |           | 0.5       | Region IV | 4.00E-04 |           |        |           |        |            |
| Hexachlorocyclopentadiene   | 7.00E-03  | IRIS   | 2.00E-05  | H      | 0.5       | Region IV | 3.50E-03  |          |           |        |           |        |            |
| Indeno[1,2,3-cd]pyrene      |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| N-nitrosodi-n-propylamine   |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |
| Naphthalene                 | 2.00E-02  | IRIS   | 9.00E-04  | IRIS   | 0.5       | Region IV | 1.00E-02  |          |           |        |           |        |            |
| Phenanthrene                | 3.00E-02  | IRIS   |           |        | 0.5       | Region IV | 1.50E-02  | 3.00E-01 | H         |        |           |        |            |
| Pyrene                      |           |        |           |        |           | 0.5       | Region IV |          |           |        |           |        |            |

E - EPA/NCEA Regional Support provisional value from Region III RBC Tables, April 2000

H - Values are published in HEAST, 1997

IRIS - Values are available in IRIS, 2000

O - Values are withdrawn from other EPA documents as presented in the Region III RBC Tables, April 1995

Region IV - Region IV default value, 1995

W - Withdrawn from IRIS or HEAST

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**Table 23**  
**Summary of Hazard and Risk Calculations**  
**Kerr McGee, Hattiesburg, MS**

| Source/Pathway                          | Potentially Exposed Population | Total Hazard Index | Total Cancer Risk | Driving Constituent | Table Referenced |
|---|--------------------------------|--------------------|-------------------|---------------------|------------------|
| Dermal Exposure to Sediment in EU1      | Visitor                        | NA                 | 4E-08             |                     | 24               |
| Oral Exposure to Sediment in EU1        | Visitor                        | NA                 | 5E-08             |                     | 25               |
|   | Sub-Total                      | NA                 | 8E-08             |                     |                  |
| Dermal Exposure to Surface Water in EU1 | Visitor                        | NA                 | 4E-07             |                     | 26               |
| Oral Exposure to Surface Water in EU1   | Visitor                        | NA                 | 9E-09             |                     | 27               |
|   | Sub-Total                      | NA                 | 4E-07             |                     |                  |
| Dermal Exposure to Surface Soil in EU2  | Visitor                        | NA                 | 3E-08             |                     | 28               |
| Oral Exposure to Surface Soil in EU2    | Visitor                        | NA                 | 6E-07             |                     | 29               |
|   | Sub-Total                      | NA                 | 6E-07             |                     |                  |
| Dermal Exposure to Surface Soil in EU3  | Visitor                        | NA                 | 4E-09             |                     | 30               |
| Oral Exposure to Surface Soil in EU3    | Visitor                        | NA                 | 9E-08             |                     | 31               |
|   | Sub-Total                      | NA                 | 4E-09             |                     |                  |
| Dermal Exposure to Sediment in EU4      | Visitor                        | 7E-02              | 1E-05             | cPAHs               | 32               |
| Oral Exposure to Sediment in EU4        | Visitor                        | 3E-02              | 2E-05             | cPAHs               | 33               |
|   | Sub-Total                      | 1E-01              | 1E-05             |                     |                  |
| Dermal Exposure to Surface Water in EU4 | Visitor                        | 2E-04              | 9E-07             |                     | 34               |
| Oral Exposure to Surface Water in EU4   | Visitor                        | 2E-05              | 2E-08             |                     | 35               |
|   | Sub-Total                      | 3E-04              | 9E-07             |                     |                  |
| Dermal Exposure to Surface Soil in EU4  | Visitor                        | 4E-03              | 3E-06             | *                   | 36               |
| Oral Exposure to Surface Soil in EU4    | Visitor                        | 3E-02              | 3E-06             | cPAHs               | 37               |
|   | Sub-Total                      | 3E-02              | 5E-06             |                     |                  |
| Dermal Exposure to Surface Soil in EU5  | Visitor                        | NA                 | 3E-07             |                     | 38               |
| Oral Exposure to Surface Soil in EU5    | Visitor                        | NA                 | 6E-06             | Benzo(a)pyrene      | 39               |
|   | Sub-Total                      | NA                 | 3E-07             |                     |                  |
| <b>Visitor Total:</b>                   |                                | <b>1E-01</b>       | <b>2E-05</b>      |                     |                  |



**Table 23**  
**Summary of Hazard and Risk Calculations**  
**Kerr McGee, Hattiesburg, MS**

| Source/Pathway                          | Potentially Exposed Population | Total Hazard Index | Total Cancer Risk | Driving Constituent | Table Referenced |
|---|--------------------------------|--------------------|-------------------|---------------------|------------------|
| Dermal Exposure to Sediment in EU1      | Maintenance Worker             | NA                 | 8E-10             |                     | 40               |
| Oral Exposure to Sediment in EU1        | Maintenance Worker             | NA                 | 1E-08             |                     | 41               |
|   | Sub-Total                      | NA                 | 1E-08             |                     |                  |
| Dermal Exposure to Surface Water in EU1 | Maintenance Worker             | NA                 | 9E-08             |                     | 42               |
| Oral Exposure to Surface Water in EU1   | Maintenance Worker             | NA                 | 2E-09             |                     | 43               |
|   | Sub-Total                      | NA                 | 9E-08             |                     |                  |
| Dermal Exposure to Surface Soil in EU2  | Maintenance Worker             | NA                 | 5E-07             |                     | 44               |
| Oral Exposure to Surface Soil in EU2    | Maintenance Worker             | NA                 | 7E-06             | cPAHs               | 45               |
|   | Sub-Total                      | NA                 | 7E-06             |                     |                  |
| Dermal Exposure to Sediment in EU4      | Maintenance Worker             | 2E-05              | 2E-07             |                     | 46               |
| Oral Exposure to Sediment in EU4        | Maintenance Worker             | 3E-03              | 4E-06             | Benzo(a)pyrene      | 47               |
|   | Sub-Total                      | 3E-03              | 4E-06             |                     |                  |
| Dermal Exposure to Surface Water in EU4 | Maintenance Worker             | 2E-05              | 2E-07             |                     | 48               |
| Oral Exposure to Surface Water in EU4   | Maintenance Worker             | 2E-06              | 6E-09             |                     | 49               |
|   | Sub-Total                      | 2E-05              | 2E-07             |                     |                  |
| Dermal Exposure to Surface Soil in EU4  | Maintenance Worker             | 2E-02              | 8E-05             | Benzo(a)pyrene      | 50               |
| Oral Exposure to Surface Soil in EU4    | Maintenance Worker             | 1E-01              | 1E-03             | cPAHs               | 51               |
|   | Sub-Total                      | 1E-01              | 1E-03             |                     |                  |
| Dermal Exposure to Surface Soil in EU5  | Maintenance Worker             | NA                 | 6E-06             | Benzo(a)pyrene      | 52               |
| Oral Exposure to Surface Soil in EU5    | Maintenance Worker             | NA                 | 9E-05             | cPAHs               | 53               |
|   | Sub-Total                      | NA                 | 1E-04             |                     |                  |
| Maintenance Worker Total:               |                                | 1E-01              | 1E-03             |                     |                  |



**Table 23**  
**Summary of Hazard and Risk Calculations**  
**Kerr McGee, Hattiesburg, MS**

| Source/Pathway                          | Potentially Exposed Population | Total Hazard Index | Total Cancer Risk | Driving Constituent | Table Referenced |
|---|--------------------------------|--------------------|-------------------|---------------------|------------------|
| Dermal Exposure to Sediment in EU1      | Construction Worker            | NA                 | 5E-10             |                     | 54               |
| Oral Exposure to Sediment in EU1        | Construction Worker            | NA                 | 9E-09             |                     | 55               |
|   | Sub-Total                      | NA                 | 1E-08             |                     |                  |
| Dermal Exposure to Surface Water in EU1 | Construction Worker            | NA                 | 1E-08             |                     | 56               |
| Oral Exposure to Surface Water in EU1   | Construction Worker            | NA                 | 4E-10             |                     | 57               |
|   | Sub-Total                      | NA                 | 1E-08             |                     |                  |
| Dermal Exposure to Soil in EU2          | Construction Worker            | NA                 | 8E-08             |                     | 58               |
| Oral Exposure to Soil in EU2            | Construction Worker            | NA                 | 4E-07             |                     | 59               |
| Inhalation of Fugitive Dust in EU2      | Construction Worker            | NA                 | 1E-08             |                     | 60               |
|   | Sub-Total                      | NA                 | 5E-07             |                     |                  |
| Dermal Exposure to Sediment in EU4      | Construction Worker            | NA                 | 2E-07             |                     | 61               |
| Oral Exposure to Sediment in EU4        | Construction Worker            | NA                 | 3E-06             | Benzo(a)pyrene      | 62               |
|   | Sub-Total                      | NA                 | 3E-06             |                     |                  |
| Dermal Exposure to Surface Water in EU4 | Construction Worker            | 9E-07              | 3E-08             |                     | 63               |
| Oral Exposure to Surface Water in EU4   | Construction Worker            | 5E-07              | 3E-08             |                     | 64               |
|   | Sub-Total                      | 1E-06              | 3E-08             |                     |                  |
| Dermal Exposure to Soil in EU4          | Construction Worker            | NA                 | 8E-06             | Benzo(a)pyrene      | 65               |
| Oral Exposure to Soil in EU4            | Construction Worker            | NA                 | 4E-05             | cPAHs               | 66               |
| Inhalation of Fugitive Dust in EU4      | Construction Worker            | NA                 | 1E-06             | Benzo(a)pyrene      | 67               |
|   | Sub-Total                      | NA                 | 5E-05             |                     |                  |
| Dermal Exposure to Soil in EU5          | Construction Worker            | 2E-04              | 7E-07             |                     | 68               |
| Oral Exposure to Soil in EU5            | Construction Worker            | 3E-03              | 3E-06             | Benzo(a)pyrene      | 69               |
| Inhalation of Fugitive Dust in EU5      | Construction Worker            | NA                 | 1E-07             |                     | 70               |
|   | Sub-Total                      | 3E-03              | 4E-06             |                     |                  |
| <b>Construction Worker Total:</b>       |                                | <b>3E-03</b>       | <b>5E-05</b>      |                     |                  |

|   |                         |              |              |       |    |
|---|-------------------------|--------------|--------------|-------|----|
| Dermal Exposure to Sediment in EU6      | Child Off-Site Resident | NA           | 2E-05        | cPAHs | 71 |
| Oral Exposure to Sediment in EU6        | Child Off-Site Resident | NA           | 7E-05        | cPAHs | 72 |
|   | Sub-Total               | NA           | 9E-05        |       |    |
| Dermal Exposure to Sediment in EU6      | Adult Off-Site Resident | NA           | 4E-05        | cPAHs | 73 |
| Oral Exposure to Sediment in EU6        | Adult Off-Site Resident | 1E-04        | 3E-05        | cPAHs | 74 |
|   | Sub-Total               | 1E-04        | 7E-05        |       |    |
| Dermal Exposure to Surface Water in EU6 | Child Off-Site Resident | NA           | 2E-06        | *     | 75 |
| Oral Exposure to Surface Water in EU6   | Child Off-Site Resident | NA           | 5E-07        |       | 76 |
|   | Sub-Total               | NA           | 2E-06        |       |    |
| Dermal Exposure to Surface Water in EU6 | Adult Off-Site Resident | NA           | 5E-06        | *     | 77 |
| Oral Exposure to Surface Water in EU6   | Adult Off-Site Resident | NA           | 8E-08        |       | 78 |
|   | Sub-Total               | NA           | 5E-06        |       |    |
| <b>Off-Site Resident Total:</b>         |                         | <b>1E-04</b> | <b>2E-04</b> |       |    |

\*Estimated carcinogenic risk level is below *de minimis* level as no single constituent exceeded  $1 \times 10^{-6}$  and the cumulative site carcinogenic risk is below  $1 \times 10^{-4}$  (Section 501, MCEQ, 1999).



Table 24

Dermal Exposure to EU1 Sediment by an Adolescent Visitor (Aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |             |                        |
|---|------------------------------|-------------|------------------------|
|   | BW*AT                        |             |                        |
| Cs - Concentration in sediment =                            | mg/kg                        | chem. spec. |                        |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day         | 3945        | calculated             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>              | 12768.3     | USEPA 1997, EFH        |
| Fs - Fraction of skin surface area available for exposure = |                              | 30.9%       | USEPA 1997, EFH        |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>           | 0.33        | USEPA 1997, EFH        |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                              | 0.03        | USEPA 1995, Region III |
| EF - Exposure frequency =                                   | days/year                    | 12          | reasonable assumption  |
| ED - Exposure duration =                                    | years                        | 10          | USEPA 1995, Region IV  |
| CF - Conversion factor =                                    | kg/mg                        | 1.00E-06    |                        |
| BW - Body weight =  | kg                           | 45          | USEPA 1995, Region IV  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                         | 3650        | USEPA 1991, HHEM       |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                         | 25550       | USEPA 1991, HHEM       |

| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|---------------------------------|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| <b>Semivolatiles</b>   |                                 |                                |                              |              |   |                                   |             |
| Benzo(a)anthracene     | 5.90E-01                        | 1.68E-08                       | NA                           | NA           | 2.41E-09                                | 1.46E+00                          | 3.51E-09    |
| Benzo(a)pyrene         | 3.90E-01                        | 1.11E-08                       | NA                           | NA           | 1.59E-09                                | 1.46E+01                          | 2.32E-08    |
| Benzo(b)fluoranthene   | 5.80E-01                        | 1.66E-08                       | NA                           | NA           | 2.36E-09                                | 1.46E+00                          | 3.45E-09    |
| Benzo(k)fluoranthene   | 1.90E-01                        | 5.42E-09                       | NA                           | NA           | 7.75E-10                                | 1.46E-01                          | 1.13E-10    |
| Chrysene               | 5.30E-01                        | 1.51E-08                       | NA                           | NA           | 2.16E-09                                | 1.46E-02                          | 3.15E-11    |
| Dibenz(a,h)anthracene  | 6.20E-02                        | 1.77E-09                       | NA                           | NA           | 2.53E-10                                | 1.46E+01                          | 3.69E-09    |
| Indeno(1,2,3-cd)pyrene | 2.20E-01                        | 6.28E-09                       | NA                           | NA           | 8.97E-10                                | 1.46E+00                          | 1.31E-09    |

NA - Not Available

Total Cancer Risk = 3.53E-08



**Table 27****Oral Exposure to EU1 Surface Water by an Adolescent Visitor (aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | $\frac{C_{sw} \cdot IngR \cdot EF \cdot ED \cdot ET}{BW \cdot AT}$ |  |  |  |  |
|--|-----------|--|--|--|--|--|
| C <sub>sw</sub> - Concentration in surface water =   | mg/L      | see below  |  |  |  |  |
| IngR - Ingestion rate for surface water =            | L/hour    | 0.01   |  |  |  | USEPA 1995, Region IV                  |
| EF - Exposure frequency =                            | days/year | 12   |  |  |  | reasonable assumption                  |
| ED - Exposure duration =                             | years     | 10   |  |  |  | USEPA 1995, Region IV                  |
| ET - Exposure time =                                 | hrs/day   | 1  |  |  |  | USEPA 1992, Dermal Exposure Assessment |
| BW - Body weight =                                   | kg        | 45   |  |  |  | USEPA 1995, Region IV                  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650   |  |  |  | USEPA 1991, HHEM                       |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550  |  |  |  | USEPA 1991, HHEM                       |

| Constituent            | Concentration in Surface Water mg/L | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average                         | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-------------------------------------|--------------------------------|----------------------------|--------------|---------------------------------|--|-------------|
|                        |                                     |                                |                            |              | Lifetime Daily Intake mg/kg-day |  |             |
| <b>Semivolatiles</b>   |                                     |                                |                            |              |                                 |  |             |
| Benzo(a)anthracene     | 1.00E-03                            | 7.31E-09                       | NA                         | NA           | 1.04E-09                        | 7.30E-01                               | 7.62E-10    |
| Benzo(a)pyrene         | 5.00E-04                            | 3.65E-09                       | NA                         | NA           | 5.22E-10                        | 7.30E+00                               | 3.81E-09    |
| Benzo(b)fluoranthene   | 5.00E-04                            | 3.65E-09                       | NA                         | NA           | 5.22E-10                        | 7.30E-01                               | 3.81E-10    |
| Benzo(k)fluoranthene   | 5.00E-04                            | 3.65E-09                       | NA                         | NA           | 5.22E-10                        | NA                                     | NA          |
| Chrysene               | 5.00E-04                            | 3.65E-09                       | NA                         | NA           | 5.22E-10                        | 7.30E-02                               | 3.81E-11    |
| Dibenz(a,h)anthracene  | 5.00E-04                            | 3.65E-09                       | NA                         | NA           | 5.22E-10                        | 7.30E+00                               | 3.81E-09    |
| Indeno(1,2,3-cd)pyrene | 5.00E-04                            | 3.65E-09                       | NA                         | NA           | 5.22E-10                        | 7.30E-01                               | 3.81E-10    |

NA - Not Applicable

Total Cancer Risk = 9.18E-09



Table 28

Dermal Exposure to EU2 Surface Soil (0-1') by an Adolescent Visitor (aged 10-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                      | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |                        |                 |                     |                         |             |
|---|----------------------|------------------------------|------------------------|-----------------|---------------------|-------------------------|-------------|
|   |                      | BW*AT                        |                        |                 |                     |                         |             |
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.                  |                        |                 |                     |                         |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 3052                         | calculated             |                 |                     |                         |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 12768.3                      | USEPA 1997, EFH        |                 |                     |                         |             |
| Fs - Fraction of skin surface area available for exposure = |                      | 23.9%                        | USEPA 1997, EFH        |                 |                     |                         |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.026                        | USEPA 1997, EFH        |                 |                     |                         |             |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                      | 0.03                         | USEPA 1995, Region III |                 |                     |                         |             |
| EF - Exposure frequency =                                   | days/year            | 12                           | reasonable assumption  |                 |                     |                         |             |
| ED - Exposure duration =                                    | years                | 10                           | USEPA 1995, Region IV  |                 |                     |                         |             |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06                     |                        |                 |                     |                         |             |
| BW - Body weight =  | kg                   | 45                           | USEPA 1995, Region IV  |                 |                     |                         |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 3650                         | USEPA 1991, HHEM       |                 |                     |                         |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550                        | USEPA 1991, HHEM       |                 |                     |                         |             |
| Average   |                      |                              |                        |                 |                     |                         |             |
| Concentration in  |                      | Average                      | Dermal Chronic         | Lifetime Daily  | Cancer Slope        |                         |             |
| Constituent   | Soil<br>mg/kg        | Daily Intake<br>mg/kg-day    | RfD<br>mg/kg-day       | Hazard<br>Index | Intake<br>mg/kg-day | Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>  |                      |                              |                        |                 |                     |                         |             |
| Benzo(a)anthracene  | 6.70E+00             | 1.17E-08                     | NA                     | NA              | 1.66E-09            | 1.46E+00                | 2.43E-09    |
| Benzo(a)pyrene  | 5.08E+00             | 8.83416E-09                  | NA                     | NA              | 1.26E-09            | 1.46E+01                | 1.84E-08    |
| Benzo(b)fluoranthene  | 9.20E+00             | 1.60E-08                     | NA                     | NA              | 2.29E-09            | 1.46E+00                | 3.34E-09    |
| Benzo(k)fluoranthene  | 2.93E+00             | 5.10E-09                     | NA                     | NA              | 7.28E-10            | 1.46E-01                | 1.06E-10    |
| Chrysene  | 8.00E+00             | 1.39121E-08                  | NA                     | NA              | 1.99E-09            | 1.46E-02                | 2.90E-11    |
| Dibenz(a,h)anthracene                                       | 4.93E-01             | 8.57E-10                     | NA                     | NA              | 1.22E-10            | 1.46E+01                | 1.79E-09    |
| Indeno(1,2,3-cd)pyrene                                      | 3.70E+00             | 6.43E-09                     | NA                     | NA              | 9.19E-10            | 1.46E+00                | 1.34E-09    |

NA - Not Available

Total Cancer Risk = 2.75E-08



**Table 29**

**Oral Exposure to EU2 Surface Soil (0-1') by an Adolescent Visitor (aged 10-16 years)**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) = $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                             |                                |                            |              |   |  |             |
|--|-----------------------------|--------------------------------|----------------------------|--------------|---|--|-------------|
| Constituent  | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>   |                             |                                |                            |              |   |  |             |
| Benzo(a)anthracene   | 6.70E+00                    | 4.89E-07                       | NA                         | NA           | 6.99E-08                                | 7.30E-01                               | 5.10E-08    |
| Benzo(a)pyrene   | 5.08E+00                    | 3.71E-07                       | NA                         | NA           | 5.30E-08                                | 7.30E+00                               | 3.87E-07    |
| Benzo(b)fluoranthene   | 9.20E+00                    | 6.72E-07                       | NA                         | NA           | 9.60E-08                                | 7.30E-01                               | 7.01E-08    |
| Benzo(k)fluoranthene   | 2.93E+00                    | 2.14E-07                       | NA                         | NA           | 3.06E-08                                | 7.30E-02                               | 2.23E-09    |
| Chrysene   | 8.00E+00                    | 5.84E-07                       | NA                         | NA           | 8.35E-08                                | 7.30E-03                               | 6.10E-10    |
| Dibenz(a,h)anthracene  | 4.93E-01                    | 3.60E-08                       | NA                         | NA           | 5.15E-09                                | 7.30E+00                               | 3.76E-08    |
| Indeno(1,2,3-cd)pyrene   | 3.70E+00                    | 2.70E-07                       | NA                         | NA           | 3.86E-08                                | 7.30E-01                               | 2.82E-08    |

NA - Not Applicable

Total Cancer Risk = 5.77E-07



Table 30

*Dermal Exposure to EU3 Surface Soil (0-1') by an Adolescent Visitor (aged 10-16 years)*

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                                   | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                    |                 |  |   |             |
|---|-----------------------------------|---|------------------------------------|-----------------|--|---|-------------|
| Cs - Concentration in soil =                                | mg/kg                             | chem. spec.   |                                    |                 |  |   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day              | 3052  | calculated                         |                 |  |   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                   | 12768.3   | USEPA 1997, EFH                    |                 |  |   |             |
| Fs - Fraction of skin surface area available for exposure = |                                   | 23.9%   | USEPA 1997, EFH                    |                 |  |   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                | 0.026   | USEPA 1997, EFH                    |                 |  |   |             |
| ABS <sub>bap</sub> - Absorption - cPAHs =                   |                                   | 0.03  | USEPA 1995, Region III             |                 |  |   |             |
| EF - Exposure frequency =                                   | days/year                         | 12  | reasonable assumption              |                 |  |   |             |
| ED - Exposure duration =                                    | years                             | 10  | USEPA 1995, Region IV              |                 |  |   |             |
| CF - Conversion factor =                                    | kg/mg                             | 1.00E-06  |                                    |                 |  |   |             |
| BW - Body weight =  | kg                                | 45  | USEPA 1995, Region IV              |                 |  |   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                              | 3650  | USEPA 1991, HHEM                   |                 |  |   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                              | 25550   | USEPA 1991, HHEM                   |                 |  |   |             |
| Constituent   | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day                | Dermal<br>Chronic RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
|   |                                   |   |                                    |                 |  |   |             |
| <b>Semivolatiles</b>  |                                   |   |                                    |                 |  |   |             |
| Benzo(a)anthracene  | 5.40E-01                          | 9.39E-10  | NA                                 | NA              | 1.34E-10   | 1.46E+00                                | 1.96E-10    |
| Benzo(a)pyrene  | 7.10E-01                          | 1.23E-09  | NA                                 | NA              | 1.76E-10   | 1.46E+01                                | 2.58E-09    |
| Benzo(b)fluoranthene  | 1.40E+00                          | 2.43E-09  | NA                                 | NA              | 3.48E-10   | 1.46E+00                                | 5.08E-10    |
| Benzo(k)fluoranthene  | 4.90E-01                          | 8.52E-10  | NA                                 | NA              | 1.22E-10   | 1.46E-01                                | 1.78E-11    |
| Chrysene  | 8.70E-01                          | 1.51E-09  | NA                                 | NA              | 2.16E-10   | 1.46E-02                                | 3.16E-12    |
| Dibenz(a,h)anthracene                                       | 1.60E-01                          | 2.78E-10  | NA                                 | NA              | 3.97E-11   | 1.46E+01                                | 5.80E-10    |
| Indeno(1,2,3-cd)pyrene                                      | 6.00E-01                          | 1.04E-09  | NA                                 | NA              | 1.49E-10   | 1.46E+00                                | 2.18E-10    |

NA - Not Available

Total Cancer Risk = 4.10E-09



**Table 31****Oral Exposure to EU3 Surface Soil by an Adolescent Visitor (aged 10-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | $\frac{\text{Cd} * \text{Ingr} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                       |  |  |
|--|-----------|---|-----------------------|--|--|
| Cd - Concentration in sediment =                     | mg/kg     | see below   |                       |  |  |
| Ingr - Ingestion rate for soil =                     | mg/day    | 100   | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 12  | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 10  | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06  |                       |  |  |
| ME <sub>s</sub> - Matrix effect - PAHs =             |           | 1   | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 45  | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650  | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550   | USEPA 1991, HHEM      |  |  |

| Constituent            | Concentration in<br>Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily<br>Intake | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------------|--------------------------------------|----------------------------------|-----------------|-------------------------------------|--|-------------|
|                        |                                   |                                      |                                  |                 | mg/kg-day                           |  |             |
| <b>Semivolatiles</b>   |                                   |                                      |                                  |                 |                                     |  |             |
| Benzo(a)anthracene     | 5.40E-01                          | 3.95E-08                             | NA                               | NA              | 5.64E-09                            | 7.30E-01                                     | 4.11E-09    |
| Benzo(a)pyrene         | 7.10E-01                          | 5.19E-08                             | NA                               | NA              | 7.41E-09                            | 7.30E+00                                     | 5.41E-08    |
| Benzo(b)fluoranthene   | 1.40E+00                          | 1.02E-07                             | NA                               | NA              | 1.46E-08                            | 7.30E-01                                     | 1.07E-08    |
| Benzo(k)fluoranthene   | 4.90E-01                          | 3.58E-08                             | NA                               | NA              | 5.11E-09                            | 7.30E-02                                     | 3.73E-10    |
| Chrysene               | 8.70E-01                          | 6.36E-08                             | NA                               | NA              | 9.08E-09                            | 7.30E-03                                     | 6.63E-11    |
| Dibenz(a,h)anthracene  | 1.60E-01                          | 1.17E-08                             | NA                               | NA              | 1.67E-09                            | 7.30E+00                                     | 1.22E-08    |
| Indeno(1,2,3-cd)pyrene | 6.00E-01                          | 4.38E-08                             | NA                               | NA              | 6.26E-09                            | 7.30E-01                                     | 4.57E-09    |

NA - Not Applicable

Total Cancer Risk = 8.61E-08



**Table 32****Dermal Exposure to EU4 Sediment by an Adolescent Visitor (Aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u><math>\frac{\text{Cs} * \text{SA} * \text{AH} * \text{ABS} * \text{EF} * \text{ED} * \text{CF}}{\text{BW} * \text{AT}}</math></u> |                                |                              |              |   |                                   |             |
|---|--|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| Cs - Concentration in sediment =                            | mg/kg  | chem. spec.                    |                              |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day   | 3945                           | calculated                   |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>  | 12768.3                        | USEPA 1997, EFH              |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |  | 30.9%                          | USEPA 1997, EFH              |              |   |                                   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.33                           | USEPA 1997, EFH              |              |   |                                   |             |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |  | 0.03                           | USEPA 1995, Region III       |              |   |                                   |             |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |  | 0.1                            | USEPA 1995, Region III       |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year  | 12                             | reasonable assumption        |              |   |                                   |             |
| ED - Exposure duration =                                    | years  | 10                             | USEPA 1995, Region IV        |              |   |                                   |             |
| CF - Conversion factor =                                    | kg/mg  | 1.00E-06                       |                              |              |   |                                   |             |
| BW - Body weight =  | kg   | 45                             | USEPA 1995, Region IV        |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days   | 3650                           | USEPA 1991, HHM              |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days   | 25550                          | USEPA 1991, HHM              |              |   |                                   |             |
| Constituent   | Concentration in Sediment mg/kg  | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |  |                                |                              |              |   |                                   |             |
| Benzo(a)anthracene  | 3.30E+02   | 9.42E-06                       | NA                           | NA           | 1.35E-06                                | 1.46E+00                          | 1.96E-06    |
| Benzo(a)pyrene  | 1.30E+02   | 3.71E-06                       | NA                           | NA           | 5.30E-07                                | 1.46E+01                          | 7.74E-06    |
| Benzo(b)fluoranthene  | 1.80E+02   | 5.14E-06                       | NA                           | NA           | 7.34E-07                                | 1.46E+00                          | 1.07E-06    |
| Benzo(k)fluoranthene  | 6.40E+01   | 1.83E-06                       | NA                           | NA           | 2.61E-07                                | 1.46E-01                          | 3.81E-08    |
| Carbazole   | 5.90E+02   | 5.61E-05                       | NA                           | NA           | 8.02E-06                                | 2.00E-02                          | 1.60E-07    |
| Chrysene  | 2.90E+02   | 8.28E-06                       | NA                           | NA           | 1.18E-06                                | 1.46E-02                          | 1.73E-08    |
| Dibenz(a,h)anthracene                                       | 1.20E+01   | 3.42E-07                       | NA                           | NA           | 4.89E-08                                | 1.46E+01                          | 7.14E-07    |
| Dibenzofuran  | 9.40E+02   | 8.94E-05                       | 2.00E-03                     | 4.47E-02     | 1.28E-05                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene                                      | 4.70E+01   | 1.34E-06                       | NA                           | NA           | 1.92E-07                                | 1.46E+00                          | 2.80E-07    |
| Naphthalene   | 3.00E+03   | 2.85E-04                       | 1.00E-02                     | 2.85E-02     | 4.08E-05                                | NA                                | NA          |
| Phenanthrene  | 3.20E+03   | 3.04E-04                       | NA                           | NA           | 4.35E-05                                | NA                                | NA          |

NA - Not Available

Total Hazard Index = 7.32E-02

Total Cancer Risk = 1.20E-05



**Table 33****Oral Exposure to EU4 Sediment by an Adolescent Visitor (Aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u><math>\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}</math></u> |                       |  |  |
|--|-----------|---|-----------------------|--|--|
| Cd - Concentration in sediment =                     | mg/kg     | see below   |                       |  |  |
| IngR - Ingestion rate for sediment =                 | mg/day    | 100   | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 12  | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 10  | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06  |                       |  |  |
| ME - Matrix effect =                                 |           | 1   | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 45  | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650  | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550   | USEPA 1991, HHEM      |  |  |

| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|---------------------------------|--------------------------------|----------------------------|--------------|---|--|-------------|
|                        |                                 |                                |                            |              | mg/kg-day                               | 1/(mg/kg-day)                          |             |
| <b>Semivolatiles</b>   |                                 |                                |                            |              |   |  |             |
| Benzo(a)anthracene     | 3.30E+02                        | 2.41E-05                       | NA                         | NA           | 3.44E-06                                | 7.30E-01                               | 2.51E-06    |
| Benzo(a)pyrene         | 1.30E+02                        | 9.50E-06                       | NA                         | NA           | 1.36E-06                                | 7.30E+00                               | 9.90E-06    |
| Benzo(b)fluoranthene   | 1.80E+02                        | 1.32E-05                       | NA                         | NA           | 1.88E-06                                | 7.30E-01                               | 1.37E-06    |
| Benzo(k)fluoranthene   | 6.40E+01                        | 4.68E-06                       | NA                         | NA           | 6.68E-07                                | 7.30E-02                               | 4.88E-08    |
| Carbazole              | 5.90E+02                        | 4.31E-05                       | NA                         | NA           | 6.16E-06                                | 2.00E-02                               | 1.23E-07    |
| Chrysene               | 2.90E+02                        | 2.12E-05                       | NA                         | NA           | 3.03E-06                                | 7.30E-03                               | 2.21E-08    |
| Dibenz(a,h)anthracene  | 1.20E+01                        | 8.77E-07                       | NA                         | NA           | 1.25E-07                                | 7.30E+00                               | 9.14E-07    |
| Dibenzofuran           | 9.40E+02                        | 6.87E-05                       | 4.00E-03                   | 1.72E-02     | 9.81E-06                                | NA                                     | NA          |
| Indeno(1,2,3-cd)pyrene | 4.70E+01                        | 3.43E-06                       | NA                         | NA           | 4.91E-07                                | 7.30E-01                               | 3.58E-07    |
| Naphthalene            | 3.00E+03                        | 2.19E-04                       | 2.00E-02                   | 1.10E-02     | 3.13E-05                                | NA                                     | NA          |
| Phenanthrene           | 3.20E+03                        | 2.34E-04                       | NA                         | NA           | 3.34E-05                                | NA                                     | NA          |

NA - Not Applicable

Total Hazard Index = 2.81E-02

Total Cancer Risk = 1.53E-05



**Table 34**

**Dermal Exposure to EU4 Surface Water by an Adolescent Visitor (aged 7-16 years)**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =   | $C_w * SA * K_p * ABS * ET * EF * ED * CF$ |           |  |  |  |
|--|--|-----------|--|--|--|
|  | BW * AT                                    |           |  |  |  |
| $C_w$ - Concentration in surface water =                       | mg/L                                       | see below |  |  |  |
| SA - Surface area available for exposure =                     | cm <sup>2</sup>                            | 3945      | calculated                             |  |  |
| $SA_t$ - Total skin surface area =                             | cm <sup>2</sup>                            | 12768.3   | USEPA 1997, EFH                        |  |  |
| $F_s$ - Fraction of skin surface area available for exposure = |  | 30.9%     | USEPA 1997, EFH                        |  |  |
| $K_p$ - Dermal permeability constant =                         | cm/hr                                      | see below |  |  |  |
| $ABS_p$ - Absorption - cPAHs =                                 |  | 0.03      | USEPA 1995, Region III                 |  |  |
| $ABS_o$ - Absorption - other SVOCs =                           |  | 0.1       | USEPA 1995, Region III                 |  |  |
| ET - Exposure time =   | hrs/day                                    | 1         | USEPA 1992, Dermal Exposure Assessment |  |  |
| EF - Exposure frequency =                                      | days/year                                  | 12        | reasonable assumption                  |  |  |
| ED - Exposure duration =                                       | years                                      | 10        | USEPS 1995, Region IV                  |  |  |
| CF - Conversion factor =                                       | L/cm <sup>3</sup>                          | 1.00E-03  |  |  |  |
| BW - Body weight =   | kg   | 45        | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =           | days                                       | 3650      | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =              | days                                       | 25550     | USEPA 1991, HHEM                       |  |  |

| Constituent                | Concentration in Surface Water |          | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day |          | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|----------------------------|--------------------------------|----------|--------------------------------|------------------------------|----------|--------------|---|-----------------------------------|-------------|
|                            | Water mg/L                     | Kp cm/hr |                                | Chronic                      | RfD      |              |   |                                   |             |
| <b>Semivolatiles</b>       |                                |          |                                |                              |          |              |   |                                   |             |
| Benzo(a)anthracene         | 5.00E-03                       | 8.10E-01 | 3.50E-07                       | NA                           | NA       | NA           | 5.00E-08                                | 1.46E+00                          | 7.30E-08    |
| Benzo(a)pyrene             | 5.00E-04                       | 1.20E+00 | 5.19E-08                       | NA                           | NA       | NA           | 7.41E-09                                | 1.46E+01                          | 1.08E-07    |
| Benzo(b)fluoranthene       | 1.20E-02                       | 1.20E+00 | 1.25E-06                       | NA                           | NA       | NA           | 1.78E-07                                | 1.46E+00                          | 2.60E-07    |
| Benzo(k)fluoranthene       | 2.00E-03                       | 4.48E+01 | 7.74E-06                       | NA                           | NA       | NA           | 1.11E-06                                | 1.46E-01                          | 1.62E-07    |
| Bis(2-ethylhexyl)phthalate | 3.00E-03                       | 3.30E-02 | 2.85E-08                       | 1.00E-02                     | 2.85E-06 | 2.85E-06     | 4.08E-09                                | 1.40E-02                          | 5.71E-11    |
| Carbazole                  | 1.00E-02                       | 3.57E-02 | 1.03E-07                       | NA                           | NA       | NA           | 1.47E-08                                | 2.00E-02                          | 2.94E-10    |
| Chrysene                   | 6.00E-03                       | 8.10E-01 | 4.20E-07                       | NA                           | NA       | NA           | 6.00E-08                                | 1.46E-02                          | 8.77E-10    |
| Dibenz(a,h)anthracene      | 5.00E-04                       | 2.70E+00 | 1.17E-07                       | NA                           | NA       | NA           | 1.67E-08                                | 1.46E+01                          | 2.43E-07    |
| Dibenzofuran               | 1.10E-02                       | 1.51E-01 | 4.79E-07                       | 2.00E-03                     | 2.40E-04 | 2.40E-04     | 6.84E-08                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene     | 5.00E-04                       | 1.90E+00 | 8.22E-08                       | NA                           | NA       | NA           | 1.17E-08                                | 1.46E+00                          | 1.71E-08    |
| Phenanthrene               | 1.70E-02                       | 2.30E-01 | 1.13E-06                       | NA                           | NA       | NA           | 1.61E-07                                | NA                                | NA          |

NA - Not Available

Total Hazard Index = 2.42E-04

Total Cancer Risk = 8.64E-07



**Table 35****Oral Exposure to EU4 Surface Water by an Adolescent Visitor (aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u><math>C_{sw} \times IngR \times EF \times ED \times ET</math></u> |  |  |  |
|--|-----------|--|--|--|--|
|  |           | <u>BW * AT</u>   |  |  |  |
| Csw - Concentration in surface water =               | mg/L      | see below  |  |  |  |
| IngR - Ingestion rate for surface water =            | L/hour    | 0.01   | USEPA 1995, Region IV                  |  |  |
| EF - Exposure frequency =                            | days/year | 12   | reasonable assumption                  |  |  |
| ED - Exposure duration =                             | years     | 10   | USEPA 1995, Region IV                  |  |  |
| ET - Exposure time =                                 | hrs/day   | 1  | USEPA 1992, Dermal Exposure Assessment |  |  |
| BW - Body weight =                                   | kg        | 45   | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650   | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550  | USEPA 1991, HHEM                       |  |  |

| Constituent                | Concentration<br>in Surface<br>Water<br>mg/L | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average                               |          | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
|----------------------------|--|--------------------------------------|----------------------------------|-----------------|---------------------------------------|----------|--|-------------|
|                            |  |                                      |                                  |                 | Lifetime Daily<br>Intake<br>mg/kg-day | Daily    |  |             |
| <b>Semivolatiles</b>       |  |                                      |                                  |                 |                                       |          |  |             |
| Benzo(a)anthracene         | 5.00E-03                                     | 3.65E-08                             | NA                               | NA              | 5.22E-09                              | 7.30E-01 | 3.81E-09                                     |             |
| Benzo(a)pyrene             | 5.00E-04                                     | 3.65E-09                             | NA                               | NA              | 5.22E-10                              | 7.30E+00 | 3.81E-09                                     |             |
| Benzo(b)fluoranthene       | 1.20E-02                                     | 8.77E-08                             | NA                               | NA              | 1.25E-08                              | 7.30E-01 | 9.14E-09                                     |             |
| Benzo(k)fluoranthene       | 2.00E-03                                     | 1.46E-08                             | NA                               | NA              | 2.09E-09                              | 7.30E-02 | 1.52E-10                                     |             |
| Bis(2-ethylhexyl)phthalate | 3.00E-03                                     | 2.19E-08                             | 2.00E-02                         | 1.10E-06        | 3.13E-09                              | 1.40E-02 | 4.38E-11                                     |             |
| Carbazole                  | 1.00E-02                                     | 7.31E-08                             | NA                               | NA              | 1.04E-08                              | 2.00E-02 | 2.09E-10                                     |             |
| Chrysene                   | 6.00E-03                                     | 4.38E-08                             | NA                               | NA              | 6.26E-09                              | 7.30E-03 | 4.57E-11                                     |             |
| Dibenz(a,h)anthracene      | 5.00E-04                                     | 3.65E-09                             | NA                               | NA              | 5.22E-10                              | 7.30E+00 | 3.81E-09                                     |             |
| Dibenzofuran               | 1.10E-02                                     | 8.04E-08                             | 4.00E-03                         | 2.01E-05        | 1.15E-08                              | NA       | NA   |             |
| Indeno(1,2,3-cd)pyrene     | 5.00E-04                                     | 3.65E-09                             | NA                               | NA              | 5.22E-10                              | 7.30E-01 | 3.81E-10                                     |             |
| Phenanthrene               | 1.70E-02                                     | 1.24E-07                             | NA                               | NA              | 1.77E-08                              | NA       | NA   |             |

NA - Not Applicable

Total Hazard Index = 2.12E-05

Total Cancer Risk = 2.14E-08



Table 36

Dermal Exposure to EU4 Surface Soil (0-1') by an Adolescent Visitor (Aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |             |                        |
|---|------------------------------|-------------|------------------------|
|   | BW*AT                        |             |                        |
| Cs - Concentration in soil =                                | mg/kg                        | chem. spec. |                        |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day         | 3052        | calculated             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>              | 12768.3     | USEPA 1997, EFH        |
| Fs - Fraction of skin surface area available for exposure = |                              | 23.9%       | USEPA 1997, EFH        |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>           | 0.026       | USEPA 1997, EFH        |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                              | 0.03        | USEPA 1995, Region III |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |                              | 0.1         | USEPA 1995, Region III |
| EF - Exposure frequency =                                   | days/year                    | 12          | reasonable assumption  |
| ED - Exposure duration =                                    | years                        | 10          | USEPA 1995, Region IV  |
| CF - Conversion factor =                                    | kg/mg                        | 1.00E-06    |                        |
| BW - Body weight =  | kg                           | 45          | USEPA 1995, Region IV  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                         | 3650        | USEPA 1991, HHEM       |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                         | 25550       | USEPA 1991, HHEM       |

| Constituent            | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day | RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |          |          | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------|--------------------------------|---------------|--------------|---|----------|----------|-----------------------------------|-------------|
|                        |                             |                                |               |              | Lifetime                                | Daily    | Chronic  |                                   |             |
| <b>Semivolatiles</b>   |                             |                                |               |              |   |          |          |                                   |             |
| Benzo(a)anthracene     | 9.30E+02                    | 1.62E-06                       | NA            | NA           | 2.31E-07                                | 1.46E+00 | 3.37E-07 |                                   |             |
| Benzo(a)pyrene         | 5.00E+02                    | 8.70E-07                       | NA            | NA           | 1.24E-07                                | 1.46E+01 | 1.81E-06 |                                   |             |
| Benzo(b)fluoranthene   | 5.30E+02                    | 9.22E-07                       | NA            | NA           | 1.32E-07                                | 1.46E+00 | 1.92E-07 |                                   |             |
| Benzo(k)fluoranthene   | 2.90E+02                    | 5.04E-07                       | NA            | NA           | 7.20E-08                                | 1.46E-01 | 1.05E-08 |                                   |             |
| Carbazole              | 2.30E+02                    | 1.33E-06                       | NA            | NA           | 1.90E-07                                | 2.00E-02 | 3.81E-09 |                                   |             |
| Chrysene               | 6.90E+02                    | 1.20E-06                       | NA            | NA           | 1.71E-07                                | 1.46E-02 | 2.50E-09 |                                   |             |
| Dibenz(a,h)anthracene  | 6.40E+01                    | 1.11E-07                       | NA            | NA           | 1.59E-08                                | 1.46E+01 | 2.32E-07 |                                   |             |
| Fluoranthene           | 4.60E+03                    | 2.67E-05                       | 2.00E-02      | 1.33E-03     | 3.81E-06                                | NA       | NA       |                                   |             |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                    | 4.35E-07                       | NA            | NA           | 6.21E-08                                | 1.46E+00 | 9.07E-08 |                                   |             |
| Naphthalene            | 2.20E+03                    | 1.28E-05                       | 1.00E-02      | 1.28E-03     | 1.82E-06                                | NA       | NA       |                                   |             |
| Phenanthrene           | 6.40E+03                    | 3.71E-05                       | NA            | NA           | 5.30E-06                                | NA       | NA       |                                   |             |
| Pyrene                 | 4.40E+03                    | 2.55E-05                       | 1.50E-02      | 1.70E-03     | 3.64E-06                                | NA       | NA       |                                   |             |

NA - Not Available

Total Hazard Index = 4.31E-03

Total Cancer Risk = 2.68E-06



**Table 37****Oral Exposure to EU4 Surface Soil (0-1') by an Adolescent Visitor (Aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u><math>\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}</math></u> |                       |  |  |
|--|-----------|---|-----------------------|--|--|
| Cd - Concentration in soil =                         | mg/kg     | see below   |                       |  |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 100   | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 12  | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 10  | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06  |                       |  |  |
| ME - Matrix effect =                                 |           | 1   | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 45  | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650  | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550   | USEPA 1991, HHEM      |  |  |

| Constituent            | Concentration in<br>Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average                               |  |             |
|------------------------|-----------------------------------|--------------------------------------|----------------------------------|-----------------|---------------------------------------|--|-------------|
|                        |                                   |                                      |                                  |                 | Lifetime Daily<br>Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>   |                                   |                                      |                                  |                 |                                       |  |             |
| Benzo(a)anthracene     | 9.30E+02                          | 6.79E-05                             | NA                               | NA              | 9.71E-06                              | 7.30E-01                                     | 7.09E-06    |
| Benzo(a)pyrene         | 5.00E+02                          | 3.65E-05                             | NA                               | NA              | 5.22E-06                              | 7.30E+00                                     | 3.81E-05    |
| Benzo(b)fluoranthene   | 5.30E+02                          | 3.87E-05                             | NA                               | NA              | 5.53E-06                              | 7.30E-01                                     | 4.04E-06    |
| Benzo(k)fluoranthene   | 2.90E+02                          | 2.12E-05                             | NA                               | NA              | 3.03E-06                              | 7.30E-02                                     | 2.21E-07    |
| Carbazole              | 2.30E+02                          | 1.68E-05                             | NA                               | NA              | 2.40E-06                              | 2.00E-02                                     | 4.80E-08    |
| Chrysene               | 6.90E+02                          | 5.04E-05                             | NA                               | NA              | 7.20E-06                              | 7.30E-03                                     | 5.26E-08    |
| Dibenz(a,h)anthracene  | 6.40E+01                          | 4.68E-06                             | NA                               | NA              | 6.68E-07                              | 7.30E+00                                     | 4.88E-06    |
| Fluoranthene           | 4.60E+03                          | 3.36E-04                             | 4.00E-02                         | 8.40E-03        | 4.80E-05                              | NA   | NA          |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                          | 1.83E-05                             | NA                               | NA              | 2.61E-06                              | 7.30E-01                                     | 1.90E-06    |
| Naphthalene            | 2.20E+03                          | 1.61E-04                             | 2.00E-02                         | 8.04E-03        | 2.30E-05                              | NA   | NA          |
| Phenanthrene           | 3.20E+03                          | 2.34E-04                             | NA                               | NA              | 3.34E-05                              | NA   | NA          |
| Pyrene                 | 4.40E+03                          | 3.21E-04                             | 3.00E-02                         | 1.07E-02        | 4.59E-05                              | NA   | NA          |

NA - Not Applicable

Total Hazard Index = 2.72E-02

Total Cancer Risk = 5.63E-05



Table 38

Dermal Exposure to EU5 Surface Soil (0-1') by an Adolescent Visitor (Aged 7-16 years)  
Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                          |                        |  |                                      |             |          |
|---|---|--------------------------|------------------------|--|--------------------------------------|-------------|----------|
| Cs - Concentration in soil =                                | mg/kg   | chem. spec.              |                        |  |                                      |             |          |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                                | 3052                     | calculated             |  |                                      |             |          |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                                     | 12768.3                  | USEPA 1997, EFH        |  |                                      |             |          |
| Fs - Fraction of skin surface area available for exposure = |   | 23.9%                    | USEPA 1997, EFH        |  |                                      |             |          |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                                  | 0.026                    | USEPA 1997, EFH        |  |                                      |             |          |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |   | 0.03                     | USEPA 1995, Region III |  |                                      |             |          |
| EF - Exposure frequency =                                   | days/year   | 12                       | reasonable assumption  |  |                                      |             |          |
| ED - Exposure duration =                                    | years   | 10                       | USEPA 1995, Region IV  |  |                                      |             |          |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                 |                        |  |                                      |             |          |
| BW - Body weight =  | kg  | 45                       | USEPA 1995, Region IV  |  |                                      |             |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 3650                     | USEPA 1991, HHEM       |  |                                      |             |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                    | USEPA 1991, HHEM       |  |                                      |             |          |
| Dermal  |   |                          |                        |  |                                      |             |          |
| Constituent   | Average Daily Intake<br>Soil mg/kg                  | Chronic RfD<br>mg/kg-day | Hazard Index           | Average Lifetime Daily Intake<br>mg/kg-day | Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |          |
| Semivolatiles   |   |                          |                        |  |                                      |             |          |
| Benzo(a)anthracene  | 8.35E+01  | 1.45E-07                 | NA                     | NA   | 2.07E-08                             | 1.46E+00    | 3.03E-08 |
| Benzo(a)pyrene  | 5.25E+01  | 9.13E-08                 | NA                     | NA   | 1.30E-08                             | 1.46E+01    | 1.90E-07 |
| Benzo(b)fluoranthene  | 7.95E+01  | 1.38E-07                 | NA                     | NA   | 1.98E-08                             | 1.46E+00    | 2.88E-08 |
| Benzo(k)fluoranthene  | 2.85E+01  | 4.96E-08                 | NA                     | NA   | 7.08E-09                             | 1.46E-01    | 1.03E-09 |
| Chrysene  | 8.25E+01  | 1.43E-07                 | NA                     | NA   | 2.05E-08                             | 1.46E-02    | 2.99E-10 |
| Dibenz(a,h)anthracene                                       | 7.45E+00  | 1.30E-08                 | NA                     | NA   | 1.85E-09                             | 1.46E+01    | 2.70E-08 |
| Indeno(1,2,3-cd)pyrene                                      | 3.10E+01  | 5.39E-08                 | NA                     | NA   | 7.70E-09                             | 1.46E+00    | 1.12E-08 |

NA - Not Available

Total Cancer Risk = 2.89E-07



**Table 39****Oral Exposure to EU5 Surface Soil (0-1') by an Adolescent Visitor (Aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u><math>Cd * IngR * EF * ED * CF * ME</math></u> |                       |  |  |
|--|-----------|---|-----------------------|--|--|
|  |           | <u>BW * AT</u>                                    |                       |  |  |
| Cd - Concentration in sediment =                     | mg/kg     | see below   |                       |  |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 100   | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 12  | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 10  | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06  |                       |  |  |
| ME - Matrix effect =                                 |           | 1   | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 45  | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650  | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550   | USEPA 1991, HHEM      |  |  |

| Constituent            | Concentration in<br>Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day |          |          | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------------|--------------------------------------|----------------------------------|-----------------|--|----------|----------|--|-------------|
|                        |                                   |                                      |                                  |                 |  |          |          |  |             |
| <b>Semivolatiles</b>   |                                   |                                      |                                  |                 |  |          |          |  |             |
| Benzo(a)anthracene     | 8.35E+01                          | 6.10E-06                             | NA                               | NA              | 8.71E-07   | 7.30E-01 | 6.36E-07 |  |             |
| Benzo(a)pyrene         | 5.25E+01                          | 3.84E-06                             | NA                               | NA              | 5.48E-07   | 7.30E+00 | 4.00E-06 |  |             |
| Benzo(b)fluoranthene   | 7.95E+01                          | 5.81E-06                             | NA                               | NA              | 8.30E-07   | 7.30E-01 | 6.06E-07 |  |             |
| Benzo(k)fluoranthene   | 2.85E+01                          | 2.08E-06                             | NA                               | NA              | 2.97E-07   | 7.30E-02 | 2.17E-08 |  |             |
| Chrysene               | 8.25E+01                          | 6.03E-06                             | NA                               | NA              | 8.61E-07   | 7.30E-03 | 6.29E-09 |  |             |
| Dibenz(a,h)anthracene  | 7.45E+00                          | 5.44E-07                             | NA                               | NA              | 7.78E-08   | 7.30E+00 | 5.68E-07 |  |             |
| Indeno(1,2,3-cd)pyrene | 3.10E+01                          | 2.26E-06                             | NA                               | NA              | 3.24E-07   | 7.30E-01 | 2.36E-07 |  |             |

NA - Not Applicable

Total Cancer Risk = 6.07E-06



**Table 40****Dermal Exposure to EU1 Sediment by a Maintenance Worker**

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                      | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                        |  |
|---|----------------------|---|------------------------|--|
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.   |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 3000  | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000   | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 15.0%   | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.038   | USEPA 1997, EFH        |  |
| ABSp - Absorption - cPAHs =                                 |                      | 0.03  | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year            | 2   | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                | 25  | USEPA 1995, Region IV  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06  |                        |  |
| BW - Body weight =  | kg                   | 70  | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 9125  | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550   | USEPA 1991, HHEM       |  |

| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|---------------------------------|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| <b>Semivolatiles</b>   |                                 |                                |                              |              |   |                                   |             |
| Benzo(a)anthracene     | 5.90E-01                        | 1.58E-10                       | NA                           | NA           | 5.64E-11                                | 1.46E+00                          | 8.24E-11    |
| Benzo(a)pyrene         | 3.90E-01                        | 1.04E-10                       | NA                           | NA           | 3.73E-11                                | 1.46E+01                          | 5.44E-10    |
| Benzo(b)fluoranthene   | 5.80E-01                        | 1.55E-10                       | NA                           | NA           | 5.55E-11                                | 1.46E+00                          | 8.10E-11    |
| Benzo(k)fluoranthene   | 1.90E-01                        | 5.09E-11                       | NA                           | NA           | 1.82E-11                                | 1.46E-01                          | 2.65E-12    |
| Chrysene               | 5.30E-01                        | 1.42E-10                       | NA                           | NA           | 5.07E-11                                | 1.46E-02                          | 7.40E-13    |
| Dibenz(a,h)anthracene  | 6.20E-02                        | 1.66E-11                       | NA                           | NA           | 5.93E-12                                | 1.46E+01                          | 8.65E-11    |
| Indeno(1,2,3-cd)pyrene | 2.20E-01                        | 5.89E-11                       | NA                           | NA           | 2.10E-11                                | 1.46E+00                          | 3.07E-11    |

NA - Not Available

Total Cancer Risk = 8.28E-10



**Table 41**  
**Oral Exposure to EU1 Sediment by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u>Cd*IngR*EF*ED*CF*ME</u><br>BW*AT |                       |  |  |
|--|-----------|-------------------------------------|-----------------------|--|--|
| Cd - Concentration in sediment =                     | mg/kg     | see below                           |                       |  |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 100                                 | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 2                                   | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 25                                  | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06                            |                       |  |  |
| ME - Matrix effect =                                 |           | 1                                   | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 70                                  | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 9125                                | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550                               | USEPA 1991, HHEM      |  |  |

| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |               |             | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|---------------------------------|--------------------------------|----------------------------|--------------|---|---------------|-------------|--|-------------|
|                        |                                 |                                |                            |              | Lifetime Daily Intake mg/kg-day         | 1/(mg/kg-day) | Cancer Risk |  |             |
| <b>Semivolatiles</b>   |                                 |                                |                            |              |   |               |             |  |             |
| Benzo(a)anthracene     | 5.90E-01                        | 4.62E-09                       | NA                         | NA           | 1.65E-09                                | 7.30E-01      | 1.20E-09    |  |             |
| Benzo(a)pyrene         | 3.90E-01                        | 3.05E-09                       | NA                         | NA           | 1.09E-09                                | 7.30E+00      | 7.96E-09    |  |             |
| Benzo(b)fluoranthene   | 5.80E-01                        | 4.54E-09                       | NA                         | NA           | 1.62E-09                                | 7.30E-01      | 1.18E-09    |  |             |
| Benzo(k)fluoranthene   | 1.90E-01                        | 1.49E-09                       | NA                         | NA           | 5.31E-10                                | 7.30E-02      | 3.88E-11    |  |             |
| Chrysene               | 5.30E-01                        | 4.15E-09                       | NA                         | NA           | 1.48E-09                                | 7.30E-03      | 1.08E-11    |  |             |
| Dibenz(a,h)anthracene  | 6.20E-02                        | 4.85E-10                       | NA                         | NA           | 1.73E-10                                | 7.30E+00      | 1.27E-09    |  |             |
| Indeno(1,2,3-cd)pyrene | 2.20E-01                        | 1.72E-09                       | NA                         | NA           | 6.15E-10                                | 7.30E-01      | 4.49E-10    |  |             |

NA - Not Applicable

Total Cancer Risk = 1.21E-08



**Table 42****Dermal Exposure to EU1 Surface Water by a Maintenance Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                   | <u><math>C_w * SA * K_p * ABS * ET * EF * ED * CF</math></u> |  |  |  |
|---|-------------------|--|--|--|--|
|   |                   | <u><math>BW * AT</math></u>                                  |  |  |  |
| Cw - Concentration in surface water =                       | mg/L              | see below  |  |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup>   | 3000   | calculated                             |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>   | 20000  | USEPA 1997, EFH                        |  |  |
| Fs - Fraction of skin surface area available for exposure = |                   | 15.0%  | USEPA 1997, EFH                        |  |  |
| Kp - Dermal permeability constant =                         | cm/hr             | see below  |  |  |  |
| ABSp - Absorption - cPAHs =                                 |                   | 0.03   | USEPA 1995, Region III                 |  |  |
| ET - Exposure time =  | hrs/day           | 1  | USEPA 1992, Dermal Exposure Assessment |  |  |
| EF - Exposure frequency =                                   | days/year         | 2  | reasonable assumption                  |  |  |
| ED - Exposure duration =                                    | years             | 25   | USEPA 1995, Region IV                  |  |  |
| CF - Conversion factor =                                    | L/cm <sup>3</sup> | 1.00E-03   |  |  |  |
| BW - Body weight =  | kg                | 70   | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days              | 9125   | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days              | 25550  | USEPA 1991, HHEM                       |  |  |

| Constituent            | Concentration in Surface Water mg/L | Kp cm/hr | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day |    | Hazard Index | Lifetime Daily Intake mg/kg-day | Average Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-------------------------------------|----------|--------------------------------|------------------------------|----|--------------|---------------------------------|---|-------------|
|                        |                                     |          |                                | NA                           | NA |              |                                 |   |             |
| <b>Semivolatiles</b>   |                                     |          |                                |                              |    |              |                                 |   |             |
| Benzo(a)anthracene     | 1.00E-03                            | 8.10E-01 | 5.71E-09                       | NA                           | NA | NA           | 2.04E-09                        | 1.46E+00                                  | 2.98E-09    |
| Benzo(a)pyrene         | 5.00E-04                            | 1.20E+00 | 4.23E-09                       | NA                           | NA | NA           | 1.51E-09                        | 1.46E+01                                  | 2.20E-08    |
| Benzo(b)fluoranthene   | 5.00E-04                            | 1.20E+00 | 4.23E-09                       | NA                           | NA | NA           | 1.51E-09                        | 1.46E+00                                  | 2.20E-09    |
| Benzo(k)fluoranthene   | 5.00E-04                            | 4.48E+01 | 1.58E-07                       | NA                           | NA | NA           | 5.64E-08                        | 1.46E-01                                  | 8.23E-09    |
| Chrysene               | 5.00E-04                            | 8.10E-01 | 2.85E-09                       | NA                           | NA | NA           | 1.02E-09                        | 1.46E-02                                  | 1.49E-11    |
| Dibenz(a,h)anthracene  | 5.00E-04                            | 2.70E+00 | 9.51E-09                       | NA                           | NA | NA           | 3.40E-09                        | 1.46E+01                                  | 4.96E-08    |
| Indeno(1,2,3-cd)pyrene | 5.00E-04                            | 1.90E+00 | 6.69E-09                       | NA                           | NA | NA           | 2.39E-09                        | 1.46E+00                                  | 3.49E-09    |

NA - Not Available

Total Cancer Risk = 8.85E-08



**Table 43**  
**Oral Exposure to EU1 Surface Water by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |  | $\frac{C_{sw} \cdot I_{ngR} \cdot E_F \cdot E_D \cdot E_T}{B_W \cdot A_T}$ |  |   |             |          |          |
|---|--|--|--|---|-------------|----------|----------|
| C <sub>sw</sub> - Concentration in surface water =              | mg/L                                   | see below  |  |   |             |          |          |
| I <sub>ngR</sub> - Ingestion rate for surface water =           | L/hour                                 | 0.01   | USEPA 1995, Region IV                  |   |             |          |          |
| E <sub>F</sub> - Exposure frequency =                           | days/year                              | 2  | reasonable assumption                  |   |             |          |          |
| E <sub>D</sub> - Exposure duration =                            | years                                  | 25   | USEPA 1995, Region IV                  |   |             |          |          |
| E <sub>T</sub> - Exposure time =                                | hrs/day                                | 1  | USEPA 1992, Dermal Exposure Assessment |   |             |          |          |
| B <sub>W</sub> - Body weight =                                  | kg                                     | 70   | USEPA 1995, Region IV                  |   |             |          |          |
| A <sub>T<sub>n</sub></sub> - Averaging time - noncarcinogenic = | days                                   | 9125   | USEPA 1991, HHEM                       |   |             |          |          |
| A <sub>T<sub>c</sub></sub> - Averaging time - carcinogenic =    | days                                   | 25550  | USEPA 1991, HHEM                       |   |             |          |          |
| Constituent   | Concentration in Surface Water<br>mg/L | Average Daily Intake<br>mg/kg-day  | Oral Chronic RfD<br>mg/kg-day          | Average                                   |             |          |          |
|   |  |  |  | Lifetime Daily Intake<br>mg/kg-day        |             |          |          |
|   |  |  | Hazard Index                           | Oral Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |          |          |
| Semivolatiles   |  |  |  |   |             |          |          |
| Benzo(a)anthracene  | 1.00E-03                               | 7.83E-10   | NA                                     | NA  | 2.80E-10    | 7.30E-01 | 2.04E-10 |
| Benzo(a)pyrene  | 5.00E-04                               | 3.91E-10   | NA                                     | NA  | 1.40E-10    | 7.30E+00 | 1.02E-09 |
| Benzo(b)fluoranthene  | 5.00E-04                               | 3.91E-10   | NA                                     | NA  | 1.40E-10    | 7.30E-01 | 1.02E-10 |
| Benzo(k)fluoranthene  | 5.00E-04                               | 3.91E-10   | NA                                     | NA  | 1.40E-10    | 7.30E-02 | 1.02E-11 |
| Chrysene  | 5.00E-04                               | 3.91E-10   | NA                                     | NA  | 1.40E-10    | 7.30E-03 | 1.02E-12 |
| Dibenz(a,h)anthracene   | 5.00E-04                               | 3.91E-10   | NA                                     | NA  | 1.40E-10    | 7.30E+00 | 1.02E-09 |
| Indeno(1,2,3-cd)pyrene  | 5.00E-04                               | 3.91E-10   | NA                                     | NA  | 1.40E-10    | 7.30E-01 | 1.02E-10 |

NA - Not Applicable

Total Cancer Risk = 2.46E-09



Table 44

*Dermal Exposure to EU2 Surface Soil (0-6') by a Maintenance Worker**Kerr McGee, Hattiesburg, MS*

| Intake (mg/kg-day) =  |                                   | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                    |                 |   |   |                |
|---|-----------------------------------|---|------------------------------------|-----------------|---|---|----------------|
| Cs - Concentration in soil =                                | mg/kg                             | chem. spec.   |                                    |                 |   |   |                |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day              | 3000  | calculated                         |                 |   |   |                |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                   | 20000   | USEPA 1997, EFH                    |                 |   |   |                |
| Fs - Fraction of skin surface area available for exposure = |                                   | 15%   | USEPA 1997, EFH                    |                 |   |   |                |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                | 0.038   | USEPA 1997, EFH                    |                 |   |   |                |
| ABSp - Absorption - cPAHs =                                 |                                   | 0.03  | USEPA 1995, Region III             |                 |   |   |                |
| EF - Exposure frequency =                                   | days/year                         | 150   | reasonable assumption              |                 |   |   |                |
| ED - Exposure duration =                                    | years                             | 25  | USEPA 1995, Region IV              |                 |   |   |                |
| CF - Conversion factor =                                    | kg/mg                             | 1.00E-06  |                                    |                 |   |   |                |
| BW - Body weight =  | kg                                | 70  | USEPA 1995, Region IV              |                 |   |   |                |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                              | 9125  | USEPA 1991, HHEM                   |                 |   |   |                |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                              | 25550   | USEPA 1991, HHEM                   |                 |   |   |                |
| Constituent   | Concentration in<br>Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day                | Dermal<br>Chronic RfD<br>mg/kg-day | Hazard<br>Index | Average Lifetime<br>Daily Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
| Semivolatiles   |                                   |   |                                    |                 |   |   |                |
| Benzo(a)anthracene  | 2.80E+00                          | 5.62E-08  | NA                                 | NA              | 2.01E-08                                      | 1.46E+00                                | 2.93E-08       |
| Benzo(a)pyrene  | 2.64E+00                          | 5.30E-08  | NA                                 | NA              | 1.89E-08                                      | 1.46E+01                                | 2.76E-07       |
| Benzo(b)fluoranthene  | 9.20E+00                          | 1.85E-07  | NA                                 | NA              | 6.60E-08                                      | 1.46E+00                                | 9.63E-08       |
| Benzo(k)fluoranthene  | 1.84E+00                          | 3.69E-08  | NA                                 | NA              | 1.32E-08                                      | 1.46E-01                                | 1.93E-09       |
| Chrysene  | 5.33E+00                          | 1.07E-07  | NA                                 | NA              | 3.82E-08                                      | 1.46E-02                                | 5.58E-10       |
| Dibenz(a,h)anthracene                                       | 2.39E-01                          | 4.80E-09  | NA                                 | NA              | 1.71E-09                                      | 1.46E+01                                | 2.50E-08       |
| Indeno(1,2,3-cd)pyrene                                      | 1.97E+00                          | 3.96E-08  | NA                                 | NA              | 1.41E-08                                      | 1.46E+00                                | 2.06E-08       |

NA - Not Available

Total Cancer Risk = 4.50E-07



**Table 45**

**Oral Exposure to EU2 Surface Soil (0-6') by a Maintenance Worker**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u>Cd*Ingr*EF*ED*CF*ME</u> |                       |  |  |
|--|-----------|----------------------------|-----------------------|--|--|
|  |           | BW*AT                      |                       |  |  |
| Cd - Concentration in sediment =                     | mg/kg     | see below                  |                       |  |  |
| Ingr - Ingestion rate for soil =                     | mg/day    | 100                        | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 150                        | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 25                         | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06                   |                       |  |  |
| ME - Matrix effect =                                 |           | 1                          | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 70                         | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 9125                       | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550                      | USEPA 1991, HHEM      |  |  |

| Constituent            | Soil mg/kg | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |          |          | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|------------|--------------------------------|----------------------------|--------------|---|----------|----------|--|-------------|
|                        |            |                                |                            |              |   |          |          |  |             |
| <b>Semivolatiles</b>   |            |                                |                            |              |   |          |          |  |             |
| Benzo(a)anthracene     | 2.80E+00   | 1.64E-06                       | NA                         | NA           | 5.87E-07                                | 7.30E-01 | 4.29E-07 |  |             |
| Benzo(a)pyrene         | 2.64E+00   | 1.55E-06                       | NA                         | NA           | 5.54E-07                                | 7.30E+00 | 4.04E-06 |  |             |
| Benzo(b)fluoranthene   | 9.20E+00   | 5.40E-06                       | NA                         | NA           | 1.93E-06                                | 7.30E-01 | 1.41E-06 |  |             |
| Benzo(k)fluoranthene   | 1.84E+00   | 1.08E-06                       | NA                         | NA           | 3.86E-07                                | 7.30E-02 | 2.82E-08 |  |             |
| Chrysene               | 5.33E+00   | 3.13E-06                       | NA                         | NA           | 1.12E-06                                | 7.30E-03 | 8.16E-09 |  |             |
| Dibenz(a,h)anthracene  | 2.39E-01   | 1.40E-07                       | NA                         | NA           | 5.01E-08                                | 7.30E+00 | 3.66E-07 |  |             |
| Indeno(1,2,3-cd)pyrene | 1.97E+00   | 1.16E-06                       | NA                         | NA           | 4.13E-07                                | 7.30E-01 | 3.02E-07 |  |             |

NA - Not Applicable

Total Cancer Risk = 6.58E-06



**Table 46**  
**Dermal Exposure to EU4 Sediment by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |                        |  |  |
|---|----------------------|------------------------------|------------------------|--|--|
|   |                      | BW*AT                        |                        |  |  |
| Cs - Concentration in sediment =                            | mg/kg                | chem. spec.                  |                        |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 3000                         | calculated             |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000                        | USEPA 1997, EFH        |  |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 15.0%                        | USEPA 1997, EFH        |  |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.038                        | USEPA 1997, EFH        |  |  |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                      | 0.03                         | USEPA 1995, Region III |  |  |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |                      | 0.1                          | USEPA 1995, Region III |  |  |
| EF - Exposure frequency =                                   | days/year            | 2                            | reasonable assumption  |  |  |
| ED - Exposure duration =                                    | years                | 25                           | USEPA 1995, Region IV  |  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06                     |                        |  |  |
| BW - Body weight =  | kg                   | 70                           | USEPA 1995, Region IV  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 9125                         | USEPA 1991, HHEM       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550                        | USEPA 1991, HHEM       |  |  |

| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Dermal                |              | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|---------------------------------|--------------------------------|-----------------------|--------------|---|-----------------------------------|-------------|
|                        |                                 |                                | Chronic RfD mg/kg-day | Hazard Index |   |                                   |             |
| <b>Semivolatiles</b>   |                                 |                                |                       |              |   |                                   |             |
| Benzo(a)anthracene     | 3.30E+02                        | 8.83E-08                       | NA                    | NA           | 3.16E-08                                | 1.46E+00                          | 4.61E-08    |
| Benzo(a)pyrene         | 1.30E+02                        | 3.48E-08                       | NA                    | NA           | 1.24E-08                                | 1.46E+01                          | 1.81E-07    |
| Benzo(b)fluoranthene   | 1.80E+02                        | 4.82E-08                       | NA                    | NA           | 1.72E-08                                | 1.46E+00                          | 2.51E-08    |
| Benzo(k)fluoranthene   | 6.40E+01                        | 1.71E-08                       | NA                    | NA           | 6.12E-09                                | 1.46E-01                          | 8.93E-10    |
| Carbazole              | 5.90E+02                        | 5.26E-07                       | NA                    | NA           | 1.88E-07                                | 2.00E-02                          | 3.76E-09    |
| Chrysene               | 2.90E+02                        | 7.76E-08                       | NA                    | NA           | 2.77E-08                                | 1.46E-02                          | 4.05E-10    |
| Diben(a,h)anthracene   | 1.20E+01                        | 3.21E-09                       | NA                    | NA           | 1.15E-09                                | 1.46E+01                          | 1.68E-08    |
| Dibenzofuran           | 9.40E+02                        | 8.39E-07                       | 2.00E-03              | 4.19E-04     | 3.00E-07                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene | 4.70E+01                        | 1.26E-08                       | NA                    | NA           | 4.49E-09                                | 1.46E+00                          | 6.56E-09    |
| Naphthalene            | 3.00E+03                        | 2.68E-06                       | 1.00E-02              | 2.68E-04     | 9.56E-07                                | NA                                | NA          |
| Phenanthrene           | 3.20E+03                        | 2.86E-06                       | NA                    | NA           | 1.02E-06                                | NA                                | NA          |

NA - Not Available

Total Hazard Index = 6.87E-04

Total Cancer Risk = 2.81E-07



**Table 47****Oral Exposure to EU4 Sediment by a Maintenance Worker**

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =   |                                 | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                            |              |   |  |             |
|------------------------|---------------------------------|---|----------------------------|--------------|---|--|-------------|
| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day  | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>   |                                 |   |                            |              |   |  |             |
| Benzo(a)anthracene     | 3.30E+02                        | 2.58E-06  | NA                         | NA           | 9.23E-07                                | 7.30E-01                               | 6.73E-07    |
| Benzo(a)pyrene         | 1.30E+02                        | 1.02E-06  | NA                         | NA           | 3.63E-07                                | 7.30E+00                               | 2.65E-06    |
| Benzo(b)fluoranthene   | 1.80E+02                        | 1.41E-06  | NA                         | NA           | 5.03E-07                                | 7.30E-01                               | 3.67E-07    |
| Benzo(k)fluoranthene   | 6.40E+01                        | 5.01E-07  | NA                         | NA           | 1.79E-07                                | 7.30E-02                               | 1.31E-08    |
| Carbazole              | 5.90E+02                        | 4.62E-06  | NA                         | NA           | 1.65E-06                                | 2.00E-02                               | 3.30E-08    |
| Chrysene               | 2.90E+02                        | 2.27E-06  | NA                         | NA           | 8.11E-07                                | 7.30E-03                               | 5.92E-09    |
| Dibenz(a,h)anthracene  | 1.20E+01                        | 9.39E-08  | NA                         | NA           | 3.35E-08                                | 7.30E+00                               | 2.45E-07    |
| Dibenzofuran           | 9.40E+02                        | 7.36E-06  | 4.00E-03                   | 1.84E-03     | 2.63E-06                                | NA                                     | NA          |
| Indeno(1,2,3-cd)pyrene | 4.70E+01                        | 3.68E-07  | NA                         | NA           | 1.31E-07                                | 7.30E-01                               | 9.59E-08    |
| Naphthalene            | 3.00E+03                        | 2.35E-05  | 2.00E-02                   | 1.17E-03     | 8.39E-06                                | NA                                     | NA          |
| Phenanthrene           | 3.20E+03                        | 2.50E-05  | NA                         | NA           | 8.95E-06                                | NA                                     | NA          |

NA - Not Applicable

Total Hazard Index = 3.01E-03

Total Cancer Risk = 4.09E-06



**Table 48****Dermal Exposure to EU4 Surface Water by a Maintenance Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | Cw*SA*Kp*ABS*ET*EF*ED*CF<br>BW*AT |
|---|-----------------------------------|
| Cw - Concentration in surface water =                                   | mg/L                              |
| SA - Surface area available for exposure =                              | cm <sup>2</sup>                   |
| SA <sub>t</sub> - Total skin surface area =                             | cm <sup>2</sup>                   |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = | 15.0%                             |
| K <sub>p</sub> - Dermal permeability constant =                         | cm/hr                             |
| ABS <sub>p</sub> - Absorption - cPAHs =                                 | 0.03                              |
| ABS <sub>s</sub> - Absorption - other SVOCs =                           | 0.1                               |
| ET - Exposure time =  | hrs/day                           |
| EF - Exposure frequency =   | days/year                         |
| ED - Exposure duration =  | years                             |
| CF - Conversion factor =  | L/cm <sup>3</sup>                 |
| BW - Body weight =  | kg                                |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                    | days                              |
| AT <sub>c</sub> - Averaging time - carcinogenic =                       | days                              |

| Constituent                | Concentration in Surface |                      | Average Daily Intake<br>mg/kg-day | Dermal Chronic RfD<br>mg/kg-day | Hazard Index | Average Lifetime Daily Intake<br>mg/kg-day |                                      |          | Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
|----------------------------|--------------------------|----------------------|-----------------------------------|---------------------------------|--------------|--|--------------------------------------|----------|--------------------------------------|-------------|
|                            | Water mg/L               | K <sub>p</sub> cm/hr |                                   |                                 |              | Lifetime Daily Intake<br>mg/kg-day         | Cancer Slope Factor<br>1/(mg/kg-day) |          |                                      |             |
| <b>Semivolatiles</b>       |                          |                      |                                   |                                 |              |  |                                      |          |                                      |             |
| Benzo(a)anthracene         | 5.00E-03                 | 8.10E-01             | 2.85E-08                          | NA                              | NA           | 1.02E-08                                   | 1.46E+00                             | 1.49E-08 |                                      |             |
| Benzo(a)pyrene             | 5.00E-04                 | 1.20E+00             | 4.23E-09                          | NA                              | NA           | 1.51E-09                                   | 1.46E+01                             | 2.20E-08 |                                      |             |
| Benzo(b)fluoranthene       | 1.20E-02                 | 1.20E+00             | 1.01E-07                          | NA                              | NA           | 3.62E-08                                   | 1.46E+00                             | 5.29E-08 |                                      |             |
| Benzo(k)fluoranthene       | 2.00E-03                 | 4.48E+01             | 6.31E-07                          | NA                              | NA           | 2.25E-07                                   | 1.46E-01                             | 3.29E-08 |                                      |             |
| Bis(2-ethylhexyl)phthalate | 3.00E-03                 | 3.30E-02             | 2.32E-09                          | 1.00E-02                        | 2.32E-07     | 8.30E-10                                   | 1.40E-02                             | 1.16E-11 |                                      |             |
| Carbazole                  | 1.00E-02                 | 3.57E-02             | 8.39E-09                          | NA                              | NA           | 3.00E-09                                   | 2.00E-02                             | 6.00E-11 |                                      |             |
| Chrysene                   | 6.00E-03                 | 8.10E-01             | 3.42E-08                          | NA                              | NA           | 1.22E-08                                   | 1.46E-02                             | 1.79E-10 |                                      |             |
| Dibenz(a,h)anthracene      | 5.00E-04                 | 2.70E+00             | 9.51E-09                          | NA                              | NA           | 3.40E-09                                   | 1.46E+01                             | 4.96E-08 |                                      |             |
| Dibenzofuran               | 1.10E-02                 | 1.51E-01             | 3.90E-08                          | 2.00E-03                        | 1.95E-05     | 1.39E-08                                   | NA                                   | NA       |                                      |             |
| Indeno(1,2,3-cd)pyrene     | 5.00E-04                 | 1.90E+00             | 6.69E-09                          | NA                              | NA           | 2.39E-09                                   | 1.46E+00                             | 3.49E-09 |                                      |             |
| Phenanthrene               | 1.70E-02                 | 2.30E-01             | 9.18E-08                          | NA                              | NA           | 3.28E-08                                   | NA                                   | NA       |                                      |             |

NA - Not Available

Total Hazard Index = 1.97E-05

Total Cancer Risk = 1.76E-07



**Table 49****Oral Exposure to EU4 Surface Water by a Maintenance Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u><math>C_{sw} * IngR * EF * ED * ET</math></u> |  |  |  |  |
|--|-----------|--|--|--|--|--|
|  |           | <u><math>BW * AT</math></u>                      |  |  |  |  |
| Csw - Concentration in surface water =               | mg/L      | see below  |  |  |  |  |
| IngR - Ingestion rate for surface water =            | L/hour    | 0.01   | USEPA 1995, Region IV                  |  |  |  |
| EF - Exposure frequency =                            | days/year | 2  | reasonable assumption                  |  |  |  |
| ED - Exposure duration =                             | years     | 25   | USEPA 1995, Region IV                  |  |  |  |
| ET - Exposure time =                                 | hrs/day   | 1  | USEPA 1992, Dermal Exposure Assessment |  |  |  |
| BW - Body weight =                                   | kg        | 70   | USEPA 1995, Region IV                  |  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 9125   | USEPA 1991, HHEM                       |  |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550  | USEPA 1991, HHEM                       |  |  |  |

| Constituent                | Concentration         | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average                         | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|----------------------------|-----------------------|--------------------------------|----------------------------|--------------|---------------------------------|--|-------------|
|                            | in Surface Water mg/L |                                |                            |              | Lifetime Daily Intake mg/kg-day |  |             |
| <b>Semivolatiles</b>       |                       |                                |                            |              |                                 |  |             |
| Benzo(a)anthracene         | 5.00E-03              | 3.91E-09                       | NA                         | NA           | 1.40E-09                        | 7.30E-01                               | 1.02E-09    |
| Benzo(a)pyrene             | 5.00E-04              | 3.91E-10                       | NA                         | NA           | 1.40E-10                        | 7.30E+00                               | 1.02E-09    |
| Benzo(b)fluoranthene       | 1.20E-02              | 9.39E-09                       | NA                         | NA           | 3.35E-09                        | 7.30E-01                               | 2.45E-09    |
| Benzo(k)fluoranthene       | 2.00E-03              | 1.57E-09                       | NA                         | NA           | 5.59E-10                        | 7.30E-02                               | 4.08E-11    |
| Bis(2-ethylhexyl)phthalate | 3.00E-03              | 2.35E-09                       | 2.00E-02                   | 1.17E-07     | 8.39E-10                        | 1.40E-02                               | 1.17E-11    |
| Carbazole                  | 1.00E-02              | 7.83E-09                       | NA                         | NA           | 2.80E-09                        | 2.00E-02                               | 5.59E-11    |
| Chrysene                   | 6.00E-03              | 4.70E-09                       | NA                         | NA           | 1.68E-09                        | 7.30E-03                               | 1.22E-11    |
| Dibenz(a,h)anthracene      | 5.00E-04              | 3.91E-10                       | NA                         | NA           | 1.40E-10                        | 7.30E+00                               | 1.02E-09    |
| Dibenzofuran               | 1.10E-02              | 8.61E-09                       | 4.00E-03                   | 2.15E-06     | 3.08E-09                        | NA                                     | NA          |
| Indeno(1,2,3-cd)pyrene     | 5.00E-04              | 3.91E-10                       | NA                         | NA           | 1.40E-10                        | 7.30E-01                               | 1.02E-10    |
| Phenanthrene               | 1.70E-02              | 1.33E-08                       | NA                         | NA           | 4.75E-09                        | NA                                     | NA          |

NA - Not Applicable

Total Hazard Index = 2.27E-06

Total Cancer Risk = 5.73E-09



**Table 50**  
**Dermal Exposure to EU4 Surface Soil (0-6') by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | $C_s * SA * AH * ABS * EF * ED * CF$ |                        |  |  |
|---|----------------------|--------------------------------------|------------------------|--|--|
|   |                      | BW * AT                              |                        |  |  |
| $C_s$ - Concentration in soil =                             | mg/kg                | chem. spec.                          |                        |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 3000                                 | calculated             |  |  |
| SA <sub>s</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000                                | USEPA 1997, EFH        |  |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 15%                                  | USEPA 1997, EFH        |  |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.038                                | USEPA 1997, EFH        |  |  |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                      | 0.03                                 | USEPA 1995, Region III |  |  |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |                      | 0.1                                  | USEPA 1995, Region III |  |  |
| EF - Exposure frequency =                                   | days/year            | 150                                  | reasonable assumption  |  |  |
| ED - Exposure duration =                                    | years                | 25                                   | USEPA 1995, Region IV  |  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06                             |                        |  |  |
| BW - Body weight =  | kg                   | 70                                   | USEPA 1995, Region IV  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 9125                                 | USEPA 1991, HHEM       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550                                | USEPA 1991, HHEM       |  |  |

| Constituent            | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal                      |                 | Average                                  |   |          | Cancer Risk |
|------------------------|-----------------------------------|--------------------------------------|-----------------------------|-----------------|--|---|----------|-------------|
|                        |                                   |                                      | Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Lifetime<br>Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) |          |             |
| <b>Semivolatiles</b>   |                                   |                                      |                             |                 |  |   |          |             |
| Benzo(a)anthracene     | 9.30E+02                          | 1.87E-05                             | NA                          | NA              | 6.67E-06                                 | 1.46E+00                                | 9.74E-06 |             |
| Benzo(a)pyrene         | 5.00E+02                          | 1.00E-05                             | NA                          | NA              | 3.59E-06                                 | 1.46E+01                                | 5.23E-05 |             |
| Benzo(b)fluoranthene   | 5.30E+02                          | 1.06E-05                             | NA                          | NA              | 3.80E-06                                 | 1.46E+00                                | 5.55E-06 |             |
| Benzo(k)fluoranthene   | 2.90E+02                          | 5.82E-06                             | NA                          | NA              | 2.08E-06                                 | 1.46E-01                                | 3.04E-07 |             |
| Carbazole              | 6.20E+02                          | 4.15E-05                             | NA                          | NA              | 1.48E-05                                 | 2.00E-02                                | 2.96E-07 |             |
| Chrysene               | 6.90E+02                          | 1.39E-05                             | NA                          | NA              | 4.95E-06                                 | 1.46E-02                                | 7.22E-08 |             |
| Dibenz(a,h)anthracene  | 6.40E+01                          | 1.29E-06                             | NA                          | NA              | 4.59E-07                                 | 1.46E+01                                | 6.70E-06 |             |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                          | 5.02E-06                             | NA                          | NA              | 1.79E-06                                 | 1.46E+00                                | 2.62E-06 |             |
| Naphthalene            | 3.50E+03                          | 2.34E-04                             | 1.00E-02                    | 2.34E-02        | 8.37E-05                                 | NA                                      | NA       |             |

NA - Not Available

Total Hazard Index = 2.34E-02

Total Cancer Risk = 7.76E-05



**Table 51**

**Oral Exposure to EU4 Surface Soil (0-6') by a Maintenance Worker**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =   |                             | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                            |              |   |  |             |
|------------------------|-----------------------------|---|----------------------------|--------------|---|--|-------------|
| Constituent            | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day  | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles          |                             |   |                            |              |   |  |             |
| Benzo(a)anthracene     | 9.30E+02                    | 5.46E-04  | NA                         | NA           | 1.95E-04                                | 7.30E-01                               | 1.42E-04    |
| Benzo(a)pyrene         | 5.00E+02                    | 2.94E-04  | NA                         | NA           | 1.05E-04                                | 7.30E+00                               | 7.65E-04    |
| Benzo(b)fluoranthene   | 5.30E+02                    | 3.11E-04  | NA                         | NA           | 1.11E-04                                | 7.30E-01                               | 8.11E-05    |
| Benzo(k)fluoranthene   | 2.90E+02                    | 1.70E-04  | NA                         | NA           | 6.08E-05                                | 7.30E-02                               | 4.44E-06    |
| Carbazole              | 6.20E+02                    | 3.64E-04  | NA                         | NA           | 1.30E-04                                | 2.00E-02                               | 2.60E-06    |
| Chrysene               | 6.90E+02                    | 4.05E-04  | NA                         | NA           | 1.45E-04                                | 7.30E-03                               | 1.06E-06    |
| Dibenz(a,h)anthracene  | 6.40E+01                    | 3.76E-05  | NA                         | NA           | 1.34E-05                                | 7.30E+00                               | 9.80E-05    |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                    | 1.47E-04  | NA                         | NA           | 5.24E-05                                | 7.30E-01                               | 3.83E-05    |
| Naphthalene            | 3.50E+03                    | 2.05E-03  | 2.00E-02                   | 1.03E-01     | 7.34E-04                                | NA                                     | NA          |

NA - Not Applicable

Total Hazard Index = 1.03E-01

Total Cancer Risk = 1.13E-03



**Table 52**  
**Dermal Exposure to EU5 Surface Soil (0-6') by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                             | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                              |              |   |                                   |             |
|---|-----------------------------|---|------------------------------|--------------|---|-----------------------------------|-------------|
| Cs - Concentration in soil =                                | mg/kg                       | chem. spec.   |                              |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day        | 3000  | calculated                   |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>             | 20000   | USEPA 1997, EFH              |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |                             | 15%   | USEPA 1997, EFH              |              |   |                                   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>          | 0.038   | USEPA 1997, EFH              |              |   |                                   |             |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                             | 0.03  | USEPA 1995, Region III       |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year                   | 150   | reasonable assumption        |              |   |                                   |             |
| ED - Exposure duration =                                    | years                       | 25  | USEPA 1995, Region IV        |              |   |                                   |             |
| CF - Conversion factor =                                    | kg/mg                       | 1.00E-06  |                              |              |   |                                   |             |
| BW - Body weight =  | kg                          | 70  | USEPA 1995, Region IV        |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                        | 9125  | USEPA 1991, HHEM             |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                        | 25550   | USEPA 1991, HHEM             |              |   |                                   |             |
| Constituent   | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day                      | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |                             |   |                              |              |   |                                   |             |
| Benzo(a)anthracene  | 7.77E+01                    | 1.56E-06  | NA                           | NA           | 5.57E-07                                | 1.46E+00                          | 8.13E-07    |
| Benzo(a)pyrene  | 4.10E+01                    | 8.23E-07  | NA                           | NA           | 2.94E-07                                | 1.46E+01                          | 4.29E-06    |
| Benzo(b)fluoranthene  | 7.95E+01                    | 1.60E-06  | NA                           | NA           | 5.70E-07                                | 1.46E+00                          | 8.32E-07    |
| Benzo(k)fluoranthene  | 1.97E+01                    | 3.96E-07  | NA                           | NA           | 1.41E-07                                | 1.46E-01                          | 2.06E-08    |
| Chrysene  | 8.25E+01                    | 1.66E-06  | NA                           | NA           | 5.92E-07                                | 1.46E-02                          | 8.64E-09    |
| Dibenz(a,h)anthracene                                       | 2.04E+00                    | 4.10E-08  | NA                           | NA           | 1.46E-08                                | 1.46E+01                          | 2.14E-07    |
| Indeno(1,2,3-cd)pyrene                                      | 1.71E+01                    | 3.43E-07  | NA                           | NA           | 1.23E-07                                | 1.46E+00                          | 1.79E-07    |

NA - Not Available

Total Cancer Risk = 6.36E-06



**Table 53****Oral Exposure to EU5 Surface Soil (0-6') by a Maintenance Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u>Cd*IngR*EF*ED*CF*ME</u> |                       |  |
|--|-----------|----------------------------|-----------------------|--|
|  |           | <u>BW*AT</u>               |                       |  |
| Cd - Concentration in sediment =                     | mg/kg     | see below                  |                       |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 100                        | USEPA 1997, EFH       |  |
| EF - Exposure frequency =                            | days/year | 150                        | reasonable assumption |  |
| ED - Exposure duration =                             | years     | 25                         | USEPA 1995, Region IV |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06                   |                       |  |
| ME - Matrix effect =                                 |           | 1                          | Magee, et al., 1996   |  |
| BW - Body weight =                                   | kg        | 70                         | USEPA 1995, Region IV |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 9125                       | USEPA 1991, HHEM      |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550                      | USEPA 1991, HHEM      |  |

| Constituent            | Concentration in<br>Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average Lifetime<br>Daily Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------------|--------------------------------------|----------------------------------|-----------------|---|--|-------------|
| <b>Semivolatiles</b>   |                                   |                                      |                                  |                 |   |  |             |
| Benzo(a)anthracene     | 7.77E+01                          | 4.56E-05                             | NA                               | NA              | 1.63E-05                                      | 7.30E-01                                     | 1.19E-05    |
| Benzo(a)pyrene         | 4.10E+01                          | 2.41E-05                             | NA                               | NA              | 8.60E-06                                      | 7.30E+00                                     | 6.28E-05    |
| Benzo(b)fluoranthene   | 7.95E+01                          | 4.67E-05                             | NA                               | NA              | 1.67E-05                                      | 7.30E-01                                     | 1.22E-05    |
| Benzo(k)fluoranthene   | 1.97E+01                          | 1.16E-05                             | NA                               | NA              | 4.13E-06                                      | 7.30E-02                                     | 3.02E-07    |
| Chrysene               | 8.25E+01                          | 4.84E-05                             | NA                               | NA              | 1.73E-05                                      | 7.30E-03                                     | 1.26E-07    |
| Dibenz(a,h)anthracene  | 2.04E+00                          | 1.20E-06                             | NA                               | NA              | 4.28E-07                                      | 7.30E+00                                     | 3.12E-06    |
| Indeno(1,2,3-cd)pyrene | 1.71E+01                          | 1.00E-05                             | NA                               | NA              | 3.59E-06                                      | 7.30E-01                                     | 2.62E-06    |

NA - Not Applicable

Total Cancer Risk = 9.30E-05



**Table 55**

**Oral Exposure to EU1 Sediment by a Construction Worker**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                       |  |  |
|--|-----------|---|-----------------------|--|--|
| Cd - Concentration in sediment =                     | mg/kg     | see below   |                       |  |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 480   | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 8   | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 1   | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06  |                       |  |  |
| ME - Matrix effect =                                 |           | 1   | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 70  | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 365   | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550   | USEPA 1991, HHEM      |  |  |

| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |  |             |
|------------------------|---------------------------------|--------------------------------|----------------------------|--------------|---|--|-------------|
|                        |                                 |                                |                            |              | Lifetime Daily Intake mg/kg-day         | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>   |                                 |                                |                            |              |   |  |             |
| Benzo(a)anthracene     | 5.90E-01                        | 8.87E-08                       | NA                         | NA           | 1.27E-09                                | 7.30E-01                               | 9.25E-10    |
| Benzo(a)pyrene         | 3.90E-01                        | 5.86E-08                       | NA                         | NA           | 8.37E-10                                | 7.30E+00                               | 6.11E-09    |
| Benzo(b)fluoranthene   | 5.80E-01                        | 8.72E-08                       | NA                         | NA           | 1.25E-09                                | 7.30E-01                               | 9.09E-10    |
| Benzo(k)fluoranthene   | 1.90E-01                        | 2.86E-08                       | NA                         | NA           | 4.08E-10                                | 7.30E-02                               | 2.98E-11    |
| Chrysene               | 5.30E-01                        | 7.97E-08                       | NA                         | NA           | 1.14E-09                                | 7.30E-03                               | 8.31E-12    |
| Dibenz(a,h)anthracene  | 6.20E-02                        | 9.32E-09                       | NA                         | NA           | 1.33E-10                                | 7.30E+00                               | 9.72E-10    |
| Indeno(1,2,3-cd)pyrene | 2.20E-01                        | 3.31E-08                       | NA                         | NA           | 4.72E-10                                | 7.30E-01                               | 3.45E-10    |

NA - Not Applicable

Total Cancer Risk = 9.30E-09



**Table 56****Dermal Exposure to EU1 Surface Water by a Construction Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | $\frac{C_w \cdot SA \cdot K_p \cdot ABS \cdot ET \cdot EF \cdot ED \cdot CF}{BW \cdot AT}$ |           |  |                                 |              |   |                                   |             |
|---|--|-----------|--|---------------------------------|--------------|---|-----------------------------------|-------------|
| Cw - Concentration in surface water =                       | mg/L   | see below |  |                                 |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup>  | 3000      | calculated                             |                                 |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>  | 20000     | USEPA 1997, EFH                        |                                 |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |  | 15.0%     | USEPA 1997, EFH                        |                                 |              |   |                                   |             |
| Kp - Dermal permeability constant =                         | cm/hr  | see below |  |                                 |              |   |                                   |             |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |  | 0.03      | USEPA 1995, Region III                 |                                 |              |   |                                   |             |
| ET - Exposure time =  | hrs/day  | 1         | USEPA 1992, Dermal Exposure Assessment |                                 |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year  | 8         | reasonable assumption                  |                                 |              |   |                                   |             |
| ED - Exposure duration =                                    | years  | 1         | reasonable assumption                  |                                 |              |   |                                   |             |
| CF - Conversion factor =                                    | L/cm <sup>3</sup>  | 1.00E-03  |  |                                 |              |   |                                   |             |
| BW - Body weight =  | kg   | 70        | USEPA 1995, Region IV                  |                                 |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days   | 365       | USEPA 1991, HHEM                       |                                 |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days   | 25550     | USEPA 1991, HHEM                       |                                 |              |   |                                   |             |
| Constituent   | Concentration in Surface Water mg/L  | Kp cm/hr  | Average Daily Intake mg/kg-day         | Dermal Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |  |           |  |                                 |              |   |                                   |             |
| Benzo(a)anthracene  | 1.00E-03   | 8.10E-01  | 2.28E-08                               | NA                              | NA           | 3.26E-10                                | 1.46E+00                          | 4.76E-10    |
| Benzo(a)pyrene  | 5.00E-04   | 1.20E+00  | 1.69E-08                               | NA                              | NA           | 2.42E-10                                | 1.46E+01                          | 3.53E-09    |
| Benzo(b)fluoranthene  | 5.00E-04   | 1.20E+00  | 1.69E-08                               | NA                              | NA           | 2.42E-10                                | 1.46E+00                          | 3.53E-10    |
| Benzo(k)fluoranthene  | 5.00E-04   | 4.48E+01  | 6.31E-07                               | NA                              | NA           | 9.02E-09                                | 1.46E-01                          | 1.32E-09    |
| Chrysene  | 5.00E-04   | 8.10E-01  | 1.14E-08                               | NA                              | NA           | 1.63E-10                                | 1.46E-02                          | 2.38E-12    |
| Dibenz(a,h)anthracene                                       | 5.00E-04   | 2.70E+00  | 3.80E-08                               | NA                              | NA           | 5.43E-10                                | 1.46E+01                          | 7.93E-09    |
| Indeno(1,2,3-cd)pyrene                                      | 5.00E-04   | 1.90E+00  | 2.68E-08                               | NA                              | NA           | 3.82E-10                                | 1.46E+00                          | 5.58E-10    |

NA - Not Available

Total Cancer Risk = 1.42E-08



**Table 57**  
**Oral Exposure to EU1 Surface Water by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |  | <u>Csw*IngR*EF*ED*ET</u>          |  |              |  |          |   |             |  |
|--|--|-----------------------------------|--|--------------|--|----------|---|-------------|--|
|  |  | BW*AT                             |  |              |  |          |   |             |  |
| Csw - Concentration in surface water =               | mg/L                                   | see below                         |  |              |  |          |   |             |  |
| IngR - Ingestion rate for surface water =            | L/hour                                 | 0.01                              | USEPA 1995, Region IV                  |              |  |          |   |             |  |
| EF - Exposure frequency =                            | days/year                              | 8                                 | reasonable assumption                  |              |  |          |   |             |  |
| ED - Exposure duration =                             | years                                  | 1                                 | USEPA 1995, Region IV                  |              |  |          |   |             |  |
| ET - Exposure time =                                 | hrs/day                                | 1                                 | USEPA 1992, Dermal Exposure Assessment |              |  |          |   |             |  |
| BW - Body weight =                                   | kg                                     | 70                                | USEPA 1995, Region IV                  |              |  |          |   |             |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                                   | 365                               | USEPA 1991, HHEM                       |              |  |          |   |             |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                                   | 25550                             | USEPA 1991, HHEM                       |              |  |          |   |             |  |
| Constituent  | Concentration in Surface Water<br>mg/L | Average Daily Intake<br>mg/kg-day | Oral Subchronic RfD<br>mg/kg-day       | Hazard Index | Average Lifetime Daily Intake<br>mg/kg-day |          | Oral Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |  |
|  |  |                                   |  |              |  |          |   |             |  |
| <b>Semivolatiles</b>                                 |  |                                   |  |              |  |          |   |             |  |
| Benzo(a)anthracene                                   | 1.00E-03                               | 3.13E-09                          | NA                                     | NA           | 4.47E-11                                   | 7.30E-01 | 3.27E-11                                  |             |  |
| Benzo(a)pyrene                                       | 5.00E-04                               | 1.57E-09                          | NA                                     | NA           | 2.24E-11                                   | 7.30E+00 | 1.63E-10                                  |             |  |
| Benzo(b)fluoranthene                                 | 5.00E-04                               | 1.57E-09                          | NA                                     | NA           | 2.24E-11                                   | 7.30E-01 | 1.63E-11                                  |             |  |
| Benzo(k)fluoranthene                                 | 5.00E-04                               | 1.57E-09                          | NA                                     | NA           | 2.24E-11                                   | 7.30E-02 | 1.63E-12                                  |             |  |
| Chrysene   | 5.00E-04                               | 1.57E-09                          | NA                                     | NA           | 2.24E-11                                   | 7.30E-03 | 1.63E-13                                  |             |  |
| Dibenz(a,h)anthracene                                | 5.00E-04                               | 1.57E-09                          | NA                                     | NA           | 2.24E-11                                   | 7.30E+00 | 1.63E-10                                  |             |  |
| Indeno(1,2,3-cd)pyrene                               | 5.00E-04                               | 1.57E-09                          | NA                                     | NA           | 2.24E-11                                   | 7.30E-01 | 1.63E-11                                  |             |  |

NA - Not Applicable

Total Cancer Risk = 3.94E-10



**Table 58**

**Dermal Exposure to EU2 Soil (0-10') by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |                        |  |
|---|----------------------|------------------------------|------------------------|--|
|   |                      | BW*AT                        |                        |  |
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.                  |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 5560                         | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000                        | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 27.8%                        | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.1                          | USEPA 1997, EFH        |  |
| ABS <sub>b</sub> - Absorption - cPAHs =                     |                      | 0.03                         | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year            | 80                           | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                | 1                            | reasonable assumption  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06                     |                        |  |
| BW - Body weight =  | kg                   | 70                           | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 365                          | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550                        | USEPA 1991, HHEM       |  |

| Constituent            | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal<br>Subchronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily<br>Cancer Slope |                         |          | Cancer<br>Risk |
|------------------------|-----------------------------------|--------------------------------------|--|-----------------|---|-------------------------|----------|----------------|
|                        |                                   |                                      |  |                 | Intake<br>mg/kg-day                       | Factor<br>1/(mg/kg-day) |          |                |
| <b>Semivolatiles</b>   |                                   |                                      |  |                 |   |                         |          |                |
| Benzo(a)anthracene     | 6.70E+00                          | 3.50E-07                             | NA                                       | NA              | 5.00E-09                                  | 1.46E+00                | 7.30E-09 |                |
| Benzo(a)pyrene         | 5.20E+00                          | 2.72E-07                             | NA                                       | NA              | 3.88E-09                                  | 1.46E+01                | 5.66E-08 |                |
| Benzo(b)fluoranthene   | 9.20E+00                          | 4.80E-07                             | NA                                       | NA              | 6.86E-09                                  | 1.46E+00                | 1.00E-08 |                |
| Benzo(k)fluoranthene   | 3.60E+00                          | 1.88E-07                             | NA                                       | NA              | 2.69E-09                                  | 1.46E-01                | 3.92E-10 |                |
| Chrysene               | 5.33E+00                          | 2.78E-07                             | NA                                       | NA              | 3.98E-09                                  | 1.46E-02                | 5.81E-11 |                |
| Dibenz(a,h)anthracene  | 2.39E-01                          | 1.25E-08                             | NA                                       | NA              | 1.78E-10                                  | 1.46E+01                | 2.60E-09 |                |
| Indeno(1,2,3-cd)pyrene | 3.70E+00                          | 1.93E-07                             | NA                                       | NA              | 2.76E-09                                  | 1.46E+00                | 4.03E-09 |                |

NA - Not Available

Total Cancer Risk = 8.10E-08



Table 59

Oral Exposure to EU2 Soil (0-10') by a Construction Worker

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =                                 | <u><math>Cd * IngR * EF * ED * CF * ME</math></u> |           |                       |
|--|---|-----------|-----------------------|
|  | BW  | AT        |                       |
| Cd - Concentration in soil =                         | mg/kg   | see below |                       |
| IngR <sub>a</sub> - Ingestion rate for soil =        | mg/day  | 480       | USEPA 1997, EFH       |
| IngR <sub>b</sub> - Ingestion rate for soil =        | mg/day  | 100       | USEPA 1997, EFH       |
| EF <sub>a</sub> - Exposure frequency =               | days/year   | 10        | reasonable assumption |
| EF <sub>b</sub> - Exposure frequency =               | days/year   | 70        | reasonable assumption |
| ED - Exposure duration =                             | years   | 1         | USEPA 1995, Region IV |
| CF - Conversion factor =                             | kg/mg   | 1.00E-06  |                       |
| ME - Matrix effect =                                 |   | 1         | Magee, et al., 1996   |
| BW - Body weight =                                   | kg  | 70        | USEPA 1995, Region IV |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days  | 365       | USEPA 1991, HHEM      |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days  | 25550     | USEPA 1991, HHEM      |

## Exposure Level A

| Constituent            | Concentration in<br>Soil<br>mg/kg | Average Daily       |                          | Oral Chronic RfD<br>mg/kg-day | Hazard<br>Index | Average Lifetime Daily<br>Intake<br>mg/kg-day | Oral Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------------|---------------------|--------------------------|-------------------------------|-----------------|---|--|-------------|
|                        |                                   | Intake<br>mg/kg-day | Chronic RfD<br>mg/kg-day |                               |                 |   |  |             |
| <b>Semivolatiles</b>   |                                   |                     |                          |                               |                 |   |  |             |
| Benzo(a)anthracene     | 6.70E+00                          | 1.26E-06            | NA                       | NA                            | NA              | 1.80E-08                                      | 7.30E-01                                     | 1.31E-08    |
| Benzo(a)pyrene         | 5.20E+00                          | 9.77E-07            | NA                       | NA                            | NA              | 1.40E-08                                      | 7.30E+00                                     | 1.02E-07    |
| Benzo(b)fluoranthene   | 9.20E+00                          | 1.73E-06            | NA                       | NA                            | NA              | 2.47E-08                                      | 7.30E-01                                     | 1.80E-08    |
| Benzo(k)fluoranthene   | 3.60E+00                          | 6.76E-07            | NA                       | NA                            | NA              | 9.66E-09                                      | 7.30E-02                                     | 7.05E-10    |
| Chrysene               | 5.33E+00                          | 1.00E-06            | NA                       | NA                            | NA              | 1.43E-08                                      | 7.30E-03                                     | 1.04E-10    |
| Dibenz(a,h)anthracene  | 2.39E-01                          | 4.49E-08            | NA                       | NA                            | NA              | 6.41E-10                                      | 7.30E+00                                     | 4.68E-09    |
| Indeno(1,2,3-cd)pyrene | 3.70E+00                          | 6.95E-07            | NA                       | NA                            | NA              | 9.93E-09                                      | 7.30E-01                                     | 7.25E-09    |

NA - Not Applicable

Cancer Risk = 1.46E-07

## Exposure Level B

| Constituent            | Concentration in<br>Soil<br>mg/kg | Average Daily       |                          | Oral Chronic RfD<br>mg/kg-day | Hazard<br>Index | Average Lifetime Daily<br>Intake<br>mg/kg-day | Oral Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------------|---------------------|--------------------------|-------------------------------|-----------------|---|--|-------------|
|                        |                                   | Intake<br>mg/kg-day | Chronic RfD<br>mg/kg-day |                               |                 |   |  |             |
| <b>Semivolatiles</b>   |                                   |                     |                          |                               |                 |   |  |             |
| Benzo(a)anthracene     | 6.70E+00                          | 1.84E-06            | NA                       | NA                            | NA              | 2.62E-08                                      | 7.30E-01                                     | 1.91E-08    |
| Benzo(a)pyrene         | 5.20E+00                          | 1.42E-06            | NA                       | NA                            | NA              | 2.04E-08                                      | 7.30E+00                                     | 1.49E-07    |
| Benzo(b)fluoranthene   | 9.20E+00                          | 2.52E-06            | NA                       | NA                            | NA              | 3.60E-08                                      | 7.30E-01                                     | 2.63E-08    |
| Benzo(k)fluoranthene   | 3.60E+00                          | 9.86E-07            | NA                       | NA                            | NA              | 1.41E-08                                      | 7.30E-02                                     | 1.03E-09    |
| Chrysene               | 5.33E+00                          | 1.46E-06            | NA                       | NA                            | NA              | 2.09E-08                                      | 7.30E-03                                     | 1.52E-10    |
| Dibenz(a,h)anthracene  | 2.39E-01                          | 6.55E-08            | NA                       | NA                            | NA              | 9.35E-10                                      | 7.30E+00                                     | 6.83E-09    |
| Indeno(1,2,3-cd)pyrene | 3.70E+00                          | 1.01E-06            | NA                       | NA                            | NA              | 1.45E-08                                      | 7.30E-01                                     | 1.06E-08    |

NA - Not Applicable

Cancer Risk = 2.13E-07

Total Cancer Risk = 3.58E-07



Table 60

*Exposure to Construction Workers from Inhalation of Fugitive Dust in EU2*  
*Kerr McGee, Hattiesburg, MS*

| Intake (mg/kg-day) = $C_a * InhR * EF * ED * RF$ |   | $C_a = \text{Concentration in Air (mg/m}^3\text{)} = E_i / (Hb * W * V)$ |   | $E_i = \text{Emission Rate of Component (mg/sec)} = \text{see below}$ |   |
|--|---|--|---|---|---|
| Chemical   | Concentration in air =<br>mg/m <sup>3</sup> | BW * AT<br>mg/m <sup>3</sup> shift                                       | Concentration in Air<br>in Air<br>mg/m <sup>3</sup> | Average Daily<br>Intake<br>mg/kg-day                                  | Inhalation<br>Subchronic RTD<br>mg/kg-day |
| Benzo(a)anthracene                               | 6.70E+00                                    | 7.33E-03   | 6.50E-06  | 3.05E-07  | NA  |
| Benzo(a)pyrene                                   | 5.20E+00                                    | 5.69E-03   | 5.05E-06  | 2.37E-07  | NA  |
| Benzo(b)fluoranthene                             | 9.20E+00                                    | 1.01E-02   | 8.93E-06  | 4.19E-07  | NA  |
| Benzo(k)fluoranthene                             | 3.60E+00                                    | 3.94E-03   | 3.49E-06  | 1.64E-07  | NA  |
| Chrysene   | 5.33E+00                                    | 5.83E-03   | 5.17E-06  | 2.43E-07  | NA  |
| Dibenz(a,h)anthracene                            | 2.39E-01                                    | 2.61E-04   | 2.32E-07  | 1.09E-08  | NA  |
| Indeno(1,2,3-cd)pyrene                           | 3.70E+00                                    | 4.05E-03   | 3.59E-06  | 1.69E-07  | NA  |
| NA - Not Available                               |   |  |   |   | Total Cancer Risk: 1.37E-08               |

Intake (mg/kg-day) =  $C_a * InhR * EF * ED * RF$   
 $C_a$  - Concentration in air =  
 InhR - Inhalation Rate =  
 EF - Exposure Frequency =  
 ED - Exposure Duration =  
 $RF_s$  - Retention Factor - semivolatiles =  
 $AT_b$  - Averaging Time noncarcinogenic =  
 $AT_c$  - Averaging Time carcinogenic =  
 BW - Body Weight =  
 $E_i$  - Emission Rate (mg/sec) =  $C_s * (PERv + PERe)$   
 $C_s$  - Concentration in soil = mg/kg chem.spec.  
 BW \* AT  
 mg/m<sup>3</sup> shift  
 m'shift/year  
 years  
 0.75  
 days  
 days  
 kg  
 mg/kg chem.spec.  
 USEPA 1995, Region IV  
 reasonable assumption  
 reasonable assumption  
 ICRP, 1968  
 USEPA 1991, HHSEM  
 USEPA 1991, HHSEM  
 USEPA 1995, Region IV  
 $V$  - Wind speed (m/sec) = 4.69  
 Length (downwind distance) (m) = 50  
 $r$  - Roughness Ht. (m) = 0.20  
 $z$  - downwind distance (m) = 50  
 $z = 6.25[r(Hb/r) * Ln(Hb/r) - 1.58 * Hb/r + 1.58]$

Table 61

Dermal Exposure to EU4 Sediment by a Construction Worker

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                                       | <u><math>\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}</math></u> |                                |                 |                                       |   |          |
|---|---------------------------------------|---|--------------------------------|-----------------|---------------------------------------|---|----------|
| Cs - Concentration in sediment =                            | mg/kg                                 | chem. spec.   |                                |                 |                                       |   |          |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                  | 3000  | calculated                     |                 |                                       |   |          |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                       | 20000   | USEPA 1997, EFH                |                 |                                       |   |          |
| Fs - Fraction of skin surface area available for exposure = |                                       | 15.0%   | USEPA 1997, EFH                |                 |                                       |   |          |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                    | 0.13  | USEPA 1997, EFH                |                 |                                       |   |          |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                                       | 0.03  | USEPA 1995, Region III         |                 |                                       |   |          |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |                                       | 0.1   | USEPA 1995, Region III         |                 |                                       |   |          |
| EF - Exposure frequency =                                   | days/year                             | 8   | reasonable assumption          |                 |                                       |   |          |
| ED - Exposure duration =                                    | years                                 | 1   | reasonable assumption          |                 |                                       |   |          |
| CF - Conversion factor =                                    | kg/mg                                 | 1.00E-06  |                                |                 |                                       |   |          |
| BW - Body weight =  | kg                                    | 70  | USEPA 1995, Region IV          |                 |                                       |   |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                                  | 365   | USEPA 1991, HHEM               |                 |                                       |   |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                                  | 25550   | USEPA 1991, HHEM               |                 |                                       |   |          |
| Constituent   | Concentration<br>in Sediment<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day                                  | Dermal                         | Hazard<br>Index | Average                               | Cancer Slope<br>Factor<br>1/(mg/kg-day) |          |
|   |                                       |   | Subchronic<br>RfD<br>mg/kg-day |                 | Lifetime Daily<br>Intake<br>mg/kg-day |   |          |
| Semivolatiles   |                                       |   |                                |                 |                                       |   |          |
| Benzo(a)anthracene  | 3.30E+02                              | 1.21E-06  | NA                             | NA              | 1.73E-08                              | 1.46E+00                                | 2.52E-08 |
| Benzo(a)pyrene  | 1.30E+02                              | 4.76E-07  | NA                             | NA              | 6.80E-09                              | 1.46E+01                                | 9.93E-08 |
| Benzo(b)fluoranthene  | 1.80E+02                              | 6.59E-07  | NA                             | NA              | 9.42E-09                              | 1.46E+00                                | 1.38E-08 |
| Benzo(k)fluoranthene  | 6.40E+01                              | 2.34E-07  | NA                             | NA              | 3.35E-09                              | 1.46E-01                                | 4.89E-10 |
| Carbazole   | 5.90E+02                              | 7.20E-06  | NA                             | NA              | 1.03E-07                              | 2.00E-02                                | 2.06E-09 |
| Chrysene  | 2.90E+02                              | 1.06E-06  | NA                             | NA              | 1.52E-08                              | 1.46E-02                                | 2.22E-10 |
| Dibenz(a,h)anthracene                                       | 1.20E+01                              | 4.40E-08  | NA                             | NA              | 6.28E-10                              | 1.46E+01                                | 9.17E-09 |
| Dibenzofuran  | 9.40E+02                              | 1.15E-05  | NA                             | NA              | 1.64E-07                              | NA                                      | NA       |
| Indeno(1,2,3-cd)pyrene                                      | 4.70E+01                              | 1.72E-07  | NA                             | NA              | 2.46E-09                              | 1.46E+00                                | 3.59E-09 |
| Naphthalene   | 3.00E+03                              | 3.66E-05  | NA                             | NA              | 5.23E-07                              | NA                                      | NA       |
| Phenanthrene  | 3.20E+03                              | 3.91E-05  | NA                             | NA              | 5.58E-07                              | NA                                      | NA       |

NA - Not Available

Total Cancer Risk = 1.54E-07



**Table 62**  
**Oral Exposure to EU4 Sediment by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |                                 | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                               |              |   |  |          |
|--|---------------------------------|---|-------------------------------|--------------|---|--|----------|
| Cd - Concentration in sediment =                     | mg/kg                           | see below   |                               |              |   |  |          |
| IngR - Ingestion rate for sediment =                 | mg/day                          | 480   |                               |              |   | USEPA 1997, EFH                        |          |
| EF - Exposure frequency =                            | days/year                       | 8   |                               |              |   | reasonable assumption                  |          |
| ED - Exposure duration =                             | years                           | 1   |                               |              |   | USEPA 1995, Region IV                  |          |
| CF - Conversion factor =                             | kg/mg                           | 1.00E-06  |                               |              |   |  |          |
| ME - Matrix effect =                                 |                                 | 1   |                               |              |   | Magee, et al., 1996                    |          |
| BW - Body weight =                                   | kg                              | 70  |                               |              |   | USEPA 1995, Region IV                  |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                            | 365   |                               |              |   | USEPA 1991, HHEM                       |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                            | 25550   |                               |              |   | USEPA 1991, HHEM                       |          |
| Constituent  | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day  | Oral Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) |          |
| <b>Semivolatiles</b>                                 |                                 |   |                               |              |   | Cancer Risk                            |          |
| Benzo(a)anthracene                                   | 3.30E+02                        | 4.96E-05  | NA                            | NA           | 7.09E-07                                | 7.30E-01                               | 5.17E-07 |
| Benzo(a)pyrene                                       | 1.30E+02                        | 1.95E-05  | NA                            | NA           | 2.79E-07                                | 7.30E+00                               | 2.04E-06 |
| Benzo(b)fluoranthene                                 | 1.80E+02                        | 2.71E-05  | NA                            | NA           | 3.86E-07                                | 7.30E-01                               | 2.82E-07 |
| Benzo(k)fluoranthene                                 | 6.40E+01                        | 9.62E-06  | NA                            | NA           | 1.37E-07                                | 7.30E-02                               | 1.00E-08 |
| Carbazole  | 5.90E+02                        | 8.87E-05  | NA                            | NA           | 1.27E-06                                | 2.00E-02                               | 2.53E-08 |
| Chrysene   | 2.90E+02                        | 4.36E-05  | NA                            | NA           | 6.23E-07                                | 7.30E-03                               | 4.55E-09 |
| Dibenz(a,h)anthracene                                | 1.20E+01                        | 1.80E-06  | NA                            | NA           | 2.58E-08                                | 7.30E+00                               | 1.88E-07 |
| Dibenzofuran   | 9.40E+02                        | 1.41E-04  | NA                            | NA           | 2.02E-06                                | NA                                     | NA       |
| Indeno(1,2,3-cd)pyrene                               | 4.70E+01                        | 7.06E-06  | NA                            | NA           | 1.01E-07                                | 7.30E-01                               | 7.37E-08 |
| Naphthalene  | 3.00E+03                        | 4.51E-04  | NA                            | NA           | 6.44E-06                                | NA                                     | NA       |
| Phenanthrene   | 3.20E+03                        | 4.81E-04  | NA                            | NA           | 6.87E-06                                | NA                                     | NA       |

NA - Not Applicable

Total Cancer Risk = 3.14E-06



Table 63

Dermal Exposure to EU4 Surface Water by a Construction Worker

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                   | $C_w * SA * K_p * ABS * ET * EF * ED * CF$ |           |  |  |  |  |
|---|-------------------|--|-----------|--|--|--|--|
|   |                   | BW * AT                                    |           |  |  |  |  |
| Cw - Concentration in surface water =                       | mg/L              |  | see below |  |  |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup>   | 3000                                       |           | calculated                             |  |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>   | 20000                                      |           | USEPA 1997, EFH                        |  |  |  |
| Fs - Fraction of skin surface area available for exposure = |                   | 15.0%                                      |           | USEPA 1997, EFH                        |  |  |  |
| Kp - Dermal permeability constant =                         | cm/hr             | see below                                  |           |  |  |  |  |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                   | 0.03                                       |           | USEPA 1995, Region III                 |  |  |  |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |                   | 0.1  |           | USEPA 1995, Region III                 |  |  |  |
| ET - Exposure time =  | hrs/day           | 1  |           | USEPA 1992, Dermal Exposure Assessment |  |  |  |
| EF - Exposure frequency =                                   | days/year         | 8  |           | reasonable assumption                  |  |  |  |
| ED - Exposure duration =                                    | years             | 1  |           | reasonable assumption                  |  |  |  |
| CF - Conversion factor =                                    | L/cm <sup>2</sup> | 1.00E-03                                   |           |  |  |  |  |
| BW - Body weight =  | kg                | 70   |           | USEPA 1995, Region IV                  |  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days              | 365  |           | USEPA 1991, HHEM                       |  |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days              | 25550                                      |           | USEPA 1991, HHEM                       |  |  |  |

| Constituent                | Concentration in Surface Water mg/L | Kp cm/hr | Average Daily Intake mg/kg-day | Dermal RfD mg/kg-day |          | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|----------------------------|-------------------------------------|----------|--------------------------------|----------------------|----------|--------------|---|-----------------------------------|-------------|
|                            |                                     |          |                                | Subchronic           | Chronic  |              |   |                                   |             |
| <b>Semivolatiles</b>       |                                     |          |                                |                      |          |              |   |                                   |             |
| Benzo(a)anthracene         | 5.00E-03                            | 8.10E-01 | 1.14E-07                       | NA                   | NA       | NA           | 1.63E-09                                | 1.46E+00                          | 2.38E-09    |
| Benzo(a)pyrene             | 5.00E-04                            | 1.20E+00 | 1.69E-08                       | NA                   | NA       | NA           | 2.42E-10                                | 1.46E+01                          | 3.53E-09    |
| Benzo(b)fluoranthene       | 1.20E-02                            | 1.20E+00 | 4.06E-07                       | NA                   | NA       | NA           | 5.80E-09                                | 1.46E+00                          | 8.46E-09    |
| Benzo(k)fluoranthene       | 2.00E-03                            | 4.48E+01 | 2.52E-06                       | NA                   | NA       | NA           | 3.60E-08                                | 1.46E-01                          | 5.26E-09    |
| Bis(2-ethylhexyl)phthalate | 3.00E-03                            | 3.30E-02 | 9.30E-09                       | 1.00E-02             | 9.30E-07 | 1.33E-10     | 1.40E-02                                | 1.86E-12                          |             |
| Carbazole                  | 1.00E-02                            | 3.57E-02 | 3.36E-08                       | NA                   | NA       | NA           | 4.80E-10                                | 2.00E-02                          | 9.59E-12    |
| Chrysene                   | 6.00E-03                            | 8.10E-01 | 1.37E-07                       | NA                   | NA       | NA           | 1.96E-09                                | 1.46E-02                          | 2.86E-11    |
| Dibenz(a,h)anthracene      | 5.00E-04                            | 2.70E+00 | 3.80E-08                       | NA                   | NA       | NA           | 5.43E-10                                | 1.46E+01                          | 7.93E-09    |
| Dibenzofuran               | 1.10E-02                            | 1.51E-01 | 1.56E-07                       | NA                   | NA       | NA           | 2.23E-09                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene     | 5.00E-04                            | 1.90E+00 | 2.68E-08                       | NA                   | NA       | NA           | 3.82E-10                                | 1.46E+00                          | 5.58E-10    |
| Phenanthrene               | 1.70E-02                            | 2.30E-01 | 3.67E-07                       | NA                   | NA       | NA           | 5.25E-09                                | NA                                | NA          |

NA - Not Available

Total Hazard Index = 9.30E-07

Total Cancer Risk = 2.82E-08



**Table 64**  
**Oral Exposure to EU4 Surface Water by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                                       | $\frac{C_{sw} \cdot I_{ngR} \cdot E_F \cdot E_D \cdot E_T}{B_W \cdot A_T}$ |                            |                     |                                      |                                 |          |
|---|---------------------------------------|--|----------------------------|---------------------|--------------------------------------|---------------------------------|----------|
| C <sub>sw</sub> - Concentration in surface water =              |                                       | mg/L   |                            |                     |                                      |                                 |          |
| I <sub>ngR</sub> - Ingestion rate for surface water =           |                                       | L/hour   |                            |                     |                                      |                                 |          |
| E <sub>F</sub> - Exposure frequency =                           |                                       | days/year  |                            |                     |                                      |                                 |          |
| E <sub>D</sub> - Exposure duration =                            |                                       | years  |                            |                     |                                      |                                 |          |
| E <sub>T</sub> - Exposure time =                                |                                       | hrs/day  |                            |                     |                                      |                                 |          |
| B <sub>W</sub> - Body weight =                                  |                                       | kg   |                            |                     |                                      |                                 |          |
| A <sub>T<sub>n</sub></sub> - Averaging time - noncarcinogenic = |                                       | days   |                            |                     |                                      |                                 |          |
| A <sub>T<sub>c</sub></sub> - Averaging time - carcinogenic =    |                                       | days   |                            |                     |                                      |                                 |          |
|   |                                       |  |                            |                     |                                      |                                 |          |
| <b>Constituent</b>  | <b>Concentration in Surface Water</b> | <b>Average Daily Intake</b>  | <b>Oral Subchronic RfD</b> | <b>Hazard Index</b> | <b>Average Lifetime Daily Intake</b> | <b>Oral Cancer Slope Factor</b> |          |
|   | mg/L                                  | mg/kg-day  | mg/kg-day                  |                     | mg/kg-day                            | 1/(mg/kg-day)                   |          |
| Semivolatiles   |                                       |  |                            |                     |                                      |                                 |          |
| Benzo(a)anthracene  | 5.00E-03                              | 1.57E-08   | NA                         | NA                  | 2.24E-10                             | 7.30E-01                        | 1.63E-10 |
| Benzo(a)pyrene  | 5.00E-04                              | 1.57E-09   | NA                         | NA                  | 2.24E-11                             | 7.30E+00                        | 1.63E-10 |
| Benzo(b)fluoranthene  | 1.20E-02                              | 3.76E-08   | NA                         | NA                  | 5.37E-10                             | 7.30E-01                        | 3.92E-10 |
| Benzo(k)fluoranthene  | 2.00E-03                              | 6.26E-09   | NA                         | NA                  | 8.95E-11                             | 7.30E-02                        | 6.53E-12 |
| Bis(2-ethylhexyl)phthalate                                      | 3.00E-03                              | 9.39E-09   | 2.00E-02                   | 4.70E-07            | 1.34E-10                             | 1.40E-02                        | 1.88E-12 |
| Carbazole   | 1.00E-02                              | 3.13E-08   | NA                         | NA                  | 4.47E-10                             | 2.00E-02                        | 8.95E-12 |
| Chrysene  | 6.00E-03                              | 1.88E-08   | NA                         | NA                  | 2.68E-10                             | 7.30E-03                        | 1.96E-12 |
| Dibenz(a,h)anthracene   | 5.00E-04                              | 1.57E-09   | NA                         | NA                  | 2.24E-11                             | 7.30E+00                        | 1.63E-10 |
| Dibenzofuran  | 1.10E-02                              | 3.44E-08   | NA                         | NA                  | 4.92E-10                             | NA                              | NA       |
| Indeno(1,2,3-cd)pyrene  | 5.00E-04                              | 1.57E-09   | NA                         | NA                  | 2.24E-11                             | 7.30E-01                        | 1.63E-11 |
| Phenanthrene  | 1.70E-02                              | 5.32E-08   | NA                         | NA                  | 7.60E-10                             | NA                              | NA       |

NA - Not Applicable

Total Hazard Index = 4.70E-07

Total Cancer Risk = 9.17E-10



**Table 65**

*Dermal Exposure to EU4 Soil (0-20') by a Construction Worker  
Kerr McGee, Hattiesburg, MS*

| Intake (mg/kg-day) =  | $\frac{\text{Cs} * \text{SA} * \text{AH} * \text{ABS} * \text{EF} * \text{ED} * \text{CF}}{\text{BW} * \text{AT}}$ |                                |                                 |              |   |                                   |             |
|---|--|--------------------------------|---------------------------------|--------------|---|-----------------------------------|-------------|
| Cs - Concentration in soil =                                | mg/kg  | chem. spec.                    |                                 |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day   | 5560                           | calculated                      |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>  | 20000                          | USEPA 1997, EFH                 |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |  | 27.8%                          | USEPA 1997, EFH                 |              |   |                                   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.1                            | USEPA 1997, EFH                 |              |   |                                   |             |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |  | 0.03                           | USEPA 1995, Region III          |              |   |                                   |             |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |  | 0.1                            | USEPA 1995, Region III          |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year  | 80                             | reasonable assumption           |              |   |                                   |             |
| ED - Exposure duration =                                    | years  | 1                              | reasonable assumption           |              |   |                                   |             |
| CF - Conversion factor =                                    | kg/mg  | 1.00E-06                       |                                 |              |   |                                   |             |
| BW - Body weight =  | kg   | 70                             | USEPA 1995, Region IV           |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days   | 365                            | USEPA 1991, HHEM                |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days   | 25550                          | USEPA 1991, HHEM                |              |   |                                   |             |
|   |  |                                |                                 |              |   |                                   |             |
| Constituent   | Concentration in Soil mg/kg  | Average Daily Intake mg/kg-day | Dermal Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |  |                                |                                 |              |   |                                   |             |
| Benzo(a)anthracene  | 9.30E+02   | 4.86E-05                       | NA                              | NA           | 6.94E-07                                | 1.46E+00                          | 1.01E-06    |
| Benzo(a)pyrene  | 5.00E+02   | 2.61E-05                       | NA                              | NA           | 3.73E-07                                | 1.46E+01                          | 5.45E-06    |
| Benzo(b)fluoranthene  | 5.30E+02   | 2.77E-05                       | NA                              | NA           | 3.95E-07                                | 1.46E+00                          | 5.77E-07    |
| Benzo(k)fluoranthene  | 2.90E+02   | 1.51E-05                       | NA                              | NA           | 2.16E-07                                | 1.46E-01                          | 3.16E-08    |
| Carbazole   | 6.20E+02   | 1.08E-04                       | NA                              | NA           | 1.54E-06                                | 2.00E-02                          | 3.08E-08    |
| Chrysene  | 6.90E+02   | 3.60E-05                       | NA                              | NA           | 5.15E-07                                | 1.46E-02                          | 7.52E-09    |
| Dibenz(a,h)anthracene                                       | 6.40E+01   | 3.34E-06                       | NA                              | NA           | 4.78E-08                                | 1.46E+01                          | 6.97E-07    |
| Indeno(1,2,3-cd)pyrene                                      | 2.50E+02   | 1.31E-05                       | NA                              | NA           | 1.87E-07                                | 1.46E+00                          | 2.72E-07    |
| Naphthalene   | 3.50E+03   | 6.09E-04                       | NA                              | NA           | 8.70E-06                                | NA                                | NA          |

NA - Not Available

Total Cancer Risk = 8.08E-06



**Table 66**  
**Oral Exposure to EU4 Soil (0-20') by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 | <u><math>\frac{\text{Cd} \cdot \text{IngR} \cdot \text{EF} \cdot \text{ED} \cdot \text{CF} \cdot \text{ME}}{\text{BW} \cdot \text{AT}}</math></u> |           |                       |
|--|---|-----------|-----------------------|
| Cd - Concentration in soil =                         | mg/kg   | see below |                       |
| IngR <sub>a</sub> - Ingestion rate for soil =        | mg/day  | 480       | USEPA 1997, EFH       |
| IngR <sub>b</sub> - Ingestion rate for soil =        | mg/day  | 100       | USEPA 1997, EFH       |
| EF <sub>a</sub> - Exposure frequency =               | days/year   | 10        | reasonable assumption |
| EF <sub>b</sub> - Exposure frequency =               | days/year   | 70        | reasonable assumption |
| ED - Exposure duration =                             | years   | 1         | USEPA 1995, Region IV |
| CF - Conversion factor =                             | kg/mg   | 1.00E-06  |                       |
| ME - Matrix effect =                                 |   | 1         | Magee, et al., 1996   |
| BW - Body weight =                                   | kg  | 70        | USEPA 1995, Region IV |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days  | 365       | USEPA 1991, HHEM      |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days  | 25550     | USEPA 1991, HHEM      |

**Exposure Level A**

| Constituent            | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day | Oral Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------|--------------------------------|-------------------------------|--------------|---|--|-------------|
| <b>Semivolatiles</b>   |                             |                                |                               |              |   |  |             |
| Benzo(a)anthracene     | 9.30E+02                    | 1.75E-04                       | NA                            | NA           | 2.50E-06                                | 7.30E-01                               | 1.82E-06    |
| Benzo(a)pyrene         | 5.00E+02                    | 9.39E-05                       | NA                            | NA           | 1.34E-06                                | 7.30E+00                               | 9.80E-06    |
| Benzo(b)fluoranthene   | 5.30E+02                    | 9.96E-05                       | NA                            | NA           | 1.42E-06                                | 7.30E-01                               | 1.04E-06    |
| Benzo(k)fluoranthene   | 2.90E+02                    | 5.45E-05                       | NA                            | NA           | 7.78E-07                                | 7.30E-02                               | 5.68E-08    |
| Carbazole              | 6.20E+02                    | 1.16E-04                       | NA                            | NA           | 1.66E-06                                | 2.00E-02                               | 3.33E-08    |
| Chrysene               | 6.90E+02                    | 1.30E-04                       | NA                            | NA           | 1.85E-06                                | 7.30E-03                               | 1.35E-08    |
| Dibenz(a,h)anthracene  | 6.40E+01                    | 1.20E-05                       | NA                            | NA           | 1.72E-07                                | 7.30E+00                               | 1.25E-06    |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                    | 4.70E-05                       | NA                            | NA           | 6.71E-07                                | 7.30E-01                               | 4.90E-07    |
| Naphthalene            | 3.50E+03                    | 6.58E-04                       | NA                            | NA           | 9.39E-06                                | NA                                     | NA          |

NA - Not Applicable

Cancer Risk = 1.45E-05

**Exposure Level B**

| Constituent            | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day | Oral Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------|--------------------------------|-------------------------------|--------------|---|--|-------------|
| <b>Semivolatiles</b>   |                             |                                |                               |              |   |  |             |
| Benzo(a)anthracene     | 9.30E+02                    | 2.55E-04                       | NA                            | NA           | 3.64E-06                                | 7.30E-01                               | 2.66E-06    |
| Benzo(a)pyrene         | 5.00E+02                    | 1.37E-04                       | NA                            | NA           | 1.96E-06                                | 7.30E+00                               | 1.43E-05    |
| Benzo(b)fluoranthene   | 5.30E+02                    | 1.45E-04                       | NA                            | NA           | 2.07E-06                                | 7.30E-01                               | 1.51E-06    |
| Benzo(k)fluoranthene   | 2.90E+02                    | 7.95E-05                       | NA                            | NA           | 1.14E-06                                | 7.30E-02                               | 8.29E-08    |
| Carbazole              | 6.20E+02                    | 1.70E-04                       | NA                            | NA           | 2.43E-06                                | 2.00E-02                               | 4.85E-08    |
| Chrysene               | 6.90E+02                    | 1.89E-04                       | NA                            | NA           | 2.70E-06                                | 7.30E-03                               | 1.97E-08    |
| Dibenz(a,h)anthracene  | 6.40E+01                    | 1.75E-05                       | NA                            | NA           | 2.50E-07                                | 7.30E+00                               | 1.83E-06    |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                    | 6.85E-05                       | NA                            | NA           | 9.78E-07                                | 7.30E-01                               | 7.14E-07    |
| Naphthalene            | 3.50E+03                    | 9.59E-04                       | NA                            | NA           | 1.37E-05                                | NA                                     | NA          |

NA - Not Applicable

Cancer Risk = 2.12E-05

Total Cancer Risk = 3.57E-05

*Table 67  
Exposure to Construction Workers from Inhalation of Fugitive Dust in EU4  
Kerr McGee, Hattiesburg, MS*

NA - Not Available

**Table 68**  
**Dermal Exposure to EU5 Soil (0-20') by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                                   | $\frac{\text{Cs} * \text{SA} * \text{AH} * \text{ABS} * \text{EF} * \text{ED} * \text{CF}}{\text{BW} * \text{AT}}$ |  |                                   |  |   |                |
|---|-----------------------------------|--|--|-----------------------------------|--|---|----------------|
| Cs - Concentration in soil =                                | mg/kg                             | chem. spec.  |  |                                   |  |   |                |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day              | 5560   | calculated                               |                                   |  |   |                |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                   | 20000  | USEPA 1997, EFH                          |                                   |  |   |                |
| Fs - Fraction of skin surface area available for exposure = |                                   | 27.8%  | USEPA 1997, EFH                          |                                   |  |   |                |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                | 0.1  | USEPA 1997, EFH                          |                                   |  |   |                |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                                   | 0.03   | USEPA 1995, Region III                   |                                   |  |   |                |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |                                   | 0.1  | USEPA 1995, Region III                   |                                   |  |   |                |
| EF - Exposure frequency =                                   | days/year                         | 80   | reasonable assumption                    |                                   |  |   |                |
| ED - Exposure duration =                                    | years                             | 1  | reasonable assumption                    |                                   |  |   |                |
| CF - Conversion factor =                                    | kg/mg                             | 1.00E-06   |  |                                   |  |   |                |
| BW - Body weight =  | kg                                | 70   | USEPA 1995, Region IV                    |                                   |  |   |                |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                              | 365  | USEPA 1991, HHEM                         |                                   |  |   |                |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                              | 25550  | USEPA 1991, HHEM                         |                                   |  |   |                |
| Constituent   | Concentration in<br>Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day   | Dermal<br>Subchronic<br>RfD<br>mg/kg-day | Lifetime Daily<br>Hazard<br>Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
| Semivolatiles   |                                   |  |  |                                   |  |   |                |
| Benzo(a)anthracene  | 8.35E+01                          | 4.36E-06   | NA                                       | NA                                | 6.23E-08   | 1.46E+00                                | 9.10E-08       |
| Benzo(a)pyrene  | 4.42E+01                          | 2.31E-06   | NA                                       | NA                                | 3.30E-08   | 1.46E+01                                | 4.81E-07       |
| Benzo(b)fluoranthene  | 7.95E+01                          | 4.15E-06   | NA                                       | NA                                | 5.93E-08   | 1.46E+00                                | 8.66E-08       |
| Benzo(k)fluoranthene  | 1.68E+01                          | 8.77E-07   | NA                                       | NA                                | 1.25E-08   | 1.46E-01                                | 1.83E-09       |
| Chrysene  | 8.25E+01                          | 4.31E-06   | NA                                       | NA                                | 6.16E-08   | 1.46E-02                                | 8.99E-10       |
| Dibenz(a,h)anthracene                                       | 1.53E+00                          | 7.99E-08   | NA                                       | NA                                | 1.14E-09   | 1.46E+01                                | 1.67E-08       |
| Fluorene  | 2.24E+02                          | 3.90E-05   | 2.00E-01                                 | 1.95E-04                          | 5.57E-07   | NA                                      | NA             |
| Indeno(1,2,3-cd)pyrene                                      | 1.32E+01                          | 6.89E-07   | NA                                       | NA                                | 9.85E-09   | 1.46E+00                                | 1.44E-08       |

NA - Not Available

Total Hazard Index = 1.95E-04

Total Cancer Risk = 6.93E-07



**Table 69****Oral Exposure to EU5 Soil (0-20') by a Construction Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 | <u>Cd*IngR*EF*ED*CF*ME</u><br>BW*AT |           |                       |
|--|-------------------------------------|-----------|-----------------------|
| Cd - Concentration in sediment =                     | mg/kg                               | see below |                       |
| IngR <sub>a</sub> - Ingestion rate for soil =        | mg/day                              | 480       | USEPA 1997, EFH       |
| IngR <sub>b</sub> - Ingestion rate for soil =        | mg/day                              | 100       | USEPA 1997, EFH       |
| EF <sub>a</sub> - Exposure frequency =               | days/year                           | 10        | reasonable assumption |
| EF <sub>b</sub> - Exposure frequency =               | days/year                           | 70        | reasonable assumption |
| ED - Exposure duration =                             | years                               | 1         | USEPA 1995, Region IV |
| CF - Conversion factor =                             | kg/mg                               | 1.00E-06  |                       |
| ME - Matrix effect =                                 |                                     | 1         | Magee, et al., 1996   |
| BW - Body weight =                                   | kg                                  | 70        | USEPA 1995, Region IV |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                                | 365       | USEPA 1991, HHEM      |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                                | 25550     | USEPA 1991, HHEM      |

**Exposure Level A**

| Constituent            | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average Lifetime<br>Daily Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------------|--------------------------------------|----------------------------------|-----------------|---|--|-------------|
| <b>Semivolatiles</b>   |                                   |                                      |                                  |                 |   |  |             |
| Benzo(a)anthracene     | 8.35E+01                          | 1.57E-05                             | NA                               | NA              | 2.24E-07                                      | 7.30E-01                                     | 1.64E-07    |
| Benzo(a)pyrene         | 4.42E+01                          | 8.30E-06                             | NA                               | NA              | 1.19E-07                                      | 7.30E+00                                     | 8.66E-07    |
| Benzo(b)fluoranthene   | 7.95E+01                          | 1.49E-05                             | NA                               | NA              | 2.13E-07                                      | 7.30E-01                                     | 1.56E-07    |
| Benzo(k)fluoranthene   | 1.68E+01                          | 3.16E-06                             | NA                               | NA              | 4.51E-08                                      | 7.30E-02                                     | 3.29E-09    |
| Chrysene               | 8.25E+01                          | 1.55E-05                             | NA                               | NA              | 2.21E-07                                      | 7.30E-03                                     | 1.62E-09    |
| Dibenz(a,h)anthracene  | 1.53E+00                          | 2.87E-07                             | NA                               | NA              | 4.11E-09                                      | 7.30E+00                                     | 3.00E-08    |
| Fluorene               | 2.24E+02                          | 4.21E-05                             | 4.00E-02                         | 1.05E-03        | 6.01E-07                                      | NA   | NA          |
| Indeno(1,2,3-cd)pyrene | 1.32E+01                          | 2.48E-06                             | NA                               | NA              | 3.54E-08                                      | 7.30E-01                                     | 2.59E-08    |

NA - Not Applicable

Hazard Index = 1.05E-03

Cancer Risk = 1.25E-06

**Exposure Level B**

| Constituent            | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average Lifetime<br>Daily Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------------|--------------------------------------|----------------------------------|-----------------|---|--|-------------|
| <b>Semivolatiles</b>   |                                   |                                      |                                  |                 |   |  |             |
| Benzo(a)anthracene     | 8.35E+01                          | 2.29E-05                             | NA                               | NA              | 3.27E-07                                      | 7.30E-01                                     | 2.39E-07    |
| Benzo(a)pyrene         | 4.42E+01                          | 1.21E-05                             | NA                               | NA              | 1.73E-07                                      | 7.30E+00                                     | 1.26E-06    |
| Benzo(b)fluoranthene   | 7.95E+01                          | 2.18E-05                             | NA                               | NA              | 3.11E-07                                      | 7.30E-01                                     | 2.27E-07    |
| Benzo(k)fluoranthene   | 1.68E+01                          | 4.60E-06                             | NA                               | NA              | 6.58E-08                                      | 7.30E-02                                     | 4.80E-09    |
| Chrysene               | 8.25E+01                          | 2.26E-05                             | NA                               | NA              | 3.23E-07                                      | 7.30E-03                                     | 2.36E-09    |
| Dibenz(a,h)anthracene  | 1.53E+00                          | 4.19E-07                             | NA                               | NA              | 5.99E-09                                      | 7.30E+00                                     | 4.37E-08    |
| Fluorene               | 2.24E+02                          | 6.14E-05                             | 4.00E-02                         | 1.53E-03        | 8.77E-07                                      | NA   | NA          |
| Indeno(1,2,3-cd)pyrene | 1.32E+01                          | 3.62E-06                             | NA                               | NA              | 5.17E-08                                      | 7.30E-01                                     | 3.77E-08    |

NA - Not Applicable

Hazard Index = 1.53E-03

Cancer Risk = 1.82E-06

Total Hazard Index = 2.59E-03

Total Cancer Risk = 3.06E-06



Table 70

*Exposure to Construction Workers from Inhalation of Fugitive Dust in EU5  
Kerr McGee, Hattiesburg, MS*

| Intake (mg/kg-day) =  | Ca*inhR*EF*ED*RF<br>BW*AT   | Ca = Concentration in Air (mg/m <sup>3</sup> ) = Ei / (Hb * W * V)   |   |  |             |          |                    |          |
|---|---|--|---|--|-------------|----------|--------------------|----------|
| Ca - Concentration in air =<br>InhR - Inhalation Rate =<br>EF - Exposure Frequency =<br>ED - Exposure Duration =<br>RF <sub>s</sub> - Retention Factor - semivolatiles =<br>AT <sub>n</sub> - Average Time noncarcinogenic =<br>AT <sub>c</sub> - Averaging Time carcinogenic =<br>BW - Body Weight = | mg/m <sup>3</sup><br>m <sup>3</sup> /shift<br>20<br>shifts/year<br>80<br>years<br>1<br>0.75<br>days<br>365<br>days<br>25550<br>kg<br>70 | see below<br>USEPA 1995, Region IV<br>reasonable assumption<br>reasonable assumption<br>ICRP, 1968<br>USEPA 1991, HHEM<br>USEPA 1991, HHEM<br>USEPA 1995, Region IV<br>$\tau = 6.25t[Hb/\tau + \ln(Hb/\tau) + 1.58]$ |   |  |             |          |                    |          |
| E <sub>i</sub> - Emission Rate (mg/sec) = Cs*(PERv+PERe)  |   | E <sub>i</sub> - Emission Rate of Component (mg/sec) = see below   |   |  |             |          |                    |          |
| Cs - Concentration in soil =<br>mg/kg   | mg/kg<br>see below  | Hb - Downwind Ht (m) = 4.81<br>W - Width (m) = 50<br>V - Wind speed (m/sec) = 4.69<br>Length (downwind distance) (m) = 50<br>r - Roughness Ht. (m) = 0.20<br>z - downwind distance (m) = 50                          |   |  |             |          |                    |          |
| Concentration in<br>Soil<br>mg/kg   | Emission<br>Rate<br>mg/sec  | Concentration in<br>Air<br>mg/m <sup>3</sup>   |   |  |             |          |                    |          |
| Average Daily<br>Intake<br>mg/kg-day  | Inhalation<br>Subchronic RD<br>mg/kg-day  | Hazard<br>Index  | Average Lifetime<br>Daily Intake<br>mg/kg-day | Inhalation Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |          |                    |          |
| Chemicals   | Semivolatiles   |  |   |  |             |          |                    |          |
| Benzof[a]anthracene   | 8.35E+01  | 9.13E-02   | 8.0E-05                                       | 3.81E-06   | NA          | 5.44E-08 | 3.10E-01           | 1.69E-08 |
| Benzof[aj]pyrene  | 4.42E+01  | 4.83E-02   | 4.29E-05                                      | 2.01E-06   | NA          | 2.88E-08 | 3.10E+00           | 8.92E-08 |
| Benzof[b]fluoranthene   | 7.95E+01  | 8.69E-02   | 7.71E-05                                      | 3.62E-06   | NA          | 5.18E-08 | 3.10E-01           | 1.60E-08 |
| Benzof[k]fluoranthene   | 1.68E+01  | 1.84E-02   | 1.63E-05                                      | 7.66E-07   | NA          | 1.09E-08 | 3.10E-02           | 3.39E-10 |
| Chrysene  | 8.25E+01  | 9.02E-02   | 8.01E-05                                      | 3.76E-06   | NA          | 5.37E-08 | 3.10E-03           | 1.67E-10 |
| Dibenz(a,h)anthracene   | 1.53E+00  | 1.67E-03   | 1.48E-06                                      | 6.97E-08   | NA          | 9.96E-10 | 3.10E+00           | 3.09E-09 |
| Fluorene  | 2.24E+02  | 2.45E-01   | 2.17E-04                                      | 1.02E-05   | NA          | 1.46E-07 | NA                 | NA       |
| Indeno(1,2,3-cd)pyrene  | 1.32E+01  | 1.44E-02   | 1.28E-05                                      | 6.02E-07   | NA          | 8.59E-09 | 3.10E-01           | 2.66E-09 |
| NA - Not Available  |   |  |   |  |             |          | Total Cancer Risk: | 1.28E-07 |

**Table 71****Dermal Exposure to EU6 Sediment by a Child Resident (Aged 1 to 6 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |                        |                               |          |
|---|----------------------|------------------------------|------------------------|-------------------------------|----------|
|   |                      | BW*AT                        |                        |                               |          |
| Cs - Concentration in sediment =                            | mg/kg                | chem. spec.                  |                        |                               |          |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 2229                         | calculated             |                               |          |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 7213                         | USEPA 1997, EFH        |                               |          |
| Fs - Fraction of skin surface area available for exposure = |                      | 30.9%                        | USEPA 1997, EFH        |                               |          |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.33                         | USEPA 1997, EFH        |                               |          |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                      | 0.03                         | USEPA 1995, Region III |                               |          |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |                      | 0.1                          | USEPA 1995, Region III |                               |          |
| EF - Exposure frequency =                                   | days/year            | 40                           | reasonable assumption  |                               |          |
| ED - Exposure duration =                                    | years                | 6                            | USEPA 1995, Region IV  |                               |          |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06                     |                        |                               |          |
| BW - Body weight =  | kg                   | 15                           | USEPA 1995, Region IV  |                               |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 2190                         | USEPA 1991, HHEM       |                               |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550                        | USEPA 1991, HHEM       |                               |          |
| Constituent   |                      | Dermal                       |                        | Average Lifetime Daily Intake |          |
| Concentration in Sediment                                   |                      | Subchronic RID               |                        | Factor                        |          |
| mg/kg   |                      | mg/kg-day                    |                        | 1/(mg/kg-day)                 |          |
| Constituent   |                      | Hazard Index                 |                        | Cancer Risk                   |          |
| Semivolatiles   |                      |                              |                        |                               |          |
| 2-Nitroaniline  | 4.00E-01             | 2.15E-07                     | NA                     | NA                            | NA       |
| 2-Nitrophenol   | 8.00E-01             | 4.30E-07                     | NA                     | NA                            | NA       |
| 3-Nitroaniline  | 8.00E-01             | 4.30E-07                     | NA                     | NA                            | NA       |
| 4-Bromophenylphenylether                                    | 8.00E-01             | 4.30E-07                     | NA                     | NA                            | NA       |
| 4-Chloro-3-methylphenol                                     | 8.00E-01             | 4.30E-07                     | NA                     | NA                            | NA       |
| 4-Chlorophenylphenylether                                   | 4.00E-01             | 2.15E-07                     | NA                     | NA                            | NA       |
| 4-Nitroaniline  | 8.00E-01             | 4.30E-07                     | NA                     | NA                            | NA       |
| Benzo(a)anthracene  | 1.00E+02             | 1.61E-05                     | NA                     | NA                            | 1.46E+00 |
| Benzo(a)pyrene  | 4.90E+01             | 7.90E-06                     | NA                     | NA                            | 1.46E+01 |
| Benzo(b)fluoranthene  | 7.80E+01             | 1.26E-05                     | NA                     | NA                            | 1.46E+00 |
| Benzo(k)fluoranthene  | 2.30E+01             | 3.71E-06                     | NA                     | NA                            | 1.46E+01 |
| Bis(2-chloroethoxy)methane                                  | 8.00E-01             | 4.30E-07                     | NA                     | NA                            | 3.68E-08 |
| Bis(2-chloroethyl)ether                                     | 4.00E-01             | 2.15E-07                     | NA                     | NA                            | 1.84E-08 |
| Carbazole   | 1.00E+02             | 5.37E-05                     | NA                     | NA                            | 4.61E-06 |
| Chrysene  | 7.60E+01             | 1.23E-05                     | NA                     | NA                            | 1.05E-06 |
| Dibenz(a,h)anthracene                                       | 9.60E+00             | 1.55E-06                     | NA                     | NA                            | 1.33E-07 |
| Hexachlorobenzene   | 4.00E-01             | 2.15E-07                     | NA                     | NA                            | 1.84E-08 |
| Hexachlorocyclopentadiene                                   | 2.00E+00             | 1.07E-06                     | NA                     | NA                            | 9.21E-08 |
| Indeno(1,2,3-cd)pyrene                                      | 3.90E+01             | 6.29E-06                     | NA                     | NA                            | 5.39E-07 |
| N-nitrosodi-n-propylamine                                   | 4.00E-01             | 2.15E-07                     | NA                     | NA                            | 1.84E-08 |

NA - Not Available

Total Cancer Risk = 1.65E-05



**Table 72****Oral Exposure to EU6 Sediment by a Child Resident (Aged 1 to 6 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u>Cd*IngR*EF*ED*CF*ME</u><br>BW*AT |                       |  |  |  |
|--|-----------|-------------------------------------|-----------------------|--|--|--|
| Cd - Concentration in sediment =                     | mg/kg     | see below                           |                       |  |  |  |
| IngR - Ingestion rate for sediment =                 | mg/day    | 200                                 | USEPA 1997, EFH       |  |  |  |
| EF - Exposure frequency =                            | days/year | 40                                  | reasonable assumption |  |  |  |
| ED - Exposure duration =                             | years     | 6                                   | USEPA 1995, Region IV |  |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06                            |                       |  |  |  |
| ME - Matrix effect =                                 |           | 1                                   | Magee, et al., 1996   |  |  |  |
| BW - Body weight =                                   | kg        | 15                                  | USEPA 1995, Region IV |  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 2190                                | USEPA 1991, HHEM      |  |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550                               | USEPA 1991, HHEM      |  |  |  |

| Constituent                | Concentration in | Average Daily    | Oral                     | Average      |                                 | Cancer Risk |          |
|----------------------------|------------------|------------------|--------------------------|--------------|---------------------------------|-------------|----------|
|                            | Sediment mg/kg   | Intake mg/kg-day | Subchronic RfD mg/kg-day | Hazard Index | Lifetime Daily Intake mg/kg-day |             |          |
| <b>Semivolatiles</b>       |                  |                  |                          |              |                                 |             |          |
| 2-Nitroaniline             | 4.00E-01         | 5.84E-07         | NA                       | NA           | 5.01E-08                        | NA          |          |
| 2-Nitrophenol              | 8.00E-01         | 1.17E-06         | NA                       | NA           | 1.00E-07                        | NA          |          |
| 3-Nitroaniline             | 8.00E-01         | 1.17E-06         | NA                       | NA           | 1.00E-07                        | NA          |          |
| 4-Bromophenylphenoxyether  | 8.00E-01         | 1.17E-06         | NA                       | NA           | 1.00E-07                        | NA          |          |
| 4-Chloro-3-methylphenol    | 8.00E-01         | 1.17E-06         | NA                       | NA           | 1.00E-07                        | NA          |          |
| 4-Chlorophenylphenoxyether | 4.00E-01         | 5.84E-07         | NA                       | NA           | 5.01E-08                        | NA          |          |
| 4-Nitroaniline             | 8.00E-01         | 1.17E-06         | NA                       | NA           | 1.00E-07                        | NA          |          |
| Benzo(a)anthracene         | 1.00E+02         | 1.46E-04         | NA                       | NA           | 1.25E-05                        | 7.30E-01    | 9.14E-06 |
| Benzo(a)pyrene             | 4.90E+01         | 7.16E-05         | NA                       | NA           | 6.14E-06                        | 7.30E+00    | 4.48E-05 |
| Benzo(b)fluoranthene       | 7.80E+01         | 1.14E-04         | NA                       | NA           | 9.77E-06                        | 7.30E-01    | 7.13E-06 |
| Benzo(k)fluoranthene       | 2.30E+01         | 3.36E-05         | NA                       | NA           | 2.88E-06                        | 7.30E-02    | 2.10E-07 |
| Bis(2-chloroethoxy)methane | 8.00E-01         | 1.17E-06         | NA                       | NA           | 1.00E-07                        | NA          | NA       |
| Bis(2-chloroethyl)ether    | 4.00E-01         | 5.84E-07         | NA                       | NA           | 5.01E-08                        | 1.10E+00    | 5.51E-08 |
| Carbazole                  | 1.00E+02         | 1.46E-04         | NA                       | NA           | 1.25E-05                        | 2.00E-02    | 2.50E-07 |
| Chrysene                   | 7.60E+01         | 1.11E-04         | NA                       | NA           | 9.52E-06                        | 7.30E-03    | 6.95E-08 |
| Dibenz(a,h)anthracene      | 9.60E+00         | 1.40E-05         | NA                       | NA           | 1.20E-06                        | 7.30E+00    | 8.78E-06 |
| Hexachlorobenzene          | 4.00E-01         | 5.84E-07         | NA                       | NA           | 5.01E-08                        | 1.60E+00    | 8.02E-08 |
| Hexachlorocyclopentadiene  | 2.00E+00         | 2.92E-06         | NA                       | NA           | 2.50E-07                        | NA          | NA       |
| Indeno(1,2,3-cd)pyrene     | 3.90E+01         | 5.70E-05         | NA                       | NA           | 4.88E-06                        | 7.30E-01    | 3.57E-06 |
| N-nitrosodi-n-propylamine  | 4.00E-01         | 5.84E-07         | NA                       | NA           | 5.01E-08                        | 7.00E+00    | 3.51E-07 |

NA - Not Applicable

Total Cancer Risk = 7.44E-05

**Table 73****Dermal Exposure to EU6 Sediment by an Adult Resident (Aged 7 to 30 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =       |                                 | $\frac{\text{Cs} * \text{SA} * \text{AH} * \text{ABS} * \text{EF} * \text{ED} * \text{CF}}{\text{BW} * \text{AT}}$ |                              |              |   |                                   |             |
|----------------------------|---------------------------------|--|------------------------------|--------------|---|-----------------------------------|-------------|
| Constituent                | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day   | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles              |                                 |  |                              |              |   |                                   |             |
| 2-Nitroaniline             | 4.00E-01                        | 1.28E-07   | NA                           | NA           | 4.38E-08                                | NA                                | NA          |
| 2-Nitrophenol              | 8.00E-01                        | 2.55E-07   | NA                           | NA           | 8.76E-08                                | NA                                | NA          |
| 3-Nitroaniline             | 8.00E-01                        | 2.55E-07   | NA                           | NA           | 8.76E-08                                | NA                                | NA          |
| 4-Bromophenylphenoxyether  | 8.00E-01                        | 2.55E-07   | NA                           | NA           | 8.76E-08                                | NA                                | NA          |
| 4-Chloro-3-methylphenol    | 8.00E-01                        | 2.55E-07   | NA                           | NA           | 8.76E-08                                | NA                                | NA          |
| 4-Chlorophenylphenoxyether | 4.00E-01                        | 1.28E-07   | NA                           | NA           | 4.38E-08                                | NA                                | NA          |
| 4-Nitroaniline             | 8.00E-01                        | 2.55E-07   | NA                           | NA           | 8.76E-08                                | NA                                | NA          |
| Benzo(a)anthracene         | 1.00E+02                        | 9.58E-06   | NA                           | NA           | 3.28E-06                                | 1.46E+00                          | 4.79E-06    |
| Benzo(a)pyrene             | 4.90E+01                        | 4.69E-06   | NA                           | NA           | 1.61E-06                                | 1.46E+01                          | 2.35E-05    |
| Benzo(b)fluoranthene       | 7.80E+01                        | 7.47E-06   | NA                           | NA           | 2.56E-06                                | 1.46E+00                          | 3.74E-06    |
| Benzo(k)fluoranthene       | 2.30E+01                        | 2.20E-06   | NA                           | NA           | 7.55E-07                                | 1.46E-01                          | 1.10E-07    |
| Bis(2-chloroethoxy)methane | 8.00E-01                        | 2.55E-07   | NA                           | NA           | 8.76E-08                                | NA                                | NA          |
| Bis(2-chloroethyl)ether    | 4.00E-01                        | 1.28E-07   | NA                           | NA           | 4.38E-08                                | 1.10E+00                          | 4.82E-08    |
| Carbazole                  | 1.00E+02                        | 3.19E-05   | NA                           | NA           | 1.09E-05                                | 2.00E-02                          | 2.19E-07    |
| Chrysene                   | 7.60E+01                        | 7.28E-06   | NA                           | NA           | 2.50E-06                                | 1.46E-02                          | 3.64E-08    |
| Dibenz(a,h)anthracene      | 9.60E+00                        | 9.20E-07   | NA                           | NA           | 3.15E-07                                | 1.46E+01                          | 4.60E-06    |
| Hexachlorobenzene          | 4.00E-01                        | 1.28E-07   | NA                           | NA           | 4.38E-08                                | 1.60E+00                          | 7.01E-08    |
| Hexachlorocyclopentadiene  | 2.00E+00                        | 6.39E-07   | NA                           | NA           | 2.19E-07                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene     | 3.90E+01                        | 3.74E-06   | NA                           | NA           | 1.28E-06                                | 1.46E+00                          | 1.87E-06    |
| N-nitrosodi-n-propylamine  | 4.00E-01                        | 1.28E-07   | NA                           | NA           | 4.38E-08                                | 7.00E+00                          | 3.07E-07    |

NA - Not Available

Total Cancer Risk = 3.93E-05



**Table 74****Oral Exposure to EU6 Sediment by an Adult Resident (Aged 7 to 30 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |                                 | <u>Cd*Ingr*EF*ED*CF*ME</u><br>BW*AT |                            |              |   |  |
|--|---------------------------------|-------------------------------------|----------------------------|--------------|---|--|
| Cd - Concentration in sediment =                     | mg/kg                           | see below                           |                            |              |   |  |
| Ingr - Ingestion rate for sediment =                 | mg/day                          | 100                                 | USEPA 1997, EFH            |              |   |  |
| EF - Exposure frequency =                            | days/year                       | 40                                  | reasonable assumption      |              |   |  |
| ED - Exposure duration =                             | years                           | 24                                  | USEPA 1995, Region IV      |              |   |  |
| CF - Conversion factor =                             | kg/mg                           | 1.00E-06                            |                            |              |   |  |
| ME - Matrix effect =                                 |                                 | 1                                   | Magee, et al., 1996        |              |   |  |
| BW - Body weight =                                   | kg                              | 70                                  | USEPA 1995, Region IV      |              |   |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                            | 8760                                | USEPA 1991, HHEM           |              |   |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                            | 25550                               | USEPA 1991, HHEM           |              |   |  |
| Constituent  | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day      | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) |
| Semivolatiles  |                                 |                                     |                            |              |   |  |
| 2-Nitroaniline                                       | 4.00E-01                        | 6.26E-08                            | NA                         | NA           | 2.15E-08                                | NA                                     |
| 2-Nitrophenol  | 8.00E-01                        | 1.25E-07                            | NA                         | NA           | 4.29E-08                                | NA                                     |
| 3-Nitroaniline                                       | 8.00E-01                        | 1.25E-07                            | NA                         | NA           | 4.29E-08                                | NA                                     |
| 4-Bromophenylphenylether                             | 8.00E-01                        | 1.25E-07                            | NA                         | NA           | 4.29E-08                                | NA                                     |
| 4-Chloro-3-methylphenol                              | 8.00E-01                        | 1.25E-07                            | NA                         | NA           | 4.29E-08                                | NA                                     |
| 4-Chlorophenylphenylether                            | 4.00E-01                        | 6.26E-08                            | NA                         | NA           | 2.15E-08                                | NA                                     |
| 4-Nitroaniline                                       | 8.00E-01                        | 1.25E-07                            | NA                         | NA           | 4.29E-08                                | NA                                     |
| Benzo(a)anthracene                                   | 1.00E+02                        | 1.57E-05                            | NA                         | NA           | 5.37E-06                                | 7.30E-01                               |
| Benzo(a)pyrene                                       | 4.90E+01                        | 7.67E-06                            | NA                         | NA           | 2.63E-06                                | 7.30E+00                               |
| Benzo(b)fluoranthene                                 | 7.80E+01                        | 1.22E-05                            | NA                         | NA           | 4.19E-06                                | 7.30E-01                               |
| Benzo(k)fluroanthene                                 | 2.30E+01                        | 3.60E-06                            | NA                         | NA           | 1.23E-06                                | 7.30E-02                               |
| Bis(2-chloroethoxy)methane                           | 8.00E-01                        | 1.25E-07                            | NA                         | NA           | 4.29E-08                                | NA                                     |
| Bis(2-chloroethyl)ether                              | 4.00E-01                        | 6.26E-08                            | NA                         | NA           | 2.15E-08                                | 1.10E+00                               |
| Carbazole  | 1.00E+02                        | 1.57E-05                            | NA                         | NA           | 5.37E-06                                | 2.00E-02                               |
| Chrysene   | 7.60E+01                        | 1.19E-05                            | NA                         | NA           | 4.08E-06                                | 7.30E-03                               |
| Dibenz(a,h)anthracene                                | 9.60E+00                        | 1.50E-06                            | NA                         | NA           | 5.15E-07                                | 7.30E+00                               |
| Hexachlorobenzene                                    | 4.00E-01                        | 6.26E-08                            | 8.00E-04                   | 7.83E-05     | 2.15E-08                                | 1.60E+00                               |
| Hexachlorocyclopentadiene                            | 2.00E+00                        | 3.13E-07                            | 7.00E-03                   | 4.47E-05     | 1.07E-07                                | NA                                     |
| Indeno(1,2,3-cd)pyrene                               | 3.90E+01                        | 6.11E-06                            | NA                         | NA           | 2.09E-06                                | 7.30E-01                               |
| N-nitrosodi-n-propylamine                            | 4.00E-01                        | 6.26E-08                            | NA                         | NA           | 2.15E-08                                | 7.00E+00                               |

NA - Not Applicable

Total Hazard Index = 1.23E-04

Total Cancer Risk = 3.19E-05



Table 75

Dermal Exposure to EU6 Surface Water by a Child Resident (Aged 1 to 6 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cw*SA*Kp*ABS*ET*EF*ED*CF</u> |           |  |  |  |
|---|---------------------------------|-----------|--|--|--|
|   | BW*AT                           |           |  |  |  |
| Cw - Concentration in surface water =                       | mg/L                            | see below |  |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup>                 | 2229      | calculated                             |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                 | 7213      | USEPA 1997, EFH                        |  |  |
| Fs - Fraction of skin surface area available for exposure = |                                 | 30.9%     | USEPA 1997, EFH                        |  |  |
| K <sub>p</sub> - Dermal permeability constant =             | cm/hr                           | see below |  |  |  |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                                 | 0.03      | USEPA 1995, Region III                 |  |  |
| ET - Exposure time =  | hrs/day                         | 1         | USEPA 1992, Dermal Exposure Assessment |  |  |
| EF - Exposure frequency =                                   | days/year                       | 40        | reasonable assumption                  |  |  |
| ED - Exposure duration =                                    | years                           | 6         | USEPA 1995, Region IV                  |  |  |
| CF - Conversion factor =                                    | L/cm <sup>3</sup>               | 1.00E-03  |  |  |  |
| BW - Body weight =  | kg                              | 15        | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                            | 2190      | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                            | 25550     | USEPA 1991, HHEM                       |  |  |

| Constituent            | Concentration in Surface Water mg/L | K <sub>p</sub> cm/hr | Average Daily Intake mg/kg-day | Dermal Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-------------------------------------|----------------------|--------------------------------|---------------------------------|--------------|---|-----------------------------------|-------------|
| <b>Semivolatiles</b>   |                                     |                      |                                |                                 |              |   |                                   |             |
| Benzo(a)anthracene     | 5.00E-04                            | 8.10E-01             | 1.98E-07                       | NA                              | NA           | 1.70E-08                                | 1.46E+00                          | 2.48E-08    |
| Benzo(a)pyrene         | 5.00E-04                            | 1.20E+00             | 2.93E-07                       | NA                              | NA           | 2.51E-08                                | 1.46E+01                          | 3.67E-07    |
| Benzo(b)fluoranthene   | 9.00E-03                            | 1.20E+00             | 5.28E-06                       | NA                              | NA           | 4.52E-07                                | 1.46E+00                          | 6.60E-07    |
| Benzo(k)fluoranthene   | 5.00E-04                            | 4.48E+01             | 1.09E-05                       | NA                              | NA           | 9.38E-07                                | 1.46E-01                          | 1.37E-07    |
| Chrysene               | 5.00E-04                            | 8.10E-01             | 1.98E-07                       | NA                              | NA           | 1.70E-08                                | 1.46E-02                          | 2.48E-10    |
| Dibenz(a,h)anthracene  | 5.00E-04                            | 2.70E+00             | 6.59E-07                       | NA                              | NA           | 5.65E-08                                | 1.46E+01                          | 8.25E-07    |
| Indeno(1,2,3-cd)pyrene | 5.00E-04                            | 1.90E+00             | 4.64E-07                       | NA                              | NA           | 3.98E-08                                | 1.46E+00                          | 5.81E-08    |

NA - Not Available

Total Cancer Risk = 2.07E-06



**Table 76****Oral Exposure to EU6 Surface Water by a Child Resident (Aged 1 to 6 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 | <u>Csw*IngR*EF*ED*ET</u> |           |  |
|--|--------------------------|-----------|--|
|  | BW                       | AT        |  |
| Csw - Concentration in surface water =               | mg/L                     | see below |  |
| IngR - Ingestion rate for surface water =            | L/hour                   | 0.05      | USEPA 1995, Region IV                  |
| EF - Exposure frequency =                            | days/year                | 40        | reasonable assumption                  |
| ED - Exposure duration =                             | years                    | 6         | USEPA 1995, Region IV                  |
| ET - Exposure time =                                 | hrs/day                  | 1         | USEPA 1992, Dermal Exposure Assessment |
| BW - Body weight =                                   | kg                       | 15        | USEPA 1995, Region IV                  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                     | 2190      | USEPA 1991, HHEM                       |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                     | 25550     | USEPA 1991, HHEM                       |

| Constituent            | Concentration in Surface Water<br>mg/L | Average Daily Intake<br>mg/kg-day | Oral Subchronic RfD<br>mg/kg-day | Hazard Index | Average Lifetime Daily Intake |               |                          | Cancer Risk |
|------------------------|--|-----------------------------------|----------------------------------|--------------|-------------------------------|---------------|--------------------------|-------------|
|                        |  |                                   |                                  |              | mg/kg-day                     | 1/(mg/kg-day) | Oral Cancer Slope Factor |             |
| <b>Semivolatiles</b>   |  |                                   |                                  |              |                               |               |                          |             |
| Benzo(a)anthracene     | 5.00E-04                               | 1.83E-07                          | NA                               | NA           | 1.57E-08                      | 7.30E-01      | 1.14E-08                 |             |
| Benzo(a)pyrene         | 5.00E-04                               | 1.83E-07                          | NA                               | NA           | 1.57E-08                      | 7.30E+00      | 1.14E-07                 |             |
| Benzo(b)fluoranthene   | 9.00E-03                               | 3.29E-06                          | NA                               | NA           | 2.82E-07                      | 7.30E-01      | 2.06E-07                 |             |
| Benzo(k)fluoranthene   | 5.00E-04                               | 1.83E-07                          | NA                               | NA           | 1.57E-08                      | 7.30E-02      | 1.14E-09                 |             |
| Chrysene               | 5.00E-04                               | 1.83E-07                          | NA                               | NA           | 1.57E-08                      | 7.30E-03      | 1.14E-10                 |             |
| Dibenz(a,h)anthracene  | 5.00E-04                               | 1.83E-07                          | NA                               | NA           | 1.57E-08                      | 7.30E+00      | 1.14E-07                 |             |
| Indeno(1,2,3-cd)pyrene | 5.00E-04                               | 1.83E-07                          | NA                               | NA           | 1.57E-08                      | 7.30E-01      | 1.14E-08                 |             |

NA - Not Applicable

Total Cancer Risk = 4.58E-07



**Table 77****Dermal Exposure to EU6 Surface Water by an Adult Resident (Aged 7 to 30 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |  | <u>C<sub>w</sub>*S<sub>A</sub>*K<sub>p</sub>*A<sub>B</sub>S*E<sub>T</sub>*E<sub>F</sub>*E<sub>D</sub>*C<sub>F</sub></u><br>BW*AT |                                   |  |                                    |                                      |             |          |
|---|--|--|-----------------------------------|--|------------------------------------|--------------------------------------|-------------|----------|
| C <sub>w</sub> - Concentration in surface water =                       |  | mg/L   | see below                         |  |                                    |                                      |             |          |
| S <sub>A</sub> - Surface area available for exposure =                  |  | cm <sup>2</sup>  | 6180                              | calculated                             |                                    |                                      |             |          |
| S <sub>A<sub>t</sub></sub> - Total skin surface area =                  |  | cm <sup>2</sup>  | 20000                             | USEPA 1997, EFH                        |                                    |                                      |             |          |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = |  |  | 30.9%                             | USEPA 1997, EFH                        |                                    |                                      |             |          |
| K <sub>p</sub> - Dermal permeability constant =                         |  | cm/hr  | see below                         |  |                                    |                                      |             |          |
| A <sub>B</sub> S <sub>p</sub> - Absorption - cPAHs =                    |  |  | 0.03                              | USEPA 1995, Region III                 |                                    |                                      |             |          |
| E <sub>T</sub> - Exposure time =  |  | hrs/day  | 1                                 | USEPA 1992, Dermal Exposure Assessment |                                    |                                      |             |          |
| E <sub>F</sub> - Exposure frequency =                                   |  | days/year  | 40                                | reasonable assumption                  |                                    |                                      |             |          |
| E <sub>D</sub> - Exposure duration =                                    |  | years  | 24                                | USEPA 1995, Region IV                  |                                    |                                      |             |          |
| C <sub>F</sub> - Conversion factor =                                    |  | L/cm <sup>2</sup>  | 1.00E-03                          |  |                                    |                                      |             |          |
| BW - Body weight =  |  | kg   | 70                                | USEPA 1995, Region IV                  |                                    |                                      |             |          |
| A <sub>T<sub>n</sub></sub> - Averaging time - noncarcinogenic =         |  | days   | 8760                              | USEPA 1991, HHEM                       |                                    |                                      |             |          |
| A <sub>T<sub>c</sub></sub> - Averaging time - carcinogenic =            |  | days   | 25550                             | USEPA 1991, HHEM                       |                                    |                                      |             |          |
| Constituent   | Concentration in Surface Water<br>mg/L | K <sub>p</sub><br>cm/hr  | Average Daily Intake<br>mg/kg-day | Dermal Chronic RfD<br>mg/kg-day        | Average                            |                                      |             |          |
|   |  |  |                                   | Hazard Index                           | Lifetime Daily Intake<br>mg/kg-day | Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |          |
| <b>Semivolatiles</b>  |  |  |                                   |  |                                    |                                      |             |          |
| Benzo(a)anthracene  | 5.00E-04                               | 8.10E-01   | 1.18E-07                          | NA                                     | NA                                 | 4.03E-08                             | 1.46E+00    | 5.88E-08 |
| Benzo(a)pyrene  | 5.00E-04                               | 1.20E+00   | 1.74E-07                          | NA                                     | NA                                 | 5.97E-08                             | 1.46E+01    | 8.72E-07 |
| Benzo(b)fluoranthene  | 9.00E-03                               | 1.20E+00   | 3.13E-06                          | NA                                     | NA                                 | 1.07E-06                             | 1.46E+00    | 1.57E-06 |
| Benzo(k)fluoranthene  | 5.00E-04                               | 4.48E+01   | 6.50E-06                          | NA                                     | NA                                 | 2.23E-06                             | 1.46E-01    | 3.25E-07 |
| Chrysene  | 5.00E-04                               | 8.10E-01   | 1.18E-07                          | NA                                     | NA                                 | 4.03E-08                             | 1.46E-02    | 5.88E-10 |
| Dibenz(a,h)anthracene   | 5.00E-04                               | 2.70E+00   | 3.92E-07                          | NA                                     | NA                                 | 1.34E-07                             | 1.46E+01    | 1.96E-06 |
| Indeno(1,2,3-cd)pyrene  | 5.00E-04                               | 1.90E+00   | 2.76E-07                          | NA                                     | NA                                 | 9.45E-08                             | 1.46E+00    | 1.38E-07 |

NA - Not Available

Total Cancer Risk = 4.93E-06



**Table 78****Oral Exposure to EU6 Surface Water by an Adult Resident (Aged 7 to 30 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u><math>C_{sw} \times Ingr \times EF \times ED \times ET</math></u> |  |  |  |
|--|-----------|--|--|--|--|
|  |           | <u><math>BW \times AT</math></u>                                     |  |  |  |
| Csw - Concentration in surface water =               | mg/L      | see below  |  |  |  |
| Ingr - Ingestion rate for surface water =            | L/hour    | 0.01   | USEPA 1995, Region IV                  |  |  |
| EF - Exposure frequency =                            | days/year | 40   | reasonable assumption                  |  |  |
| ED - Exposure duration =                             | years     | 24   | USEPA 1995, Region IV                  |  |  |
| ET - Exposure time =                                 | hrs/day   | 1  | USEPA 1992, Dermal Exposure Assessment |  |  |
| BW - Body weight =                                   | kg        | 70   | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 8760   | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550  | USEPA 1991, HHEM                       |  |  |

| Constituent            | Concentration in Surface Water mg/L | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |       |          | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-------------------------------------|--------------------------------|----------------------------|--------------|---|-------|----------|--|-------------|
|                        |                                     |                                |                            |              | Lifetime                                | Daily | Intake   |  |             |
| <b>Semivolatiles</b>   |                                     |                                |                            |              |   |       |          |  |             |
| Benzo(a)anthracene     | 5.00E-04                            | 7.83E-09                       | NA                         | NA           | 2.68E-09                                |       | 7.30E-01 | 1.96E-09                               |             |
| Benzo(a)pyrene         | 5.00E-04                            | 7.83E-09                       | NA                         | NA           | 2.68E-09                                |       | 7.30E+00 | 1.96E-08                               |             |
| Benzo(b)fluoranthene   | 9.00E-03                            | 1.41E-07                       | NA                         | NA           | 4.83E-08                                |       | 7.30E-01 | 3.53E-08                               |             |
| Benzo(k)fluoranthene   | 5.00E-04                            | 7.83E-09                       | NA                         | NA           | 2.68E-09                                |       | 7.30E-02 | 1.96E-10                               |             |
| Chrysene               | 5.00E-04                            | 7.83E-09                       | NA                         | NA           | 2.68E-09                                |       | 7.30E-03 | 1.96E-11                               |             |
| Dibenz(a,h)anthracene  | 5.00E-04                            | 7.83E-09                       | NA                         | NA           | 2.68E-09                                |       | 7.30E+00 | 1.96E-08                               |             |
| Indeno(1,2,3-cd)pyrene | 5.00E-04                            | 7.83E-09                       | NA                         | NA           | 2.68E-09                                |       | 7.30E-01 | 1.96E-09                               |             |

NA - Not Applicable

Total Cancer Risk = 7.86E-08



**Table 79****Dermal Exposure to EU4 Sediment by an Adolescent Visitor (Aged 7-16 years)****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =   |                                 | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                              |              |   |                                   |             |
|------------------------|---------------------------------|---|------------------------------|--------------|---|-----------------------------------|-------------|
| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day                      | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles          |                                 |   |                              |              |   |                                   |             |
| Benzo(a)anthracene     | 9.00E+00                        | 2.57E-07  | NA                           | NA           | 3.67E-08                                | 1.46E+00                          | 5.36E-08    |
| Benzo(a)pyrene         | 5.57E+00                        | 1.59E-07  | NA                           | NA           | 2.27E-08                                | 1.46E+01                          | 3.32E-07    |
| Benzo(b)fluoranthene   | 6.31E+00                        | 1.80E-07  | NA                           | NA           | 2.57E-08                                | 1.46E+00                          | 3.76E-08    |
| Benzo(k)fluoranthene   | 3.44E+00                        | 9.82E-08  | NA                           | NA           | 1.40E-08                                | 1.46E-01                          | 2.05E-09    |
| Carbazole              | *                               |   | NA                           | NA           |   | 2.00E-02                          |             |
| Chrysene               | 7.10E+00                        | 2.03E-07  | NA                           | NA           | 2.89E-08                                | 1.46E-02                          | 4.23E-10    |
| Dibenz(a,h)anthracene  | 8.40E-01                        | 2.40E-08  | NA                           | NA           | 3.42E-09                                | 1.46E+01                          | 5.00E-08    |
| Dibenzofuran           | *                               |   | 2.00E-03                     |              |   |                                   |             |
| Indeno(1,2,3-cd)pyrene | 3.40E+00                        | 9.70E-08  | NA                           | NA           | 1.39E-08                                | 1.46E+00                          | 2.02E-08    |
| Naphthalene            | 8.20E+00                        | 7.80E-07  | 1.00E-02                     | 7.80E-05     | 1.11E-07                                | NA                                | NA          |
| Phenanthrene           | 2.38E+01                        | 2.26E-06  | NA                           | NA           | 3.23E-07                                | NA                                | NA          |

NA - Not Available

Total Hazard Index = 7.80E-05

Total Cancer Risk = 4.95E-07

\*Constituent not present in remaining samples.



**Table 80****Oral Exposure to EU4 Sediment by an Adolescent Visitor (Aged 7-16 years)****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                                      |                                  |                 |                                       |  |                |
|--|---|--------------------------------------|----------------------------------|-----------------|---------------------------------------|--|----------------|
| Cd - Concentration in sediment =                     | mg/kg   | see below                            |                                  |                 |                                       |  |                |
| IngR - Ingestion rate for sediment =                 | mg/day  | 100                                  | USEPA 1997, EFH                  |                 |                                       |  |                |
| EF - Exposure frequency =                            | days/year   | 12                                   | reasonable assumption            |                 |                                       |  |                |
| ED - Exposure duration =                             | years   | 10                                   | USEPA 1995, Region IV            |                 |                                       |  |                |
| CF - Conversion factor =                             | kg/mg   | 1.00E-06                             |                                  |                 |                                       |  |                |
| ME - Matrix effect =                                 |   | 1                                    | Magee, et al., 1996              |                 |                                       |  |                |
| BW - Body weight =                                   | kg  | 45                                   | USEPA 1995, Region IV            |                 |                                       |  |                |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days  | 3650                                 | USEPA 1991, HHEM                 |                 |                                       |  |                |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days  | 25550                                | USEPA 1991, HHEM                 |                 |                                       |  |                |
| Average  |   |                                      |                                  |                 |                                       |  |                |
| Constituent  | Concentration<br>in Sediment<br>mg/kg   | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Lifetime Daily<br>Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
| Semivolatiles  |   |                                      |                                  |                 |                                       |  |                |
| Benzo(a)anthracene                                   | 9.00E+00  | 6.58E-07                             | NA                               | NA              | 9.39E-08                              | 7.30E-01                                     | 6.86E-08       |
| Benzo(a)pyrene                                       | 5.57E+00  | 4.07E-07                             | NA                               | NA              | 5.81E-08                              | 7.30E+00                                     | 4.24E-07       |
| Benzo(b)fluoranthene                                 | 6.31E+00  | 4.61E-07                             | NA                               | NA              | 6.59E-08                              | 7.30E-01                                     | 4.81E-08       |
| Benzo(k)fluoranthene                                 | 3.44E+00  | 2.51E-07                             | NA                               | NA              | 3.59E-08                              | 7.30E-02                                     | 2.62E-09       |
| Carbazole  | *   |                                      | NA                               | NA              |                                       | 2.00E-02                                     |                |
| Chrysene   | 7.10E+00  | 5.19E-07                             | NA                               | NA              | 7.41E-08                              | 7.30E-03                                     | 5.41E-10       |
| Dibenz(a,h)anthracene                                | 8.40E-01  | 6.14E-08                             | NA                               | NA              | 8.77E-09                              | 7.30E+00                                     | 6.40E-08       |
| Dibenzofuran   | *   |                                      | 4.00E-03                         |                 |                                       | NA   | NA             |
| Indeno(1,2,3-cd)pyrene                               | 3.40E+00  | 2.48E-07                             | NA                               | NA              | 3.55E-08                              | 7.30E-01                                     | 2.59E-08       |
| Naphthalene  | 8.20E+00  | 5.99E-07                             | 2.00E-02                         | 3.00E-05        | 8.56E-08                              | NA   | NA             |
| Phenanthrene   | 2.38E+01  | 1.74E-06                             | NA                               | NA              | 2.48E-07                              | NA   | NA             |

NA - Not Applicable

Total Hazard Index = 3.00E-05

Total Cancer Risk = 6.34E-07

\*Constituent not present in remaining samples.



**Table 81****Oral Exposure to EU4 Sediment by a Maintenance Worker****Preliminary Remediation Goal Calculation**

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =                                 |           | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                       |  |  |
|--|-----------|---|-----------------------|--|--|
| Cd - Concentration in sediment =                     | mg/kg     | see below   |                       |  |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 100   | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 2   | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 25  | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06  |                       |  |  |
| ME - Matrix effect =                                 |           | 1   | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 70  | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 9125  | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550   | USEPA 1991, HHEM      |  |  |

| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|---------------------------------|--------------------------------|----------------------------|--------------|---|--|-------------|
| <b>Semivolatiles</b>   |                                 |                                |                            |              |   |  |             |
| Benzo(a)anthracene     | 9.00E+00                        | 7.05E-08                       | NA                         | NA           | 2.52E-08                                | 7.30E-01                               | 1.84E-08    |
| Benzo(a)pyrene         | 5.57E+00                        | 4.36E-08                       | NA                         | NA           | 1.56E-08                                | 7.30E+00                               | 1.14E-07    |
| Benzo(b)fluoranthene   | 6.31E+00                        | 4.94E-08                       | NA                         | NA           | 1.76E-08                                | 7.30E-01                               | 1.29E-08    |
| Benzo(k)fluoranthene   | 3.44E+00                        | 2.69E-08                       | NA                         | NA           | 9.62E-09                                | 7.30E-02                               | 7.02E-10    |
| Carbazole              | *                               |                                | NA                         | NA           |   | 2.00E-02                               |             |
| Chrysene               | 7.10E+00                        | 5.56E-08                       | NA                         | NA           | 1.98E-08                                | 7.30E-03                               | 1.45E-10    |
| Dibenz(a,h)anthracene  | 8.40E-01                        | 6.58E-09                       | NA                         | NA           | 2.35E-09                                | 7.30E+00                               | 1.71E-08    |
| Dibenzofuran           | *                               |                                | 4.00E-03                   |              |   | NA                                     | NA          |
| Indeno(1,2,3-cd)pyrene | 3.40E+00                        | 2.66E-08                       | NA                         | NA           | 9.51E-09                                | 7.30E-01                               | 6.94E-09    |
| Naphthalene            | 8.20E+00                        | 6.42E-08                       | 2.00E-02                   | 3.21E-06     | 2.29E-08                                | NA                                     | NA          |
| Phenanthrene           | 2.38E+01                        | 1.86E-07                       | NA                         | NA           | 6.65E-08                                | NA                                     | NA          |

NA - Not Applicable

Total Hazard Index = 3.21E-06

Total Cancer Risk = 1.70E-07

\*Constituent not present in remaining samples.



**Table 82****Oral Exposure to EU4 Sediment by a Construction Worker****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =   |                                 | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                               |              |   |  |             |
|------------------------|---------------------------------|---|-------------------------------|--------------|---|--|-------------|
| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day  | Oral Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles          |                                 |   |                               |              |   |  |             |
| Benzo(a)anthracene     | 9.00E+00                        | 1.35E-06  | NA                            | NA           | 1.93E-08                                | 7.30E-01                               | 1.41E-08    |
| Benzo(a)pyrene         | 5.57E+00                        | 8.37E-07  | NA                            | NA           | 1.20E-08                                | 7.30E+00                               | 8.73E-08    |
| Benzo(b)fluoranthene   | 6.31E+00                        | 9.48E-07  | NA                            | NA           | 1.35E-08                                | 7.30E-01                               | 9.89E-09    |
| Benzo(k)fluoranthene   | 3.44E+00                        | 5.17E-07  | NA                            | NA           | 7.39E-09                                | 7.30E-02                               | 5.39E-10    |
| Carbazole              | *                               |   | NA                            | NA           |   | 2.00E-02                               |             |
| Chrysene               | 7.10E+00                        | 1.07E-06  | NA                            | NA           | 1.52E-08                                | 7.30E-03                               | 1.11E-10    |
| Dibenz(a,h)anthracene  | 8.40E-01                        | 1.26E-07  | NA                            | NA           | 1.80E-09                                | 7.30E+00                               | 1.32E-08    |
| Dibenzofuran           | *                               |   | NA                            | NA           |   | NA                                     | NA          |
| Indeno(1,2,3-cd)pyrene | 3.40E+00                        | 5.11E-07  | NA                            | NA           | 7.30E-09                                | 7.30E-01                               | 5.33E-09    |
| Naphthalene            | 8.20E+00                        | 1.23E-06  | NA                            | NA           | 1.76E-08                                | NA                                     | NA          |
| Phenanthrene           | 2.38E+01                        | 3.58E-06  | NA                            | NA           | 5.11E-08                                | NA                                     | NA          |

NA - Not Applicable

Total Cancer Risk = 1.30E-07

\*Constituent not present in remaining samples.



**Table 83**

**Dermal Exposure to EU4 Surface Soil (0-6') by a Maintenance Worker**

**Preliminary Remediation Goal Calculation**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                        |  |
|---|----------------------|---|------------------------|--|
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.   |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 3000  | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000   | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 15%   | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.038   | USEPA 1997, EFH        |  |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                      | 0.03  | USEPA 1995, Region III |  |
| ABS <sub>s</sub> - Absorption - other SVOCs =               |                      | 0.1   | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year            | 150   | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                | 25  | USEPA 1995, Region IV  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06  |                        |  |
| BW - Body weight =  | kg                   | 70  | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 9125  | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550   | USEPA 1991, HHEM       |  |

| Constituent            | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal   |                 | Average                               |   |                |
|------------------------|-----------------------------------|--------------------------------------|----------|-----------------|---------------------------------------|---|----------------|
|                        |                                   |                                      | RfD      | Hazard<br>Index | Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
| <b>Semivolatiles</b>   |                                   |                                      |          |                 |                                       |   |                |
| Benzo(a)anthracene     | 8.10E-01                          | 1.63E-08                             | NA       | NA              | 5.81E-09                              | 1.46E+00                                | 8.48E-09       |
| Benzo(a)pyrene         | 2.90E-01                          | 5.82E-09                             | NA       | NA              | 2.08E-09                              | 1.46E+01                                | 3.04E-08       |
| Benzo(b)fluoranthene   | 3.70E-01                          | 7.43E-09                             | NA       | NA              | 2.65E-09                              | 1.46E+00                                | 3.87E-09       |
| Benzo(k)fluoranthene   | 1.60E-01                          | 3.21E-09                             | NA       | NA              | 1.15E-09                              | 1.46E-01                                | 1.68E-10       |
| Carbazole              | 4.90E-01                          | 3.28E-08                             | NA       | NA              | 1.17E-08                              | 2.00E-02                                | 2.34E-10       |
| Chrysene               | 6.10E-01                          | 1.22E-08                             | NA       | NA              | 4.37E-09                              | 1.46E-02                                | 6.39E-11       |
| Dibenz(a,h)anthracene  | 1.10E-02                          | 2.21E-10                             | NA       | NA              | 7.89E-11                              | 1.46E+01                                | 1.15E-09       |
| Indeno(1,2,3-cd)pyrene | 9.40E-02                          | 1.89E-09                             | NA       | NA              | 6.74E-10                              | 1.46E+00                                | 9.84E-10       |
| Naphthalene            | 4.00E-01                          | 2.68E-08                             | 1.00E-02 | 2.68E-06        | 9.56E-09                              | NA                                      | NA             |

NA - Not Available

Total Hazard Index = 2.68E-06

Total Cancer Risk = 4.53E-08



**Table 84****Oral Exposure to EU4 Surface Soil (0-6') by a Maintenance Worker****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |                             | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                            |              |   |  |
|--|-----------------------------|---|----------------------------|--------------|---|--|
| Cd - Concentration in soil =                         | mg/kg                       | see below   |                            |              |   |  |
| IngR - Ingestion rate for soil =                     | mg/day                      | 100   | USEPA 1997, EFH            |              |   |  |
| EF - Exposure frequency =                            | days/year                   | 150   | reasonable assumption      |              |   |  |
| ED - Exposure duration =                             | years                       | 25  | USEPA 1995, Region IV      |              |   |  |
| CF - Conversion factor =                             | kg/mg                       | 1.00E-06  |                            |              |   |  |
| ME - Matrix effect =                                 |                             | 1   | Magee, et al., 1996        |              |   |  |
| BW - Body weight =                                   | kg                          | 70  | USEPA 1995, Region IV      |              |   |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                        | 9125  | USEPA 1991, HHEM           |              |   |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                        | 25550   | USEPA 1991, HHEM           |              |   |  |
| Constituent  | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day  | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) |
| Semivolatiles  |                             |   |                            |              |   |  |
| Benzo(a)anthracene                                   | 8.10E-01                    | 4.76E-07  | NA                         | NA           | 1.70E-07                                | 7.30E-01                               |
| Benzo(a)pyrene                                       | 2.90E-01                    | 1.70E-07  | NA                         | NA           | 6.08E-08                                | 7.30E+00                               |
| Benzo(b)fluoranthene                                 | 3.70E-01                    | 2.17E-07  | NA                         | NA           | 7.76E-08                                | 7.30E-01                               |
| Benzo(k)fluoranthene                                 | 1.60E-01                    | 9.39E-08  | NA                         | NA           | 3.35E-08                                | 7.30E-02                               |
| Carbazole  | 4.90E-01                    | 2.88E-07  | NA                         | NA           | 1.03E-07                                | 2.00E-02                               |
| Chrysene   | 6.10E-01                    | 3.58E-07  | NA                         | NA           | 1.28E-07                                | 7.30E-03                               |
| Dibenz(a,h)anthracene                                | 1.10E-02                    | 6.46E-09  | NA                         | NA           | 2.31E-09                                | 7.30E+00                               |
| Indeno(1,2,3-cd)pyrene                               | 9.40E-02                    | 5.52E-08  | NA                         | NA           | 1.97E-08                                | 7.30E-01                               |
| Naphthalene  | 4.00E-01                    | 2.35E-07  | 2.00E-02                   | 1.17E-05     | 8.39E-08                                | NA                                     |

NA - Not Applicable

Total Hazard Index = 1.17E-05

Total Cancer Risk = 6.61E-07



**Table 85****Dermal Exposure to EU4 Soil (0-20') by a Construction Worker****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | $\frac{\text{Cs} * \text{SA} * \text{AH} * \text{ABS} * \text{EF} * \text{ED} * \text{CF}}{\text{BW} * \text{AT}}$ |                        |  |
|---|----------------------|--|------------------------|--|
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.  |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 5560   | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000  | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 27.8%  | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.1  | USEPA 1997, EFH        |  |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |                      | 0.03   | USEPA 1995, Region III |  |
| ABS <sub>o</sub> - Absorption - other SVOCs =               |                      | 0.1  | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year            | 80   | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                | 1  | reasonable assumption  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06   |                        |  |
| BW - Body weight =  | kg                   | 70   | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 365  | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550  | USEPA 1991, HHEM       |  |

| Constituent            | Concentration<br>in Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Dermal           |                 | Average                         |                                  |                                |
|------------------------|-----------------------------------|--------------------------------------|------------------|-----------------|---------------------------------|----------------------------------|--------------------------------|
|                        |                                   |                                      | RfD<br>mg/kg-day | Hazard<br>Index | Lifetime<br>Intake<br>mg/kg-day | Daily<br>Factor<br>1/(mg/kg-day) | Cancer Slope<br>Cancer<br>Risk |
| <b>Semivolatiles</b>   |                                   |                                      |                  |                 |                                 |                                  |                                |
| Benzo(a)anthracene     | 3.00E+01                          | 1.57E-06                             | NA               | NA              | 2.24E-08                        | 1.46E+00                         | 3.27E-08                       |
| Benzo(a)pyrene         | 1.10E+01                          | 5.74E-07                             | NA               | NA              | 8.21E-09                        | 1.46E+01                         | 1.20E-07                       |
| Benzo(b)fluoranthene   | 1.70E+01                          | 8.88E-07                             | NA               | NA              | 1.27E-08                        | 1.46E+00                         | 1.85E-08                       |
| Benzo(k)fluoranthene   | 6.00E+00                          | 3.13E-07                             | NA               | NA              | 4.48E-09                        | 1.46E-01                         | 6.54E-10                       |
| Carbazole              | 2.40E+01                          | 4.18E-06                             | NA               | NA              | 5.97E-08                        | 2.00E-02                         | 1.19E-09                       |
| Chrysene               | 2.30E+01                          | 1.20E-06                             | NA               | NA              | 1.72E-08                        | 1.46E-02                         | 2.51E-10                       |
| Dibenz(a,h)anthracene  | 1.40E+00                          | 7.31E-08                             | NA               | NA              | 1.04E-09                        | 1.46E+01                         | 1.53E-08                       |
| Indeno(1,2,3-cd)pyrene | 4.90E+00                          | 2.56E-07                             | NA               | NA              | 3.66E-09                        | 1.46E+00                         | 5.34E-09                       |
| Naphthalene            | 2.40E+02                          | 4.18E-05                             | NA               | NA              | 5.97E-07                        | NA                               | NA                             |

NA - Not Available

Total Cancer Risk = 1.94E-07



**Table 86**  
**Oral Exposure to EU4 Soil (0-20') by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

|  |   |           |                       |
|--|---|-----------|-----------------------|
| Intake (mg/kg-day) =                                 | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |           |                       |
| Cd - Concentration in soil =                         | mg/kg   | see below |                       |
| IngR <sub>a</sub> - Ingestion rate for soil =        | mg/day  | 480       | USEPA 1997, EFH       |
| IngR <sub>b</sub> - Ingestion rate for soil =        | mg/day  | 100       | USEPA 1997, EFH       |
| EF <sub>a</sub> - Exposure frequency =               | days/year   | 10        | reasonable assumption |
| EF <sub>b</sub> - Exposure frequency =               | days/year   | 70        | reasonable assumption |
| ED - Exposure duration =                             | years   | 1         | USEPA 1995, Region IV |
| CF - Conversion factor =                             | kg/mg   | 1.00E-06  |                       |
| ME - Matrix effect =                                 |   | 1         | Magee, et al., 1996   |
| BW - Body weight =                                   | kg  | 70        | USEPA 1995, Region IV |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days  | 365       | USEPA 1991, HHEM      |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days  | 25550     | USEPA 1991, HHEM      |

**Exposure Level A**

| Constituent            | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day | Oral Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------|--------------------------------|-------------------------------|--------------|---|--|-------------|
| <b>Semivolatiles</b>   |                             |                                |                               |              |   |  |             |
| Benzo(a)anthracene     | 3.00E+01                    | 5.64E-06                       | NA                            | NA           | 8.05E-08                                | 7.30E-01                               | 5.88E-08    |
| Benzo(a)pyrene         | 1.10E+01                    | 2.07E-06                       | NA                            | NA           | 2.95E-08                                | 7.30E+00                               | 2.16E-07    |
| Benzo(b)fluoranthene   | 1.70E+01                    | 3.19E-06                       | NA                            | NA           | 4.56E-08                                | 7.30E-01                               | 3.33E-08    |
| Benzo(k)fluoranthene   | 6.00E+00                    | 1.13E-06                       | NA                            | NA           | 1.61E-08                                | 7.30E-02                               | 1.18E-09    |
| Carbazole              | 2.40E+01                    | 4.51E-06                       | NA                            | NA           | 6.44E-08                                | 2.00E-02                               | 1.29E-09    |
| Chrysene               | 2.30E+01                    | 4.32E-06                       | NA                            | NA           | 6.17E-08                                | 7.30E-03                               | 4.51E-10    |
| Dibenz(a,h)anthracene  | 1.40E+00                    | 2.63E-07                       | NA                            | NA           | 3.76E-09                                | 7.30E+00                               | 2.74E-08    |
| Indeno(1,2,3-cd)pyrene | 4.90E+00                    | 9.21E-07                       | NA                            | NA           | 1.32E-08                                | 7.30E-01                               | 9.60E-09    |
| Naphthalene            | 2.40E+02                    | 4.51E-05                       | NA                            | NA           | 6.44E-07                                | NA                                     | NA          |

NA - Not Applicable

Cancer Risk = 3.48E-07

**Exposure Level B**

| Constituent            | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day | Oral Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------|--------------------------------|-------------------------------|--------------|---|--|-------------|
| <b>Semivolatiles</b>   |                             |                                |                               |              |   |  |             |
| Benzo(a)anthracene     | 3.00E+01                    | 8.22E-06                       | NA                            | NA           | 1.17E-07                                | 7.30E-01                               | 8.57E-08    |
| Benzo(a)pyrene         | 1.10E+01                    | 3.01E-06                       | NA                            | NA           | 4.31E-08                                | 7.30E+00                               | 3.14E-07    |
| Benzo(b)fluoranthene   | 1.70E+01                    | 4.66E-06                       | NA                            | NA           | 6.65E-08                                | 7.30E-01                               | 4.86E-08    |
| Benzo(k)fluoranthene   | 6.00E+00                    | 1.64E-06                       | NA                            | NA           | 2.35E-08                                | 7.30E-02                               | 1.71E-09    |
| Carbazole              | 2.40E+01                    | 6.58E-06                       | NA                            | NA           | 9.39E-08                                | 2.00E-02                               | 1.88E-09    |
| Chrysene               | 2.30E+01                    | 6.30E-06                       | NA                            | NA           | 9.00E-08                                | 7.30E-03                               | 6.57E-10    |
| Dibenz(a,h)anthracene  | 1.40E+00                    | 3.84E-07                       | NA                            | NA           | 5.48E-09                                | 7.30E+00                               | 4.00E-08    |
| Indeno(1,2,3-cd)pyrene | 4.90E+00                    | 1.34E-06                       | NA                            | NA           | 1.92E-08                                | 7.30E-01                               | 1.40E-08    |
| Naphthalene            | 2.40E+02                    | 6.58E-05                       | NA                            | NA           | 9.39E-07                                | NA                                     | NA          |

NA - Not Applicable

Cancer Risk = 5.07E-07

Total Cancer Risk = 8.54E-07

**Table 8.7**  
**Exposure to Construction Workers from Inhalation of Fugitive Dust in EU4**  
**Preliminary Remediation Goal Calculation**  
**Kerr McGee, Hattiesburg, MS**

| Chemicals              | Concentration in Soil<br>mg/kg | Emission Rate<br>mg/sec | Concentration in Air<br>mg/m <sup>3</sup> | Average Daily Intake<br>mg/kg-day | Inhalation RfD<br>mg/kg-day | Subchronic RfD<br>mg/kg-day | Hazard Index | Average Lifetime Daily Intake<br>mg/kg-day | Inhalation Factor<br>1/(mg/kg-day) | Cancer Slope Factor | Total Cancer Risk: |
|------------------------|--------------------------------|-------------------------|---|-----------------------------------|-----------------------------|-----------------------------|--------------|--|------------------------------------|---------------------|--------------------|
| <b>Semivolatile</b>    |                                |                         |   |                                   |                             |                             |              |  |                                    |                     |                    |
| Benzo(a)anthracene     | 3.00E+01                       | 3.28E-02                | 2.91E-05                                  | 1.37E-06                          | NA                          | NA                          | NA           | 1.95E-08                                   | 3.10E-01                           | 6.05E-09            |                    |
| Benzo(a)pyrene         | 1.10E+01                       | 1.20E-02                | 1.07E-05                                  | 5.01E-07                          | NA                          | NA                          | NA           | 7.16E-09                                   | 3.10E+00                           | 2.22E-08            |                    |
| Benzo(b)fluoranthene   | 1.70E+01                       | 1.86E-02                | 1.65E-05                                  | 7.75E-07                          | NA                          | NA                          | NA           | 1.11E-08                                   | 3.10E-01                           | 3.43E-09            |                    |
| Benzo(k)fluoranthene   | 6.00E+00                       | 6.56E-03                | 5.82E-06                                  | 2.73E-07                          | NA                          | NA                          | NA           | 3.91E-09                                   | 3.10E-02                           | 1.21E-10            |                    |
| Carbazole              | 2.40E+01                       | 2.62E-02                | 2.33E-05                                  | 1.09E-06                          | NA                          | NA                          | NA           | 1.56E-08                                   | NA                                 | NA                  |                    |
| Chrysene               | 2.30E+01                       | 2.51E-02                | 2.23E-05                                  | 1.05E-06                          | NA                          | NA                          | NA           | 1.50E-08                                   | 3.10E-03                           | 4.64E-11            |                    |
| Dibenz(a,h)anthracene  | 1.40E+00                       | 1.53E-03                | 1.36E-06                                  | 6.38E-08                          | NA                          | NA                          | NA           | 9.11E-10                                   | 3.10E+00                           | 2.83E-09            |                    |
| Indeno(1,2,3-cd)pyrene | 4.90E+00                       | 5.36E-03                | 4.75E-06                                  | 2.23E-07                          | NA                          | NA                          | NA           | 3.19E-09                                   | 3.10E-01                           | 9.89E-10            |                    |
| Naphthalene            | 2.40E+02                       | 2.62E-01                | 2.33E-04                                  | 1.09E-05                          | NA                          | NA                          | NA           | 1.56E-07                                   | NA                                 | NA                  |                    |
| NA - Not Available     |                                |                         |   |                                   |                             |                             |              |  |                                    |                     |                    |

Intake (mg/kg-day) =  $\frac{Ca * InhR * EF * ED * RF}{BW * AT}$

Ca - Concentration in air = mg/m<sup>3</sup>  
InhR - Inhalation Rate = m<sup>3</sup>/shift  
EF - Exposure Frequency = shifts/year  
ED - Exposure Duration = years  
RF<sub>s</sub> - Retention Factor - semivolatiles = 0.75  
AT<sub>n</sub> - Averaging Time noncarcinogenic = days  
AT<sub>c</sub> - Averaging Time carcinogenic = days  
BW - Body Weight = kg

z = 6.25[r(Hb/r \* Ln(Hb/r) - 1.58\*Hb/r + 1.58)]

E<sub>i</sub> - Emission Rate (mg/sec) = Cs\*(PERv+PERe)  
Cs - Concentration in soil = mg/kg  
see below

Ca = Concentration in Air (mg/m<sup>3</sup>) = E<sub>i</sub> / (Hb \* W \* V)  
E<sub>i</sub> - Emission Rate of Component (mg/sec) = see below  
Hb - Downwind Ht (m) = 4.81  
W - Width (m) = 50  
V - Wind speed (m/sec) = 4.69  
Length (downwind distance) (m) = 50  
r - Roughness Ht. (m) = 0.20  
z - downwind distance (m) = 50

**Table 88****Oral Exposure to EU5 Surface Soil (0-1') by an Adolescent Visitor (Aged 7-16 years)****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |                             | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                            |              |   |  |             |
|--|-----------------------------|---|----------------------------|--------------|---|--|-------------|
| Cd - Concentration in sediment =                     | mg/kg                       | see below   |                            |              |   |  |             |
| IngR - Ingestion rate for soil =                     | mg/day                      | 100   | USEPA 1997, EFH            |              |   |  |             |
| EF - Exposure frequency =                            | days/year                   | 12  | reasonable assumption      |              |   |  |             |
| ED - Exposure duration =                             | years                       | 10  | USEPA 1995, Region IV      |              |   |  |             |
| CF - Conversion factor =                             | kg/mg                       | 1.00E-06  |                            |              |   |  |             |
| ME - Matrix effect =                                 |                             | 1   | Magee, et al., 1996        |              |   |  |             |
| BW - Body weight =                                   | kg                          | 45  | USEPA 1995, Region IV      |              |   |  |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                        | 3650  | USEPA 1991, HHEM           |              |   |  |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                        | 25550   | USEPA 1991, HHEM           |              |   |  |             |
| Constituent  | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day  | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles  |                             |   |                            |              |   |  |             |
| Benzo(a)anthracene                                   | 2.90E-01                    | 2.12E-08  | NA                         | NA           | 3.03E-09                                | 7.30E-01                               | 2.21E-09    |
| Benzo(a)pyrene                                       | 3.70E-01                    | 2.70E-08  | NA                         | NA           | 3.86E-09                                | 7.30E+00                               | 2.82E-08    |
| Benzo(b)fluoranthene                                 | 7.60E-01                    | 5.55E-08  | NA                         | NA           | 7.93E-09                                | 7.30E-01                               | 5.79E-09    |
| Benzo(k)fluoranthene                                 | 4.60E-01                    | 3.36E-08  | NA                         | NA           | 4.80E-09                                | 7.30E-02                               | 3.50E-10    |
| Chrysene   | 3.70E-01                    | 2.70E-08  | NA                         | NA           | 3.86E-09                                | 7.30E-03                               | 2.82E-11    |
| Dibenz(a,h)anthracene                                | 6.60E-02                    | 4.82E-09  | NA                         | NA           | 6.89E-10                                | 7.30E+00                               | 5.03E-09    |
| Indeno(1,2,3-cd)pyrene                               | 2.90E-01                    | 2.12E-08  | NA                         | NA           | 3.03E-09                                | 7.30E-01                               | 2.21E-09    |

NA - Not Applicable

Total Cancer Risk = 4.38E-08



**Table 89****Dermal Exposure to EU5 Surface Soil (0-6') by a Maintenance Worker****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                |                              |              |   |                                   |             |
|---|---|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| Cs - Concentration in soil =                                | mg/kg   | chem. spec.                    |                              |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                                | 3000                           | calculated                   |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                                     | 20000                          | USEPA 1997, EFH              |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |   | 15%                            | USEPA 1997, EFH              |              |   |                                   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                                  | 0.038                          | USEPA 1997, EFH              |              |   |                                   |             |
| ABS <sub>p</sub> - Absorption - cPAHs =                     |   | 0.03                           | USEPA 1995, Region III       |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year   | 150                            | reasonable assumption        |              |   |                                   |             |
| ED - Exposure duration =                                    | years   | 25                             | USEPA 1995, Region IV        |              |   |                                   |             |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                       |                              |              |   |                                   |             |
| BW - Body weight =  | kg  | 70                             | USEPA 1995, Region IV        |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 9125                           | USEPA 1991, HHEM             |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                          | USEPA 1991, HHEM             |              |   |                                   |             |
| Constituent   | Concentration in Soil mg/kg                         | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |   |                                |                              |              |   |                                   |             |
| Benzo(a)anthracene  | 9.54E-02  | 1.91E-09                       | NA                           | NA           | 6.84E-10                                | 1.46E+00                          | 9.98E-10    |
| Benzo(a)pyrene  | 1.18E-01  | 2.36E-09                       | NA                           | NA           | 8.43E-10                                | 1.46E+01                          | 1.23E-08    |
| Benzo(b)fluoranthene  | 2.90E-01  | 5.82E-09                       | NA                           | NA           | 2.08E-09                                | 1.46E+00                          | 3.03E-09    |
| Benzo(k)fluoranthene  | 1.55E-01  | 3.10E-09                       | NA                           | NA           | 1.11E-09                                | 1.46E-01                          | 1.62E-10    |
| Chrysene  | 1.50E-01  | 3.01E-09                       | NA                           | NA           | 1.07E-09                                | 1.46E-02                          | 1.57E-11    |
| Dibenz(a,h)anthracene                                       | 4.40E-02  | 8.83E-10                       | NA                           | NA           | 3.15E-10                                | 1.46E+01                          | 4.60E-09    |
| Indeno(1,2,3-cd)pyrene                                      | 9.13E-02  | 1.83E-09                       | NA                           | NA           | 6.54E-10                                | 1.46E+00                          | 9.55E-10    |

NA - Not Available

Total Cancer Risk = 2.21E-08



**Table 90****Oral Exposure to EU5 Surface Soil (0-6') by a Maintenance Worker****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                                      |                                  |                 |  |  |                |
|--|---|--------------------------------------|----------------------------------|-----------------|--|--|----------------|
| Cd - Concentration in sediment =                     | mg/kg   | see below                            |                                  |                 |  |  |                |
| IngR - Ingestion rate for soil =                     | mg/day  | 100                                  | USEPA 1997, EFH                  |                 |  |  |                |
| EF - Exposure frequency =                            | days/year   | 150                                  | reasonable assumption            |                 |  |  |                |
| ED - Exposure duration =                             | years   | 25                                   | USEPA 1995, Region IV            |                 |  |  |                |
| CF - Conversion factor =                             | kg/mg   | 1.00E-06                             |                                  |                 |  |  |                |
| ME - Matrix effect =                                 |   | 1                                    | Magee, et al., 1996              |                 |  |  |                |
| BW - Body weight =                                   | kg  | 70                                   | USEPA 1995, Region IV            |                 |  |  |                |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days  | 9125                                 | USEPA 1991, HHEM                 |                 |  |  |                |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days  | 25550                                | USEPA 1991, HHEM                 |                 |  |  |                |
|  |   |                                      |                                  |                 |  |  |                |
| Constituent  | Concentration<br>in Soil<br>mg/kg   | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
| Semivolatiles  |   |                                      |                                  |                 |  |  |                |
| Benzo(a)anthracene                                   | 9.54E-02  | 5.60E-08                             | NA                               | NA              | 2.00E-08   | 7.30E-01                                     | 1.46E-08       |
| Benzo(a)pyrene                                       | 1.18E-01  | 6.91E-08                             | NA                               | NA              | 2.47E-08   | 7.30E+00                                     | 1.80E-07       |
| Benzo(b)fluoranthene                                 | 2.90E-01  | 1.70E-07                             | NA                               | NA              | 6.08E-08   | 7.30E-01                                     | 4.44E-08       |
| Benzo(k)fluoranthene                                 | 1.55E-01  | 9.08E-08                             | NA                               | NA              | 3.24E-08   | 7.30E-02                                     | 2.37E-09       |
| Chrysene   | 1.50E-01  | 8.79E-08                             | NA                               | NA              | 3.14E-08   | 7.30E-03                                     | 2.29E-10       |
| Dibenz(a,h)anthracene                                | 4.40E-02  | 2.58E-08                             | NA                               | NA              | 9.22E-09   | 7.30E+00                                     | 6.73E-08       |
| Indeno(1,2,3-cd)pyrene                               | 9.13E-02  | 5.36E-08                             | NA                               | NA              | 1.91E-08   | 7.30E-01                                     | 1.40E-08       |

NA - Not Applicable

Total Cancer Risk = 3.23E-07



**Table 91****Oral Exposure to EU5 Soil (0-20') by a Construction Worker****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 | Cd*IngR*EF*ED*CF*ME<br>BW*AT |           |                       |
|--|------------------------------|-----------|-----------------------|
| Cd - Concentration in sediment =                     | mg/kg                        | see below |                       |
| IngR <sub>a</sub> - Ingestion rate for soil =        | mg/day                       | 480       | USEPA 1997, EFH       |
| IngR <sub>b</sub> - Ingestion rate for soil =        | mg/day                       | 100       | USEPA 1997, EFH       |
| EF <sub>a</sub> - Exposure frequency =               | days/year                    | 10        | reasonable assumption |
| EF <sub>b</sub> - Exposure frequency =               | days/year                    | 70        | reasonable assumption |
| ED - Exposure duration =                             | years                        | 1         | USEPA 1995, Region IV |
| CF - Conversion factor =                             | kg/mg                        | 1.00E-06  |                       |
| ME - Matrix effect =                                 |                              | 1         | Magee, et al., 1996   |
| BW - Body weight =                                   | kg                           | 70        | USEPA 1995, Region IV |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                         | 365       | USEPA 1991, HHEM      |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                         | 25550     | USEPA 1991, HHEM      |

**Exposure Level A**

| Constituent            | Concentration in | Average Daily    | Oral Chronic  | Hazard Index | Average Lifetime | Daily                | Oral Cancer Slope |
|------------------------|------------------|------------------|---------------|--------------|------------------|----------------------|-------------------|
|                        | Soil mg/kg       | Intake mg/kg-day | RfD mg/kg-day |              | Intake mg/kg-day | Factor 1/(mg/kg-day) | Cancer Risk       |
| <b>Semivolatiles</b>   |                  |                  |               |              |                  |                      |                   |
| Benzo(a)anthracene     | 1.93E-01         | 3.62E-08         | NA            | NA           | 5.18E-10         | 7.30E-01             | 3.78E-10          |
| Benzo(a)pyrene         | 1.91E-01         | 3.59E-08         | NA            | NA           | 5.13E-10         | 7.30E+00             | 3.74E-09          |
| Benzo(b)fluoranthene   | 3.89E-01         | 7.30E-08         | NA            | NA           | 1.04E-09         | 7.30E-01             | 7.61E-10          |
| Benzo(k)fluoranthene   | 1.90E-01         | 3.58E-08         | NA            | NA           | 5.11E-10         | 7.30E-02             | 3.73E-11          |
| Chrysene               | 2.64E-01         | 4.95E-08         | NA            | NA           | 7.07E-10         | 7.30E-03             | 5.16E-12          |
| Dibenz(a,h)anthracene  | 5.15E-02         | 9.68E-09         | NA            | NA           | 1.38E-10         | 7.30E+00             | 1.01E-09          |
| Fluorene               | 2.52E-01         | 4.73E-08         | 4.00E-02      | 1.18E-06     | 6.76E-10         | NA                   | NA                |
| Indeno(1,2,3-cd)pyrene | 1.30E-01         | 2.45E-08         | NA            | NA           | 3.50E-10         | 7.30E-01             | 2.56E-10          |

NA - Not Applicable

Hazard Index = 1.18E-06

Cancer Risk = 6.19E-09

**Exposure Level B**

| Constituent            | Concentration in | Average Daily    | Oral Chronic  | Hazard Index | Average Lifetime | Daily                | Oral Cancer Slope |
|------------------------|------------------|------------------|---------------|--------------|------------------|----------------------|-------------------|
|                        | Soil mg/kg       | Intake mg/kg-day | RfD mg/kg-day |              | Intake mg/kg-day | Factor 1/(mg/kg-day) | Cancer Risk       |
| <b>Semivolatiles</b>   |                  |                  |               |              |                  |                      |                   |
| Benzo(a)anthracene     | 1.93E-01         | 5.28E-08         | NA            | NA           | 7.55E-10         | 7.30E-01             | 5.51E-10          |
| Benzo(a)pyrene         | 1.91E-01         | 5.23E-08         | NA            | NA           | 7.47E-10         | 7.30E+00             | 5.46E-09          |
| Benzo(b)fluoranthene   | 3.89E-01         | 1.06E-07         | NA            | NA           | 1.52E-09         | 7.30E-01             | 1.11E-09          |
| Benzo(k)fluoranthene   | 1.90E-01         | 5.22E-08         | NA            | NA           | 7.45E-10         | 7.30E-02             | 5.44E-11          |
| Chrysene               | 2.64E-01         | 7.22E-08         | NA            | NA           | 1.03E-09         | 7.30E-03             | 7.53E-12          |
| Dibenz(a,h)anthracene  | 5.15E-02         | 1.41E-08         | NA            | NA           | 2.02E-10         | 7.30E+00             | 1.47E-09          |
| Fluorene               | 2.52E-01         | 6.90E-08         | 4.00E-02      | 1.73E-06     | 9.86E-10         | NA                   | NA                |
| Indeno(1,2,3-cd)pyrene | 1.30E-01         | 3.57E-08         | NA            | NA           | 5.11E-10         | 7.30E-01             | 3.73E-10          |

NA - Not Applicable

Hazard Index = 1.73E-06

Cancer Risk = 9.02E-09

Total Hazard Index = 2.91E-06

Total Cancer Risk = 1.52E-08



**Table 92****Dermal Exposure to EU6 Sediment by an Adult Resident (Aged 7 to 30 years)****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =       |                                 | $\frac{Cs*SA*AH*ABS*EF*ED*CF}{BW*AT}$ |                              |              |   |                                   |             |
|----------------------------|---------------------------------|---------------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| Constituent                | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day        | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles              |                                 |                                       |                              |              |   |                                   |             |
| 2-Nitroaniline             | 4.20E-02                        | 1.34E-08                              | NA                           | NA           | 4.60E-09                                | NA                                | NA          |
| 2-Nitrophenol              | 8.40E-02                        | 2.68E-08                              | NA                           | NA           | 9.20E-09                                | NA                                | NA          |
| 3-Nitroaniline             | 8.40E-02                        | 2.68E-08                              | NA                           | NA           | 9.20E-09                                | NA                                | NA          |
| 4-Bromophenylphenylether   | 8.40E-02                        | 2.68E-08                              | NA                           | NA           | 9.20E-09                                | NA                                | NA          |
| 4-Chloro-3-methylphenol    | 8.40E-02                        | 2.68E-08                              | NA                           | NA           | 9.20E-09                                | NA                                | NA          |
| 4-Chlorophenylphenylether  | 4.20E-02                        | 1.34E-08                              | NA                           | NA           | 4.60E-09                                | NA                                | NA          |
| 4-Nitroaniline             | 8.40E-02                        | 2.68E-08                              | NA                           | NA           | 9.20E-09                                | NA                                | NA          |
| Benzo(a)anthracene         | 9.30E-01                        | 8.91E-08                              | NA                           | NA           | 3.05E-08                                | 1.46E+00                          | 4.46E-08    |
| Benzo(a)pyrene             | 9.70E-01                        | 9.29E-08                              | NA                           | NA           | 3.19E-08                                | 1.46E+01                          | 4.65E-07    |
| Benzo(b)fluoranthene       | 1.40E+00                        | 1.34E-07                              | NA                           | NA           | 4.60E-08                                | 1.46E+00                          | 6.71E-08    |
| Benzo(k)fluroanthene       | 5.00E-01                        | 4.79E-08                              | NA                           | NA           | 1.64E-08                                | 1.46E-01                          | 2.40E-09    |
| Bis(2-chloroethoxy)methane | 8.40E-02                        | 2.68E-08                              | NA                           | NA           | 9.20E-09                                | NA                                | NA          |
| Bis(2-chloroethyl)ether    | 4.20E-02                        | 1.34E-08                              | NA                           | NA           | 4.60E-09                                | 1.10E+00                          | 5.06E-09    |
| Carbazole                  | 2.20E-01                        | 7.02E-08                              | NA                           | NA           | 2.41E-08                                | 2.00E-02                          | 4.82E-10    |
| Chrysene                   | 1.30E+00                        | 1.25E-07                              | NA                           | NA           | 4.27E-08                                | 1.46E-02                          | 6.23E-10    |
| Dibenz(a,h)anthracene      | 1.50E-01                        | 1.44E-08                              | NA                           | NA           | 4.93E-09                                | 1.46E+01                          | 7.19E-08    |
| Hexachlorobenzene          | 4.20E-02                        | 1.34E-08                              | NA                           | NA           | 4.60E-09                                | 1.60E+00                          | 7.36E-09    |
| Hexachlorocyclopentadiene  | 2.10E-01                        | 6.70E-08                              | NA                           | NA           | 2.30E-08                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene     | 5.40E-01                        | 5.17E-08                              | NA                           | NA           | 1.77E-08                                | 1.46E+00                          | 2.59E-08    |
| N-nitrosodi-n-propylamine  | 4.20E-02                        | 1.34E-08                              | NA                           | NA           | 4.60E-09                                | 7.00E+00                          | 3.22E-08    |

NA - Not Available

Total Cancer Risk = 7.23E-07



**Table 93****Oral Exposure to EU6 Sediment by an Adult Resident (Aged 7 to 30 years)****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =       |                                 | <u><math>\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}</math></u> |                             |              |   |  |             |
|----------------------------|---------------------------------|---|-----------------------------|--------------|---|--|-------------|
| Constituent                | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day  | Oral Chronic Rf/D mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles              |                                 |   |                             |              |   |  |             |
| 2-Nitroaniline             | 4.20E-02                        | 6.58E-09  | NA                          | NA           | 2.25E-09                                | NA                                     | NA          |
| 2-Nitrophenol              | 8.40E-02                        | 1.32E-08  | NA                          | NA           | 4.51E-09                                | NA                                     | NA          |
| 3-Nitroaniline             | 8.40E-02                        | 1.32E-08  | NA                          | NA           | 4.51E-09                                | NA                                     | NA          |
| 4-Bromophenylphenoxyether  | 8.40E-02                        | 1.32E-08  | NA                          | NA           | 4.51E-09                                | NA                                     | NA          |
| 4-Chloro-3-methylphenol    | 8.40E-02                        | 1.32E-08  | NA                          | NA           | 4.51E-09                                | NA                                     | NA          |
| 4-Chlorophenylphenoxyether | 4.20E-02                        | 6.58E-09  | NA                          | NA           | 2.25E-09                                | NA                                     | NA          |
| 4-Nitroaniline             | 8.40E-02                        | 1.32E-08  | NA                          | NA           | 4.51E-09                                | NA                                     | NA          |
| Benzo(a)anthracene         | 9.30E-01                        | 1.46E-07  | NA                          | NA           | 4.99E-08                                | 7.30E-01                               | 3.64E-08    |
| Benzo(a)pyrene             | 9.70E-01                        | 1.52E-07  | NA                          | NA           | 5.21E-08                                | 7.30E+00                               | 3.80E-07    |
| Benzo(b)fluoranthene       | 1.40E+00                        | 2.19E-07  | NA                          | NA           | 7.51E-08                                | 7.30E-01                               | 5.49E-08    |
| Benzo(k)fluoranthene       | 5.00E-01                        | 7.83E-08  | NA                          | NA           | 2.68E-08                                | 7.30E-02                               | 1.96E-09    |
| Bis(2-chloroethoxy)methane | 8.40E-02                        | 1.32E-08  | NA                          | NA           | 4.51E-09                                | NA                                     | NA          |
| Bis(2-chloroethyl)ether    | 4.20E-02                        | 6.58E-09  | NA                          | NA           | 2.25E-09                                | 1.10E+00                               | 2.48E-09    |
| Carbazole                  | 2.20E-01                        | 3.44E-08  | NA                          | NA           | 1.18E-08                                | 2.00E-02                               | 2.36E-10    |
| Chrysene                   | 1.30E+00                        | 2.04E-07  | NA                          | NA           | 6.98E-08                                | 7.30E-03                               | 5.09E-10    |
| Dibenz(a,h)anthracene      | 1.50E-01                        | 2.35E-08  | NA                          | NA           | 8.05E-09                                | 7.30E+00                               | 5.88E-08    |
| Hexachlorobenzene          | 4.20E-02                        | 6.58E-09  | 8.00E-04                    | 8.22E-06     | 2.25E-09                                | 1.60E+00                               | 3.61E-09    |
| Hexachlorocyclopentadiene  | 2.10E-01                        | 3.29E-08  | 7.00E-03                    | 4.70E-06     | 1.13E-08                                | NA                                     | NA          |
| Indeno(1,2,3-cd)pyrene     | 5.40E-01                        | 8.45E-08  | NA                          | NA           | 2.90E-08                                | 7.30E-01                               | 2.12E-08    |
| N-nitrosodi-n-propylamine  | 4.20E-02                        | 6.58E-09  | NA                          | NA           | 2.25E-09                                | 7.00E+00                               | 1.58E-08    |

NA - Not Applicable

Total Hazard Index = 1.29E-05

Total Cancer Risk = 5.76E-07



**Table 94****Dermal Exposure to EU6 Sediment by a Child Resident (Aged 1 to 6 years)****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =       |                                 | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                          |              |   |                                   |             |
|----------------------------|---------------------------------|---|--------------------------|--------------|---|-----------------------------------|-------------|
| Constituent                | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day                      | Subchronic RID mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles              |                                 |   |                          |              |   |                                   |             |
| 2-Nitroaniline             | 4.20E-02                        | 2.26E-08  | NA                       | NA           | 1.93E-09                                | NA                                | NA          |
| 2-Nitrophenol              | 8.40E-02                        | 4.51E-08  | NA                       | NA           | 3.87E-09                                | NA                                | NA          |
| 3-Nitroaniline             | 8.40E-02                        | 4.51E-08  | NA                       | NA           | 3.87E-09                                | NA                                | NA          |
| 4-Bromophenylphenoxyether  | 8.40E-02                        | 4.51E-08  | NA                       | NA           | 3.87E-09                                | NA                                | NA          |
| 4-Chloro-3-methylphenol    | 8.40E-02                        | 4.51E-08  | NA                       | NA           | 3.87E-09                                | NA                                | NA          |
| 4-Chlorophenylphenoxyether | 4.20E-02                        | 2.26E-08  | NA                       | NA           | 1.93E-09                                | NA                                | NA          |
| 4-Nitroaniline             | 8.40E-02                        | 4.51E-08  | NA                       | NA           | 3.87E-09                                | NA                                | NA          |
| Benzo(a)anthracene         | 9.30E-01                        | 1.50E-07  | NA                       | NA           | 1.29E-08                                | 1.46E+00                          | 1.88E-08    |
| Benzo(a)pyrene             | 9.70E-01                        | 1.56E-07  | NA                       | NA           | 1.34E-08                                | 1.46E+01                          | 1.96E-07    |
| Benzo(b)fluoranthene       | 1.40E+00                        | 2.26E-07  | NA                       | NA           | 1.93E-08                                | 1.46E+00                          | 2.82E-08    |
| Benzo(k)fluoranthene       | 5.00E-01                        | 8.06E-08  | NA                       | NA           | 6.91E-09                                | 1.46E-01                          | 1.01E-09    |
| Bis(2-chloroethoxy)methane | 8.40E-02                        | 4.51E-08  | NA                       | NA           | 3.87E-09                                | NA                                | NA          |
| Bis(2-chloroethyl)ether    | 4.20E-02                        | 2.26E-08  | NA                       | NA           | 1.93E-09                                | 1.10E+00                          | 2.13E-09    |
| Carbazole                  | 2.20E-01                        | 1.18E-07  | NA                       | NA           | 1.01E-08                                | 2.00E-02                          | 2.03E-10    |
| Chrysene                   | 1.30E+00                        | 2.10E-07  | NA                       | NA           | 1.80E-08                                | 1.46E-02                          |             |
| Dibenz(a,h)anthracene      | 1.50E-01                        | 2.42E-08  | NA                       | NA           | 2.07E-09                                | 1.46E+01                          | 3.03E-08    |
| Hexachlorobenzene          | 4.20E-02                        | 2.26E-08  | NA                       | NA           | 1.93E-09                                | 1.60E+00                          | 3.10E-09    |
| Hexachlorocyclopentadiene  | 2.10E-01                        | 1.13E-07  | NA                       | NA           | 9.67E-09                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene     | 5.40E-01                        | 8.71E-08  | NA                       | NA           | 7.46E-09                                | 1.46E+00                          | 1.09E-08    |
| N-nitrosodi-n-propylamine  | 4.20E-02                        | 2.26E-08  | NA                       | NA           | 1.93E-09                                | 7.00E+00                          | 1.35E-08    |

NA - Not Available

Total Cancer Risk = 3.04E-07



**Table 95****Oral Exposure to EU6 Sediment by a Child Resident (Aged 1 to 6 years)****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

|  |  | Intake (mg/kg-day) =             | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                               |              |   |  |             |
|--|--|----------------------------------|---|-------------------------------|--------------|---|--|-------------|
|  |  | Cd - Concentration in sediment = | mg/kg   | see below                     |              |   |  |             |
| IngR - Ingestion rate for sediment =                 |  | mg/day                           | 200   | USEPA 1997, EFH               |              |   |  |             |
| EP - Exposure frequency =                            |  | days/year                        | 40  | reasonable assumption         |              |   |  |             |
| ED - Exposure duration =                             |  | years                            | 6   | USEPA 1995, Region IV         |              |   |  |             |
| CF - Conversion factor =                             |  | kg/mg                            | 1.00E-06  |                               |              |   |  |             |
| ME - Matrix effect =                                 |  |                                  | 1   | Magee, et al., 1996           |              |   |  |             |
| BW - Body weight =                                   |  | kg                               | 15  | USEPA 1995, Region IV         |              |   |  |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = |  | days                             | 2190  | USEPA 1991, HHEM              |              |   |  |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =    |  | days                             | 25550   | USEPA 1991, HHEM              |              |   |  |             |
| Constituent  |  | Concentration in Sediment mg/kg  | Average Daily Intake mg/kg-day  | Oral Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>                                 |  |                                  |   |                               |              |   |  |             |
| 2-Nitroaniline                                       |  | 4.20E-02                         | 6.14E-08  | NA                            | NA           | 5.26E-09                                | NA                                     | NA          |
| 2-Nitrophenol  |  | 8.40E-02                         | 1.23E-07  | NA                            | NA           | 1.05E-08                                | NA                                     | NA          |
| 3-Nitroaniline                                       |  | 8.40E-02                         | 1.23E-07  | NA                            | NA           | 1.05E-08                                | NA                                     | NA          |
| 4-Bromophenylphenylether                             |  | 8.40E-02                         | 1.23E-07  | NA                            | NA           | 1.05E-08                                | NA                                     | NA          |
| 4-Chloro-3-methylphenol                              |  | 8.40E-02                         | 1.23E-07  | NA                            | NA           | 1.05E-08                                | NA                                     | NA          |
| 4-Chlorophenylphenylether                            |  | 4.20E-02                         | 6.14E-08  | NA                            | NA           | 5.26E-09                                | NA                                     | NA          |
| 4-Nitroaniline                                       |  | 8.40E-02                         | 1.23E-07  | NA                            | NA           | 1.05E-08                                | NA                                     | NA          |
| Benzo(a)anthracene                                   |  | 9.30E-01                         | 1.36E-06  | NA                            | NA           | 1.16E-07                                | 7.30E-01                               | 8.50E-08    |
| Benzo(a)pyrene                                       |  | 9.70E-01                         | 1.42E-06  | NA                            | NA           | 1.21E-07                                | 7.30E+00                               | 8.87E-07    |
| Benzo(b)fluoranthene                                 |  | 1.40E+00                         | 2.05E-06  | NA                            | NA           | 1.75E-07                                | 7.30E-01                               | 1.28E-07    |
| Benzo(k)fluroanthene                                 |  | 5.00E-01                         | 7.31E-07  | NA                            | NA           | 6.26E-08                                | 7.30E-02                               | 4.57E-09    |
| Bis(2-chloroethoxy)methane                           |  | 8.40E-02                         | 1.23E-07  | NA                            | NA           | 1.05E-08                                | NA                                     | NA          |
| Bis(2-chloroethyl)ether                              |  | 4.20E-02                         | 6.14E-08  | NA                            | NA           | 5.26E-09                                | 1.10E+00                               | 5.79E-09    |
| Carbazole  |  | 2.20E-01                         | 3.21E-07  | NA                            | NA           | 2.76E-08                                | 2.00E-02                               | 5.51E-10    |
| Chrysene   |  | 1.30E+00                         | 1.90E-06  | NA                            | NA           | 1.63E-07                                | 7.30E-03                               | 1.19E-09    |
| Dibenz(a,h)anthracene                                |  | 1.50E-01                         | 2.19E-07  | NA                            | NA           | 1.88E-08                                | 7.30E+00                               | 1.37E-07    |
| Hexachlorobenzene                                    |  | 4.20E-02                         | 6.14E-08  | NA                            | NA           | 5.26E-09                                | 1.60E+00                               | 8.42E-09    |
| Hexachlorocyclopentadiene                            |  | 2.10E-01                         | 3.07E-07  | NA                            | NA           | 2.63E-08                                | NA                                     | NA          |
| Indeno(1,2,3-cd)pyrene                               |  | 5.40E-01                         | 7.89E-07  | NA                            | NA           | 6.76E-08                                | 7.30E-01                               | 4.94E-08    |
| N-nitrosodi-n-propylamine                            |  | 4.20E-02                         | 6.14E-08  | NA                            | NA           | 5.26E-09                                | 7.00E+00                               | 3.68E-08    |

NA - Not Applicable

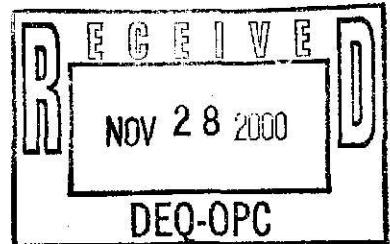
Total Cancer Risk = 1.34E-06







Setting the Standards for Innovative  
Environmental Solutions



**HUMAN HEALTH RISK ASSESSMENT  
FOR THE FORMER GULF STATES CREOSOTING FACILITY,  
HATTIESBURG, MISSISSIPPI**

November 22, 2000

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## Executive Summary

A baseline human health risk assessment (HHRA) was conducted for the Former Gulf States Creosoting facility in Hattiesburg, Mississippi. The HHRA was performed in accordance with: Mississippi Commission on Environmental Quality's (MCEQ's) *Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi* (1999); US EPA's *Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A)* (1989); US EPA Region 4 guidance entitled *Technical Services Supplemental Guidance to RAGS, Region 4 Bulletins* (1995); and other relevant US EPA guidance documents.

Creosoting constituents of potential health concern include polycyclic aromatic hydrocarbons, of which benzo(a)pyrene is the predominant contributor to potential risks. Much of the former creosoting process area is currently covered with asphalt or large building structures. Potential future exposure scenarios included a construction worker, a maintenance worker, an infrequent Site visitor, and off-Site residents. Media of concern included soils, sediment, and surface water.

Hazards posed by chemical constituents in soil for health effects other than an increased risk of cancer were well below a threshold of possible concern for each receptor evaluated in this risk assessment. Cancer risks were within the US EPA's acceptable target risk range of  $1 \times 10^{-6}$  to  $1 \times 10^{-4}$  (*i.e.*, one in one million to one in ten thousand). The added lifetime cancer risk conservatively estimated for a maintenance worker was  $1 \times 10^{-5}$  (*i.e.*, an upper bound risk of one incidence of cancer out of a population of 100,000 persons so exposed). The potential risk for a construction worker was estimated to be  $9 \times 10^{-7}$ . The estimated potential risk for an adolescent Site visitor was  $3 \times 10^{-6}$ . Lastly, the estimated risk level for an off-Site resident was  $5 \times 10^{-6}$ . While the total risk levels for adolescent visitor and off-Site resident scenarios exceeded the *de minimis* benchmark of  $1 \times 10^{-6}$ , no risk levels for any individual carcinogen exceeded the  $1 \times 10^{-6}$  for these receptors. As such, these scenarios maintain acceptable levels of risk according to MCEQ guidance (1999).

The maintenance worker risk level of  $1 \times 10^{-5}$  was attributable to sediment and surface soil exposures in EU 4 and surface soil exposures in EU 5. Dermal exposures to benzo(a)pyrene at sample locations SD-02 (130 mg/kg) and SD-12 (71 mg/kg) drove the maintenance worker risk levels for exposure to sediment in EU4. Implementing a remedy to remove, treat, or preclude contact with sediment at sample locations SD-02 and SD-12 will result in acceptable risk levels ( $1 \times 10^{-7}$ ) for sediment exposures to a maintenance worker in EU 4.

Dermal exposure to benzo(a)pyrene at sample locations GEO-48/0-1 (500 mg/kg), GEO-21/0-1 (230 mg/kg), and GEO-21/2-3 (190 mg/kg) drove the maintenance worker risk levels for exposure to surface soils (0 to 6 feet below ground surface) in EU 4. Implementing a remedy to remove, treat, or preclude contact with soils at sample locations GEO-48/0-1, GEO-21/0-1, and GEO-21/2-3 will result in acceptable risk levels ( $6 \times 10^{-7}$ ) for surface soil exposures to a maintenance worker in EU 4.

Dermal exposures to benzo(a)pyrene at sample locations GEO-33 0-1 (52.5 mg/kg) drove the maintenance worker risk levels for exposures to surface soils (0 to 6 feet below ground surface) in EU 5. Precluding direct contact with sample location GEO-33/0-1 will result in acceptable risk levels ( $4 \times 10^{-7}$ ) for surface soil exposures to a maintenance worker in EU 5. The existing pavement overlying sample location GEO-33 currently serves this purpose. Proper maintenance of the existing pavement in the vicinity of GEO-33 will ensure that direct contact with affected soils does not occur in that area. Consequently, a complete pathway between the GEO-33 sample location and potential human receptors will continue to be nonexistent thereby resulting in an acceptable level of risk to maintenance workers in EU 5.

## **1.0 Introduction**

Environmental Standards, Inc. (Environmental Standards) was retained by Kerr-McGee Chemical Corporation (Kerr-McGee) to perform a human health risk assessment (HHRA) to evaluate hazards and risks potentially posed by residual levels of chemicals present at the Former Gulf States Creosoting facility (Site). The Site, located near the intersection of US Highways 49 and 11 in Hattiesburg, Mississippi, was formerly a wood treating facility that operated between the early 1900s and 1960. In the early 1960s, the Site was redeveloped for commercial and light industrial uses (Michael Pisani & Assoc., 1997). The land on which the Site is located is a portion of the Sixteenth Section land owned by the Hattiesburg Public School District and leased to the current tenants under a 99-year lease, granted on July 7, 1947. At the time of this report, the Site, with the exception of the grassy and wooded areas in the south and southwest, respectively, was primarily used for automobile dealerships. There are no residential or institutional (*i.e.*, schools) uses of the Site (Michael Pisani & Assoc., 1997).

Operations at the Site consisted of a small-scale wood preserving process using creosote. The creosoting process was primarily confined to a 2.5-acre area in the northeast corner of the Site; this is known as the Process Area and is currently occupied by Courtesy Ford. During the redevelopment of the Site in the early 1960s, construction debris (*e.g.*, broken concrete, asphalt, etc.) appears to have been relocated to the southwestern corner of the Site along Gordon's Creek. This area is known as the Fill Area and currently remains undeveloped.

This assessment has been conducted as a result of an agreement between Kerr-McGee, the Mississippi Department of Environmental Quality (MDEQ), and the Mississippi Commission on Environmental Quality (MCEQ) pursuant to the Uncontrolled Site Voluntary Evaluation Program. The MDEQ Office of Pollution Control, Uncontrolled Sites Section has been providing oversight and review of investigations and reports relating to the former Gulf States Creosoting facility.

This report will address the potential for on-Site exposures to human receptors and off-Site exposures to humans along the northeast drainage ditch.

The primary guidance used to develop this risk assessment was the MCEQ *Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi* (1999). US EPA Region 4's *Technical Services Supplemental Guidance to RAGS: Region 4 Bulletins* (1995) were also referred to for guidance. Additional US EPA guidance documents cited herein include:

- *Guidance for Remediation of Uncontrolled Hazardous Substance Sites in Mississippi (MDEQ, 1990);*
- *Risk Assessment Guidance for Superfund, Volume 1, Human Health Evaluation Manual/ Part A (RAGS/Part A)* (US EPA, 1989);
- *Human Health Evaluation Manual, Supplemental Guidance: "Standard Default Exposure Factors"* (US EPA, 1991);
- *Exposure Factors Handbook* (US EPA, 1997);
- *Guidelines for Exposure Assessment* (US EPA 1992);
- *Dermal Exposure Assessment: Principles and Applications* (US EPA, 1992);

These documents are not listed in a hierarchical manner; other US EPA guidance documents and peer-reviewed technical papers may have also been referenced in this risk assessment report.

## **2.0 Hazard Identification and Conceptual Site Model**

As a result of the historical wood preservation process, residual levels of creosote-related chemicals are present in soils in the Process Area. Sediment and surface water in a drainage ditch along the southeast border of the Process Area also contain chemical residuals. These Site-related chemicals, mostly polycyclic aromatic hydrocarbons (PAHs) are also present in the Fill Area. Residual levels of PAHs have been found in soil in the Fill Area and in Gordon's Creek surface water and sediment.

PAH residuals have also been detected in shallow groundwater underlying the Site. Currently, there are no private water wells located on-Site that access this shallow groundwater for potable purposes. The results of a door-to-door survey conducted by Michael Pisani and Associates on October 3, 2000 indicated no private uses of shallow groundwater downgradient of the Site. For these reasons, the groundwater exposure pathway, both on- and off-Site , was considered incomplete and not evaluated in this assessment.

A conceptual site model (CSM) was developed for the Site to aid in determining the potential receptors and exposure units to be evaluated under current and future potential land use (Figure 1). These receptors were identified as infrequent Site visitors, maintenance workers, construction workers, and off-Site residents.

Under current land use assumptions, Site visitors may potentially contact residual chemicals in Gordon's Creek surface water and sediment, and/or surface soils in the Fill Area and surrounding woods, the grassy field southeast of the Fill Area, and/or the drainage ditch along side of the Process Area. Visitors may also potentially contact surface water and sediment in the Process Area drainage ditch and surface soils in the adjacent areas. The remaining affected areas of the Site are covered with either buildings or pavement precluding casual direct contact with surface soils. As a conservative measure, however, visitor exposure to soils from these paved areas was also assessed.

Under both current and future land use assumptions, a maintenance worker may contact surface soils in the Fill Area and surrounding woods, the grassy field southeast of the Fill Area, and/or the Process Area and surrounding affected areas, including the drainage ditch located to the southeast of the Process Area. Although most of the Process Area and vicinity are paved, maintenance activities may involve some shallow digging; therefore, direct contact with shallow soils in this area was assessed. Maintenance activities are not expected to occur in Gordon's Creek; however, as a conservative measure, exposure to surface water and sediment in Gordon's Creek was assessed. The remainder of the Site was relatively unaffected by historical creosoting activities.

Although there are currently no major construction activities at the Site, these types of activities may occur at some time in the future. As with the maintenance worker scenario, construction activities could potentially occur in the Fill Area and vicinity, the grassy field southeast of the Fill Area, and the Process Area and vicinity. Construction workers may be exposed to both surface and subsurface soils (down to the water table). Construction activities are not anticipated to occur in Gordon's Creek; however, construction worker exposure to surface water and sediment in Gordon's Creek was assessed as a conservative measure. The remainder of the Site was relatively unaffected by historical creosoting activities.

Areas of the Site affected by historical creosoting activities will be deed restricted prohibiting future residential development. Off-Site areas along the northeast Drainage Ditch, currently a residential neighborhood, were assessed for residential exposures to soil, sediment, and surface water.

### **3.0 Data Evaluation**

To characterize potential exposures to Site-related chemicals, the former Gulf States Creosoting facility was divided into six exposure units (EUs). Each exposure unit outlines potentially affected areas of the Site and adjacent on-Site locales that may be frequented by individuals accessing the Site for recreational or occupational purposes. The use of EUs is encouraged by the US EPA Region 4 (1995), which defines an EU as “an areal extent of a receptor’s movements during a single day... .” Each of these exposure units is depicted on Figure 2 and is discussed below.

A sixth EU was created for off-Site residential exposures to soil, surface water, and sediment along the northeast Drainage Ditch. This EU is delineated on Figure 3.

#### **3.1 Exposure Unit Delineation**

The following EUs were delineated based upon the presence of residual chemicals and the potential for receptors to contact those chemicals. Areas of the Site most affected were included in at least one of the five EUs while areas with relatively low or non-detectable concentrations of residuals were not included in an EU. By limiting Site-wide exposures to the EUs most affected by historical activities at the Site, worst-case scenarios were created.

##### **3.1.1 Exposure Unit 1**

EU1 outlines the on-Site areas in, adjacent to, and downstream of the Fill Area along Gordon’s Creek (Figure 2). EU1 includes exposures to surface water and sediment by an infrequent Site visitor, future maintenance worker and future construction worker. Although US EPA Region IV guidance indicates that “In most cases it is unnecessary to evaluate human exposures to sediments covered by surface water,” (US EPA, 1995) dermal and oral surface water exposures were conservatively assessed herein at the request of the MDEQ (2000).

Soil samples from this area were considered part of EU2 and exposures were assessed accordingly.

### 3.1.2 Exposure Unit 2

EU2 delineates the upland areas of the Fill Area and adjacent woody and grassy areas (Figure 2). Surface soils from zero to one foot and zero to six feet below ground surface [bgs] in this area were evaluated for potential visitor and future hypothetical maintenance worker scenarios, respectively. Surface and subsurface soils were also evaluated for a hypothetical future construction worker scenario. Available data for subsurface soils for a construction scenario were evaluated from the surface to the water table (approximately 10 feet bgs) as recommended by the MDEQ (2000).

### 3.1.3 Exposure Unit 3

In the southwest corner of the Site there exists a grassy field east of West Pine Street between Henson Auto Sales and Eagan Cars and Trucks. This grassy area has been defined as EU3 for purposes of this risk assessment (Figure 2). Similar to EU2, surface soil from zero to one foot and zero to six feet bgs were evaluated in EU2 for visitor and hypothetical future maintenance worker scenarios, respectively. Surface and subsurface soils in this EU were evaluated for a hypothetical future construction worker scenario. Available data for subsurface soils for a construction scenario were evaluated from the surface to the water table(approximately 20 feet bgs) as recommended by the MDEQ (2000).

### 3.1.4 Exposure Unit 4

EU 4 encompasses the grassy drainage ditch area along the fenceline behind Courtesy Ford in the northeast corner of the Site and continues parallel to the railroad tracks, and west through EU 3 and EU 2 (Figure 2). EU 4 along the southeast side of the Process Area has been widened to include soil data from that area. Receptors associated with EU 4 included Site visitor exposures via casual contact with surface soil, sediment, and surface water. Maintenance worker and construction worker scenarios were also evaluated for exposures to surface water and sediment in EU 4 as well as soils in EU 4 near the Process Area. Soils down to six feet bgs were evaluated for maintenance workers while soils down to the water table (approximately20 feet bgs) were evaluated for construction workers in this EU as requested by the MDEQ (2000).

### 3.1.5 Exposure Unit 5

EU5 outlines the Process Area and the historical drip track areas of the former Gulf States Creosoting facility (Figure 2). Surface soils from zero to six feet bgs were evaluated in EU5 for a hypothetical maintenance worker scenario. Available data for soils down to the water table (approximately 20 feet bgs) were evaluated in EU5 for a hypothetical future construction worker scenario.

### 3.1.6 Exposure Unit 6

EU6 outlines a stretch (approximately 2700 feet in length) of the northeast drainage ditch that leads from the Site into the neighboring residential area. EU6 exposures include oral and dermal exposures by off-Site residents to sediment and surface water along the northeast drainage ditch. Soil exposures were not assessed in this area for lack of soil data. Also, it was anticipated that sediment exposures in this area represent a more conservative estimate of exposure in that chemical concentrations in the exposed sediment along the drainage ditch are likely to be greater than concentrations in the surrounding soils.

## 3.2 Statistical Evaluation

Environmental samples undergo laboratory analyses that are designed to quantitate the concentrations of constituents in the various environmental media. As a result of the analytical procedures, a constituent may be detected and its concentration measured, detected but not able to be quantitated, or not detected at all in a sample. The data set for the Site contains a number of nondetections for some chemicals of potential concern (COPCs) in various samples. Assuming that the COPC is present in these samples at the achieved detection limit is biased because the chemical may be absent altogether. Assuming a concentration of zero is also flawed because the chemical could be present at a level below laboratory capabilities to detect and quantify the concentration. Consequently, in the event that an analyte identified at least once in a given medium was not detected in a given sample, it was conservatively assumed for the risk assessment purposes to be present at a concentration equivalent to one-half of the sample quantitation limit (SQL). In addition, samples labeled with an "R" (rejected) qualifier were not

included in the data analysis because those data were deemed unreliable and, therefore, unusable. Constituents that were not detected in any sample from a particular medium were eliminated from further consideration in accordance with US EPA guidelines (1989).

Site analytical data used in this assessment were collected during the Phase I (1997) and Phase II (1998) remedial investigations as well as the additional investigation conducted in 2000 at the request of the MDEQ. These data were fully validated by qualified technical professionals using standard data validation protocols, as required by the MCEQ (1999).

Previous investigations at the Site have been conducted since 1990. These investigations included the following:

- 1990 soil gas and soil sampling by Roy F. Weston
- 1991 MDEQ Site inspections and Phase II report
- 1994 Phase II Site investigation by Environmental Protection Systems (EPS)
- 1994 Site investigation by Bonner Analytical Testing Company (BATCO)
- 1994 preliminary subsurface investigation by BATCO
- 1995 three-dimension resistivity surveys by American Remediation Technology
- 1996 investigation by McLaren/Hart
- 1996 investigation by Kerr McGee Chemical Corporation

Data acquired from these historical (pre-1997) investigatory activities were not used in this assessment as they were not validated by qualified chemists and sampling locations for some of the data could not be specifically defined. These historical data were not considered valid and were, therefore, not appropriate to use in this assessment of risks. Only validated data were considered to be representative of Site conditions with a reasonable level of confidence and were used for this assessment.

The validated laboratory data from 1997, 1998, and 2000 investigations were compiled into data sets representing areas of potential exposure (EUs) for each potential receptor. Each data set was analyzed statistically using SiteStat®, a commercially available software package, to calculate the minimum, maximum, arithmetic mean, logarithmic mean, standard error of the mean, and the 95% upper confidence limit of the mean concentration (95% UCL) for each constituent based on distributional analysis of the data (*i.e.*, utilizing goodness-of-fit statistical tests to determine whether the data are distributed normally or lognormally). The data qualifier associated with the minimum and maximum detected concentrations as well as the location of the maximum detected concentration for each EU were also determined. Results of the quantitative and statistical analyses for each of the EUs discussed above are presented in Tables 1 through 18.

Standard sampling protocol requires the collection of duplicate field samples used to ensure the quality of a laboratory analysis (*i.e.*, to ensure that analytical results can be replicated). As such, duplicate sample results were provided as part of the database for the Hattiesburg Site. In accordance with US EPA guidance (1989), duplicate sample results were averaged (for any sample containing duplicates) and the average concentration was used as a single concentration for that sample in the calculation of summary statistics as discussed below.

Soils down to one foot deep were assumed to be representative of surface soils at the Site for infrequent visitor exposures. A depth of 0 to 6 feet was used to define surface soils for maintenance worker exposures. These assumptions were recommended by the MDEQ (2000). The groundwater table was considered the extent of subsurface soils as recommended by MDEQ (2000). This value (depth-to-groundwater) varies significantly across the Site and, as such, the extent of subsurface soil was EU-specific as follows:

- EU2 – soils down to 10 feet
- EU3 – soils down to 20 feet
- EU4 – soils down to 20 feet
- EU5 – soils down to 20 feet

### 3.3 Determination of Exposure-Point Concentrations

Exposure-point concentrations were determined to be the 95% UCL or the maximum concentration of a COPC in an EU, whichever was lower. This methodology is in accordance with US EPA guidance (1989). If the distribution of the concentration data was determined to be lognormal, then the lognormal 95% UCL was compared to the maximum concentration to determine the exposure-point concentration. In the event that the distribution of a chemical in any given medium could not be confidently labeled as normal or lognormal, it is termed either "unknown" or "normal/lognormal." In these cases, the lognormal 95% UCL was compared to maximum concentration when determining the exposure-point concentration. It should be noted, however, that in cases where the distribution is "unknown," the normal and lognormal 95% UCLs could not be reliably predicted. Assuming a lognormal distribution of the data increases the uncertainty associated with this step of the risk assessment process; however, hazard and risk estimates are likely to be less uncertain than if the maximum concentrations were used.

Exposure-point concentrations are provided on the statistical summary tables, Tables 1 through 18.

### 3.4 COPC Selection

Soils (both surface and subsurface) were screened according to MCEQ (1999) guidance. The first tier of the screening process compared maximum concentrations of a constituent in an EU with the Restricted Tier 1 target remediation goal (TRG) for maintenance worker and construction worker scenarios. Restricted TRGs were used because the Site is not currently used for residential purposes and the current commercial/industrial land-use is anticipated to remain into the future as a result of the implementation of deed restrictions on the impacted areas of the Site. If a maximum concentration of a constituent was less than the Restricted Tier 1 TRG, then that constituent was eliminated from further quantitative assessment.

Surface soil data (zero to one foot bgs) for the visitor scenario were screened using Unrestricted Tier 1 TRGs at the request of MDEQ (2000). If a maximum concentration of a constituent was less than the Unrestricted Tier 1 TRG, then that constituent was eliminated from further

quantitative assessment. Conversely, if the maximum concentration of a constituent exceeded the Tier 1 TRG, that constituent was retained for quantitative analysis.

If the maximum concentration of a constituent in an EU exceeded the Tier 1 TRG, then the 95% UCL of the constituent was compared to the Tier 1 TRG (Restricted or Unrestricted, depending on the exposure scenarios as described above) as part of the Tier II screening process. In the event that the concentrations of a chemical were distributed lognormally, the lognormal 95% UCL of that constituent was compared to the Tier 1 TRG. If the distribution of data of a chemical could not be positively identified as either normal or lognormal, the lognormal 95% UCL was used in the screening process. In these cases, either the maximum concentration or the lognormal 95% UCL can be conservatively used. The US EPA, however, justifies the use of an average concentration as the exposure-point concentration by explaining that toxicity criteria for both carcinogenic and non-carcinogenic effects are based on lifetime average exposures and that the "average concentration is most representative of the concentration that would be contacted at a site over time" (*Supplemental Guidance to RAGS: Calculating the Concentration Term*, 1992). Other US EPA guidance states that "...in most situations, assuming long-term contact with the maximum concentration is not reasonable" (*Risk Assessment Guidance for Superfund, Part A*, 1989). US EPA Region 4 also states that, generally, it is reasonable to assume that soil data are distributed lognormally (1995). In keeping with these guidances, the lognormal 95% UCL was considered in the screening process where the data distribution for a compound could not be defined as specifically normal or lognormal.

If the 95% UCL (or lognormal 95% UCL where appropriate) of a constituent was less than the Tier 1 TRG, then that constituent was eliminated from further quantitative analysis. If the 95% UCL (or lognormal 95% UCL where appropriate) of a constituent in soil exceeded the Tier 1 TRG, then that constituent was retained for quantitative analysis in the Site-specific risk assessment (Tier III).

MCEQ guidance (1999) does not specify screening levels for constituents in sediment or surface water; therefore, Region 4 was referred to for guidance (1995). Sediment is only found on the Site in drainage ditches that contain little to no water most of the time. US EPA Region 4 guidance states that sediments in an intermittent stream (or ditch) should be considered as surface soil for the portion of the year the stream is without water. Based on these factors and comments provided by the MDEQ (2000), the maximum detected constituent concentrations in sediment was compared to MCEQ unrestricted Tier 1 TRGs. The screening process then followed the same procedure as mentioned above for other soils.

For surface water, the maximum detected concentration of a constituent in an EU was compared to the US EPA Human Health Water Quality Standard (WQS) for consumption of water and organisms in accordance with US EPA Region 4 guidance (1995). If the maximum concentration of a constituent in surface water was less than the WQS, then that constituent was eliminated from quantitative analysis. If the maximum concentration of a constituent in surface water exceeded the WQS, then that constituent was retained for quantitative analysis.

At the request of MDEQ (2000), if any single carcinogenic polycyclic aromatic hydrocarbon (cPAH) was retained as a COPC in a medium, then all cPAHs should also be retained as COPCs in that medium. This guidance refers to the following chemicals: benzo(a)anthracene, benzo(b)fluoroanthene, benzo(k)fluoranthene, benzo(a)pyrene, chrysene, dibenzo(a,h)anthracene, and indeno(1,2,3-cd)pyrene. To establish an exposure point concentration for undetected cPAHs retained as COPCs in an EU, one-half the maximum detection limit was used.

The results of the screening process are presented on the statistical summary tables, Tables 1 through 18. The screening process eliminated detected constituents from the subsurface soil dataset down to 20 feet bgs and surface soil dataset down to 6 feet bgs in EU3. For this reason, construction worker and maintenance worker exposures to soils in EU3 were not evaluated quantitatively in this assessment.

## 4.0 Exposure Assessment

Currently, a majority of the Site is used for commercial and light industrial purposes and is paved for roads and parking lots. Unpaved areas are limited to Gordon's Creek (EU 1), the wooded portion in and around the Fill Area (EU2) and the grassy field outlined by EU 3, and the drainage ditch delineated by EU 4 (Figure 2). Since the developed and undeveloped areas of the Site vary considerably with respect to both residual chemical concentrations and land use, the Site was divided into five EUs for the exposure assessment. A sixth EU was created to assess off-Site residential exposures. Chemical data from each EU were combined with EU-specific exposure parameter values and receptor scenarios to determine the chemical intake for each receptor potentially accessing an EU for occupational, recreational, or residential purposes.

### 4.1 Receptor Identification

The following exposures pathways (indicated with an "X") have been selected for this risk assessment as reasonable and realistic scenarios under current and future land-use assumptions:

| EU/Media:                | EU1  |             | EU2  | EU3  | EU4  |      |             | EU5  | EU6  |             |
|--------------------------|------|-------------|------|------|------|------|-------------|------|------|-------------|
| Receptor/Route:          | Sed. | Surf. Water | Soil | Soil | Soil | Sed. | Surf. Water | Soil | Sed. | Surf. Water |
| <b>Visitor</b>           |      |             |      |      |      |      |             |      |      |             |
| Dermal                   | X    | X           | X    | X    | X    | X    | X           | X    |      |             |
| Oral                     | X    | X           | X    | X    | X    | X    | X           | X    |      |             |
| Inhalation               |      |             |      |      |      |      |             |      |      |             |
| <b>Maint. Worker</b>     |      |             |      |      |      |      |             |      |      |             |
| Dermal                   | X    | X           | X    | X    | X    | X    | X           | X    |      |             |
| Oral                     | X    | X           | X    | X    | X    | X    | X           | X    |      |             |
| Inhalation               |      |             |      |      |      |      |             |      |      |             |
| <b>Const. Worker</b>     |      |             |      |      |      |      |             |      |      |             |
| Dermal                   | X    | X           | X    | X    | X    | X    | X           | X    |      |             |
| Oral                     | X    | X           | X    | X    | X    | X    | X           | X    |      |             |
| Inhalation               |      |             | X    | X    | X    |      |             | X    |      |             |
| <b>Off-Site Resident</b> |      |             |      |      |      |      |             |      | X    | X           |
| Dermal                   |      |             |      |      |      |      |             |      | X    | X           |

| EU/Media:       | EU1  |             | EU2  | EU3  | EU4  |             |      | EU5  | EU6         |   |
|-----------------|------|-------------|------|------|------|-------------|------|------|-------------|---|
| Receptor/Route: | Sed. | Surf. Water | Soil | Soil | Sed. | Surf. Water | Soil | Sed. | Surf. Water |   |
| Oral            |      |             |      |      |      |             |      | X    |             | X |
| Inhalation      |      |             |      |      |      |             |      |      |             |   |

In accordance with MCEQ guidance (1999), intake of cPAH compounds via the ingestion route were evaluated qualitatively (see Section 6.1) because the published cancer slope factor for benzo(a)pyrene (IRIS, 2000) cannot be used to quantify carcinogenic risks from ingestion. Specifically, the MCEQ guidance (1999) states:

Polycyclic aromatic hydrocarbons (PAHs) – If the dermal exposure pathway is a complete pathway, the slope factor for benzo(a)pyrene provided in IRIS shall be used to estimate the carcinogenic risk. The relative potencies provided by EPA's provisional guidance (ECAO-CIN-842, 3/93) shall be followed to assess dermal carcinogenic risks from other PAHs. The same slope factor cannot be used to quantify carcinogenic risks from incident ingestion of PAHs. *Such potential ingestion risk shall be described qualitatively.* [Emphasis added]

As a result of the screening process presented in the Data Evaluation section, *only cPAHs were COPCs for following scenarios:*

| Receptor            | Exposure Unit | Media         |
|---------------------|---------------|---------------|
| Adolescent Visitor  | 2             | Surface soils |
| Adolescent Visitor  | 5             | Surface soils |
| Adolescent Visitor  | 1             | Sediments     |
| Off-Site Resident   | 6             | Surface water |
| Maintenance Worker  | 2             | Surface soils |
| Maintenance Worker  | 5             | Surface soil  |
| Construction Worker | 2             | Soils         |
| Construction Worker | 5             | Soils         |
| Construction Worker | 1             | Sediment      |

Ingestion exposures for the above tabulated scenarios, therefore, were not evaluated quantitatively because of the lack of appropriate published toxicity values for benzo(a)pyrene in accordance with MCEQ guidance (1999).

Surface water present on-Site is either ephemeral or very shallow and is conducive only to wading-type activities. Ingestion of Site surface water was considered an insignificant exposure pathway since on-Site drainage ditches "contain little or no water most of the time" (MDEQ, 2000). In addition, US EPA IV guidance indicates that "In most cases, it is unnecessary to evaluate human exposures to sediments covered by surface water" (1995). At the request of MDEQ (2000), however, dermal and oral exposures to surface water were assessed for visitors, maintenance workers, and construction workers in EUs 1 and 4. Surface water exposures were also assessed for residents in off-Site EU 6.

Each of the potential receptors is discussed below.

#### 4.1.1 Infrequent Site Visitor

Since the Site is not currently fenced or guarded, the general public has access to most areas of the Site at any given time. It is possible that an individual may use some areas of the Site, such as EU1, EU2, or EU3, for recreational purposes. For this reason, sediment and surface water exposures to visitors in EU1, and surface soil exposures in EU2 and EU3 were assessed for the visitor scenario. A strong majority of the remainder of the Site (EU5) is covered with either buildings or pavement, precluding direct contact with surface soils; however, a small exposed area encompassing a drainage ditch exists along side of the Process Area (EU4). Although this area is not attractive for recreational purposes, it is possible that an individual traversing the Site may contact surface soils, sediment, or surface water in this EU; therefore, these potential exposures were assessed. Sediment exposures in EU1 and EU4 were addressed in accordance with US EPA Region 4 guidance that recommends evaluating sediment exposures in intermittent streams. At the request of MDEQ (2000), soil exposures were assessed for visitors in EU5 regardless of the existence of buildings and pavements precluding almost all potential direct contact with soils in this area.

#### 4.1.2 Maintenance Worker

Currently, maintenance activities are most likely limited to the developed portions of the Site. Of these, the Process Area and adjacent former drip track area (EU5) were most affected by historical wood preserving processes. Although these areas are mostly paved or built upon, it is possible that maintenance activities may require some shallow digging in unpaved areas; therefore, exposures to surface soils in EU5 were assessed. As a conservative measure, surface soil data from sample locations located in paved areas were evaluated in conjunction with surface soil data from exposed areas in EU5. If the currently undeveloped portions of the Site (EU2 and EU3) become developed in the future, similar maintenance activities may be required and, therefore, exposures to surface soils in EU2 and EU3 were also assessed. The drainage ditch encompassed by EU4 requires periodic maintenance; therefore, exposures to soil, sediment, and surface water in this area were assessed. At the request of MDEQ (2000), maintenance worker exposures to surface water and sediment in EU 1 were also assessed.

#### 4.1.3 Construction Worker

Although there are currently no major construction activities at the Site, such activities may hypothetically occur in the future. Thus, exposures to surface water and sediment in EUs 1 and 4, and exposures to soil in EUs 2 through 5 were assessed herein. Construction workers may be exposed to both surface and subsurface soils during activities such as excavating. Subsurface soils, for purposes of this assessment, were defined as those soils at the water table and shallower. Since the depth to the water varies significantly across the Site, so does the definition of "subsurface" soils. Accordingly, subsurface soils were evaluated down to 10 feet for EU2 and 20 feet for EUs 3, 4, and 5.

#### 4.1.4 Future On-Site Residents

The affected areas of the Property (the Site) are currently zoned for industrial or light-commercial use, and, at the time of this report, there were no plans to develop the Site for residential housing. In fact, deed restrictions preventing residential development are in the

process of being implemented for the impacted areas on Site. Because of these deed restrictions, it is reasonable and realistic to assume that the Site will remain commercial/industrial in the future; therefore, on-Site residential exposures were not addressed in this risk assessment.

#### 4.1.5 Off-Site Residential Exposures

The northeast drainage ditch extends from the former Process Area to the northeast into a nearby residential community. Surface water and sediment data from areas along the northeast drainage ditch (EU6, Figure 3) were evaluated for off-Site residential exposures. For purposes of exposure assessment, a child resident between the ages of 1 and 6 years and an adolescent/adult resident between the ages of 7 and 30 years were evaluated. Hazards and risks for these two receptors were then combined (summed) to reflect the exposures incurred by a single individual living off-Site in the vicinity of the northeast drainage ditch for 30 years.

#### 4.2 General Intake Equation

Chemical exposure/intake is expressed as the amount of the agent at the exchange boundaries of an organism (*i.e.*, skin, lungs, gut) that is available for systemic absorption. An applied dose is defined as the amount of a chemical at the absorption barriers such as skin, lung, digestive tract, available for absorption and is (usually expressed in milligrams, or mg) absorbed per unit of body weight of the receptor (usually expressed in units of kilogram, or kg). Absorbed dose can be defined as the amount of chemical that penetrates the exchange boundaries. If the exposure occurs over time, the total exposure can be divided by the time period of interest to obtain an average exposure rate (*e.g.*, mg/kg-day). The general equation, as defined by US EPA, for estimating a time-weighted average intake is:

$$\text{Intake (mg/kg - day)} = \frac{C \times IR \times EF \times ED}{BW \times AT} \quad [\text{Equation 1}]$$

where:

- C = chemical concentration at the exposure point (*e.g.*, mg/m<sup>3</sup> air);  
IR = intake rate (*e.g.*, m<sup>3</sup>/hr);  
EF = exposure frequency (days/year);

ED = exposure duration (years);  
BW = body weight of exposed individual (kg); and  
AT = averaging time (period over which exposure is averaged, usually measured in days).

Additional parameters (*e.g.*, skin surface area) were incorporated into the above general equation to evaluate the different potential exposure routes (dermal, oral, inhalation).

Table 19 presents the general and pathway-specific exposure parameters utilized for the intake equations in this assessment.

#### 4.2.1 General Exposure Parameters

Although some of the parameters used to calculate potential exposure are pathway- or route-specific, exposure frequency (EF), exposure duration (ED), averaging time (AT; determined separately for carcinogenic and non-carcinogenic exposures), and body weight (BW) are present in each intake model. These general parameters remain consistent throughout the intake calculations for each specific receptor.

##### 4.2.1.1 Exposure Frequency

The exposure frequency (EF) describes the number of times per year an event is likely to occur. It is most often expressed in units of days/year or events/year, depending on the scenario. Variables such as weather, vacations, sick days, and institutional controls often aid in determining reasonable and realistic exposure frequencies.

The EF for an adolescent visitor was extracted from US EPA *Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A) Interim Final* (1989). This EF value of 12 days/year per EU is a reasonable estimate that assumes an adolescent would most likely be engaged in outdoor activity on the unpaved areas of the Site for one day a week during the three warmest months of the year. This value was used for soil, sediment, and surface water exposures.

The EF parameter used for construction workers was 80 days. This is a very conservative assumption in that typical construction projects, especially at industrial complexes, generally involve several phases of activity prior to completion. Examples of these activities include foundation excavation, foundation pouring, building framing, plumbing installation, electrical installation, and roofing. Generally, to complete each of these phases, a different team of specialized contractors is employed to perform the tasks for which they are most qualified. As a result, an individual may only remain at the construction site for a few weeks until his/her task has been completed and the next phase has begun. This is especially true for those activities involving direct contact with soil such as excavating and foundation pouring. Individuals performing these tasks are not usually qualified or employed to continue with the actual building processes. An 80-day EF accounts for an individual to be involved in construction activities for four entire months of the year (assuming five-day work weeks). As explained above, this is a very conservative assumption.

For surface water and sediment exposures to construction workers, an EF value of 8 days/year was used. This value represents 1/10<sup>th</sup> of the time a worker may be on-Site for construction-type activities and is conservative in that it is unlikely that construction workers would be exposed at all to Site surface water or sediment.

The EF value used for the maintenance worker scenario was 150 days/year for surface soil exposures in EUs 2, 3, and 5. This is also a conservative assumption in that the currently developed areas of the Site are covered with buildings or pavement. Maintenance activities in these areas would require little contact with the obscured surface soils. The undeveloped areas of the Site currently require little or no maintenance as they are only occasionally mowed or allowed to grow naturally. Should these areas become developed, they will most likely take on the appearance of the remainder of the Site, including industrial/commercial buildings and paved roads or parking lots. Once again, extensive direct contact with surface soils would be minimal for a maintenance worker.

For maintenance worker sediment and surface water exposures in EU 1 and 4 and surface soil exposures in EU 4, an EF value of 30 days/year was used. Historically, the northeast drainage ditch has been maintained on an as-needed basis (less than annually). Maintenance worker exposures to sediment and surface water in these areas were assessed at the request of the MDEQ (2000). An EF value of 30 days/year is amply conservative in that both Gordon's Creek (EU 1) and the northeast drainage ditch (EU 4) are currently maintained less than annually.

For residential soil exposures, an exposure frequency of 350 days/year was used in accordance with Region IV guidance. This value assumes that 15 days/year are spent away from home (US EPA, 1991).

Sediments along the bank of the northeast drainage ditch are not comparable to surface soils comprising a yard with respect to exposure. Typically, yard soils include relatively large areas where children frequently play and where surface soils are tracked into the home to become part of the household dust that can be ingested, particularly by crawling infants, on a daily basis. These are the assumptions that underlie the standard residential soil exposure algorithm and parameter values. However, it is not realistic to assume that infants, children, or adults will directly contact a relatively small area of sediments on the banks of a drainage ditch on a daily basis. A more realistic exposure scenario for this unique area under an assumption of residential land use is for a resident child to play on occasion in the drainage ditch that traverses the residential property. An exposure frequency of 40 days/year, two hours per exploring event, is conservatively plausible.

#### 4.2.1.2      Exposure Duration

The ED parameter represents the number of years during which an event is likely to occur. Factors affecting this parameter include variables such as age of receptor, population mobility, and occupational mobility. Exposure durations of less than seven years typically correspond to subchronic exposures while those greater than seven years are typically considered chronic exposures (US EPA, 1989). Toxicity indices are selected based on subchronic or chronic exposure durations.

The future construction worker scenario used an ED of one year because it is highly unlikely that a future construction worker would remain on one site for more than a year. Often, two months is considered the maximum amount of time a construction worker may reasonably remain at the same site.

The future maintenance worker ED, on the other hand, is based on occupational mobility studies. The ED of 25 years was obtained from US EPA (1991) which recommends a 95th percentile value of 25 years based on a study by the Bureau of Labor Statistics as of 1987. US EPA Region 4 also recommends a default value of 25 years for worker scenarios (1995).

The adolescent visitor scenario used an ED of 10 years. An adolescent was defined in this assessment as an individual aged seven to 16 years in accordance with US EPA Region 4 (1995); therefore, an exposure duration of 10 years was most appropriate.

An ED of 30 years (US EPA Region 4, 1995) was used for off-Site residents. This value assumes an individual spends 6 years as a child and 24 years as an adolescent/adult in the same location.

#### 4.2.1.3 Averaging Time

The averaging time (AT) parameter is the time period over which exposure is averaged. For human health cancer risk calculations, the AT<sub>c</sub> value prorates a total cumulative dose over a lifetime. As a conservative approach, the AT<sub>c</sub> value for each receptor is the product of a 365-day year and a 70-year life span, equaling 25,550 days.

The AT<sub>n</sub> used for non-carcinogenic effects is the product of a 365-day year and the exposure duration (*i.e.*, AT<sub>n</sub> = 365 days × ED). Because the ED parameter changes for each receptor, the AT<sub>n</sub> changes as well. The AT<sub>n</sub> values used for each receptor are summarized below:

Future Construction Worker - 365 days

Maintenance Worker - 9125 days

Adolescent Visitor - 3650 days  
Off-Site Child Resident – 2,190 days  
Off-Site Adult Resident – 8,760 days

#### 4.2.1.4 Body Weight

The body weight used for the adult exposures (future construction worker and maintenance worker) analyzed in this assessment was the current US EPA default value of 70 kg (US EPA, 1989; US EPA Region 4, 1995). This value was also used for the adolescent/adult off-Site resident scenario. The adolescent body weight used for the visitor scenarios was 45 kg. This value was extracted from US EPA Region 4 guidance (1995). For the child resident scenario, a body weight of 15 kg was used as recommended by US EPA (1991).

#### 4.2.2 Route-Specific Exposure Parameters

The general intake equation discussed above (Equation 1) was modified by including route-specific exposure parameters in order to calculate route-specific intake values. For dermal exposures, skin surface area, adherence factor, exposure time (surface water exposures only), and absorption factor parameters were included in the intake equation. For ingestion exposures, an ingestion rate and a matrix effect were included in the intake calculation. For inhalation exposures, an inhalation rate and a retention factor for fugitive dusts were included in the intake equation. Also, for inhalation exposures, an additional paradigm was necessary to convert soil concentrations to concentrations in air available for intake.

##### 4.2.2.1 Dermal Exposure Parameters

###### Skin Surface Area

The total skin surface area used for adult receptors in this assessment was 20,000 cm<sup>2</sup>. This is a US EPA default value extracted from the *Exposure Factors Handbook* (1997). For adolescent exposures, a value of 12,768.3 cm<sup>2</sup> was used for total skin surface area. This was a mean value calculated based on the distributions of total skin surface areas for males and females between the ages of 7 and 16 as presented in *Exposure Factors Handbook* (1997). For the off-Site child resident scenario, a skin surface area of 7,213 cm<sup>2</sup> was used. This value was based on skin surface area data for male and female children provided in *Exposure Factors Handbook* (1997).

For purposes of exposure, it was assumed that only portions of the body would be exposed to the affected media on the Site. For the construction worker scenario, it was assumed that the hands, forearms, lower legs, and face would be exposed to Site soils. These body parts comprise 27.8% of the total skin surface area, or 5560 cm<sup>2</sup>.

For maintenance worker exposures to Site soils, it was assumed that the hands, forearms, and face would be exposed. These body parts comprise 15 percent of the total skin surface area, or 3000 cm<sup>2</sup>.

For surface water and sediment exposures, exposed body parts for construction and maintenance workers included hands, forearms, and feet or 3620 cm<sup>2</sup> (18.1% of the total skin surface area).

The visitor and off-Site resident scenarios assumed that the hands, forearms, and lower legs would be exposed for contact with Site soils. These body parts comprise 23.9% of the total skin surface area, or 3052 cm<sup>2</sup> for adolescent visitors, 1724 cm<sup>2</sup> for child residents, and 4780 cm<sup>2</sup> for adult residents. For exposures to surface water and sediment, hands, forearms, feet were assumed exposed for adolescent visitor and off-Site resident scenarios. These body parts comprise 18.1 % of the total skin surface area or 2311 cm<sup>2</sup> for adolescent visitors, 1306 cm<sup>2</sup> for child residents, and 3620 cm<sup>2</sup> for adult residents.

#### Soil Adherence Factor

Until recently, the US EPA-recommended default for soil adherence on skin ranged from 0.2 to 1.0 mg/cm<sup>2</sup> for the entire exposed surface area, without consideration of the type of activity (US EPA, 1992). However, the data from which that range was derived were primarily the result of indirect measurements, artificial activities, and sampling of hands only. A more recent study has presented the results of direct measurement of soil loading on skin surfaces before and after normal occupational and recreational activities that might result in soil contact (Kissel *et al.*, 1996). A five-order of magnitude range (roughly 10<sup>-3</sup> to 10<sup>+2</sup> mg/cm<sup>2</sup>) was reported for observed activity-related

hand loadings. That report indicated that hand loadings within the range of 0.2 to 1.0 mg/cm<sup>2</sup> were produced by activities in which there was vigorous soil contact (e.g., rugby, farming); but for activities in which there was less soil contact (e.g., soccer, professional grounds maintenance), loadings substantially less than 0.2 mg/cm<sup>2</sup> were found on hands and other body parts. Kissel *et al.* (1996) concluded that, because non-hand loadings attributable to higher contact activities exceeded hand loadings resulting from lower contact activities, hand data from limited activities cannot be used as a conservative predictor of loadings that might occur on other body surfaces without regard to activity. Furthermore, because exposures are activity-dependent, dermal exposure to soil should be quantified using data describing human behavior (e.g., type of activity, frequency, duration, including interval before bathing, clothing worn, etc.).

The most recent version of the *Exposure Factors Handbook* (1997) states:

In consideration, of these general observations and the recent data from Kissel *et al.* (1996, 1997), this document recommends a new approach for estimating soil adherence to skin. First use Table 6-12 [Summary of Field Studies, Kissel *et al.*, 1996a] to select the activity which best approximates the exposure scenario of concern. Next, use Table 6-13 [Mean Soil Adherence by Activity and Body Region, Kissel *et al.*, 1996a] to select soil loadings on exposed skin surfaces which correspond to the activity of interest. This table contains soil loading estimates for various body parts. The estimates were derived from soil adherence measurements of body parts of individuals engaged in specific activities described in Table 6-12. These results provide the best estimate of central loadings, but are based on limited data. Therefore, they have a high degree of uncertainty such that considerable judgment must be used when selecting them for an assessment.

In another study that assessed the percentage of skin coverage in several soil contact trials in a greenhouse and an irrigation pipe laying trial, Kissel *et al.* (1996) concluded that adjusted loadings may be two to three orders of magnitude larger than average loadings if average loadings are small.

The activity-specific soil adherence factor for exposures to a maintenance worker was calculated based on data presented by Kissel *et al.* (1996) for grounds keepers, as presented below:

|                    |                         | Soil Adherence Factor by Body Part (mg/cm <sup>2</sup> ) |                |                 |               |               |
|--------------------|-------------------------|--|----------------|-----------------|---------------|---------------|
| Receptor           | Representative Activity | Hands  | Arms           | Lower Legs      | Face          | Feet          |
| Maintenance Worker | Grounds Keepers         | 0.030 - 0.15   | 0.0021 - 0.023 | 0.0008 - 0.0012 | 0.0021 - 0.01 | 0.004 - 0.018 |

Data for the grounds keepers were used for the maintenance worker estimates because the activities of a grounds keeper best mimic those of a maintenance worker.

Soil adherence factors were calculated by normalizing each body part-specific soil adherence value (using the mid-points of the ranges tabulated above) with regard to the percentage of total body surface area represented by the respective body part (extracted from the US EPA *Dermal Exposure Assessment: Principles and Applications* [US EPA, 1992]). The maintenance worker adherence factor for soil was calculated based upon exposure to the hands, forearms and face. Surface area percentages for the hands, forearms, and face are 5.2, 5.9, and 3.9 percent, respectively (US EPA, 1997). Those body parts comprise 15 percent of the total body surface area. The normalized values for all body parts of interest were added, and the sum was divided by the total percentage of body surface area occupied by the parts. For example, the soil adherence factor for the maintenance worker soil exposures (0.038 mg/cm<sup>2</sup>) was calculated as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.09 \times 0.052) + (0.0126 \times 0.059) + (0.006 \times 0.039)}{0.15} = 0.038$$

For sediment exposures, the soil adherence factor for maintenance workers was calculated using grounds keeper adherence data (tabulated above) for the hands, forearms, and feet. These body parts comprise 5.2, 5.9, and 7.0 percent of the total skin surface area, respectively, for a total of 18.1% of the total skin surface area. The adherence factor (0.034 mg/cm<sup>2</sup>) was calculated as follows:

$$AF \text{ (mg/cm}^2\text{)} = \frac{(0.09 \times 0.052) + (0.0126 \times 0.059) + (0.011 \times 0.07)}{0.181} = 0.034$$

The construction worker adherence factor was also calculated in this fashion. This exposure scenario assumed that the hands, forearms, lower legs, and face would be exposed to Site soils. Soil loadings for the upper torso (chest and back) were not measured by Kissel *et al.* (1996) for construction workers because this body area is generally covered. However, to account for exposure to the upper torso during the very hot months of the year, the total area of the forearms, legs, hands, and face were assumed to be completely exposed. The hands, forearms, legs, and face comprise 5.2%, 5.9%, 12.8%, and 3.9% of the total skin surface area, respectively (with the face comprising one-third the surface area of the head), for a total of 27.8% exposed surface area. The construction worker soil adherence factor was based on data from Kissel *et al.* (1996) for construction workers as follows:

|                     |                                | <b>Soil Adherence Factor by Body Part (mg/cm<sup>2</sup>)</b> |             |                   |             |
|---------------------|--------------------------------|---|-------------|-------------------|-------------|
| <b>Receptor</b>     | <b>Representative Activity</b> | <b>Hands</b>  | <b>Arms</b> | <b>Lower Legs</b> | <b>Face</b> |
| Construction Worker | Construction Worker            | 0.24  | 0.098       | 0.066             | 0.029       |

The soil adherence factor for the construction worker scenario was calculated as follows:

$$AF \text{ (mg/cm}^2\text{)} = \frac{(0.24 \times 0.052) + (0.098 \times 0.059) + (0.066 \times 0.128) + (0.029 \times 0.039)}{0.278} = 0.1$$

For sediment exposures, the soil adherence factor was calculated for the construction worker scenario using hands, forearms, and feet adherence data. Since adherence data for feet were not provided in US EPA *Exposure Factors Handbook* (1997) for the construction worker activity, the data tabulated above for grounds keepers was used for the feet adherence factor. The hands,

forearms, and feet comprise 5.2, 5.9, and 7.0 percent of the total skin surface area, respectively (totaling 18.1 percent). Thus, the adherence factor for construction workers exposed to sediment ( $0.105 \text{ mg/cm}^2$ ) was calculated as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.24 \times 0.052) + (0.098 \times 0.059) + (0.011 \times 0.07)}{0.181} = 0.105$$

The adherence factor for visitor and off-Site resident exposures to soil assumed that the forearms, hands, and lower legs would be exposed to soil or sediment. The data used in these calculation were based on data by Kissel *et al.* (1996) for soccer players (exposed to a playing field of roughly one-half grass and one-half bare earth in a light mist) as presented below:

| <b>Receptor</b>               | <b>Representative Activity</b> | <b>Soil Adherence Factor by Body Part (<math>\text{mg/cm}^2</math>)</b> |              |                   |
|-------------------------------|--------------------------------|---|--------------|-------------------|
|                               |                                | <b>Arms</b>   | <b>Hands</b> | <b>Lower Legs</b> |
| Visitor and Off-Site Resident | Soccer Players                 | 0.0029 – 0.011  | 0.019 – 0.11 | 0.0081 – 0.031    |

The forearms, hands, and lower legs comprise 5.9%, 5.2%, and 12.8% of the total skin surface area, respectively, for a total of 23.9% (US EPA *Exposure Factors Handbook*, 1997). The adherence factor was then calculated for visitor and off-Site resident dermal exposures to soil as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.00695 \times 0.059) + (0.0645 \times 0.052) + (0.0196 \times 0.128)}{0.239} = 0.026$$

A value of  $0.026 \text{ mg/cm}^2$  was used as the soil adherence factor for visitors to the Site and off-Site residents.

Soil adherence factors for sediment exposures to Site visitors and off-Site residents were calculated using adherence data for the hands, forearms, and feet. Adherence data for reed gatherers were used for these exposures to best mimic activities that may incur sediment exposures. The reed gatherers studied by Kissel *et al.* (1996) periodically visited tidal flats to collect raw materials for basket weaving. The data from Kissel *et al.* (1996) presented in *Exposure Factors Handbook* (US EPA, 1997) were as follows:

|                                 |                         | Soil Adherence Factor by Body Part (mg/cm <sup>2</sup> ) |       |      |
|---------------------------------|-------------------------|--|-------|------|
| Receptor                        | Representative Activity | Hands  | Arms  | Feet |
| Visitors and Off-Site Residents | Reed Gatherers          | 0.66   | 0.036 | 0.63 |

The hands, forearms, and feet comprises 5.2, 5.9, and 7.0 percent of the total skin surface area, respectively (totaling 18.1 percent). Thus, the adherence factor for visitors and off-Site residents exposed to sediment (0.044 mg /cm<sup>2</sup>) was calculated as follows:

$$AF \text{ (mg/cm}^2\text{)} = \frac{(0.66 \times 0.052) + (0.036 \times 0.059) + (0.63 \times 0.07)}{0.181} = 0.044$$

#### Exposure Time

To estimate intakes as a result of dermal exposure to surface water, an exposure time (ET) parameter was included in the intake formula for Site visitors and off-Site residents. The parameter value of 1.0 hour/day was estimated using best professional judgement. This value represents the amount of time a Site visitor or off-Site resident may spend exposed to surface water in any one EU.

#### Dermal Permeability Constant

The permeability constant, Kp, accounts for the movement of a constituent dissolved in water through the skin, across the stratum corneum, and into the blood stream. Kp values for the

constituents examined in this assessment for surface water exposures were obtained from US EPA *Dermal Exposure Assessment: Principles and Applications* (1992). For values not available in US EPA *Dermal Exposure Assessment* (1992), the K<sub>p</sub> value can be calculated using the equations provided by the US EPA in the same document.

#### Dermal Absorption Factor

The final parameter included in the dermal intake paradigm was a dermal absorption factor. In general, the skin provides an effective barrier to environmental toxins. For example, certain hair-coloring formulations which are vigorously rubbed onto the scalp on a daily basis contain lead acetate at concentrations up to 200,000 ppm, yet lead toxicity does not appear to result. Moore *et al.* (1980) determined that the rate of lead absorption from <sup>203</sup>Pb labeled lead acetate in cosmetic preparations containing six mmol Pb acetate/L in male volunteers over 12 hours was 0.06% during normal use of such preparations. For most inorganic salts, percutaneous (skin) absorption is considered insignificant relative to incidental ingestion (for example, US EPA, 1986). On the other hand, some drugs (*e.g.*, nicotine) are effectively administered and absorbed into the blood stream from dermal "patches."

Most dermal bioavailability data for impacted soil have been obtained in laboratory animals or in vitro test systems. This introduces a significant source of uncertainty for predicting the human response. Safety factors have sometimes been applied to dermal absorption data obtained in animals to conservatively estimate the upper-bound of likely human percutaneous uptake of a certain constituent from skin exposure. This is usually unnecessary because human skin has generally been shown, for a diverse group of constituents, to be about 10-fold less permeable than the skin of typical animal species, such as rabbits and rats (Bartek and LaBudde, 1975; Shu *et al.*, 1988).

US EPA Region III evaluated available data concerning the dermal absorption of specific constituents and classes of constituents and provided several recommendations (US EPA Region 3, 1995). For semivolatile compounds, such as *bis*(2-ethylhexyl)phthalate, the US EPA

recommends a range of 1% to 10% (US EPA, 1995). Kao *et al.* (1985) reported 2.7 percent for absorption of topically applied pure benzo(a)pyrene by human skin *in vitro*. The US EPA Region 3 recommends using 10% as a conservative assumption based on the Ryan *et al.* study (1987). In addition, US EPA Region 4 guidance (1995) states that a soil dermal absorption factor “of 1.0% for organics and 0.1% for inorganics should be used as defaults in determining the uptake associated with dermal exposure” (see the Dermal Contact subsection of Exposure Assessment section of the 1995 guidance). For the purpose of this risk assessment, an ABS of 3% for benzo(a)pyrene and of 10% for other semivolatile organic compounds were conservatively assumed for dermal absorption, in keeping with US EPA Region 3’s recommendations.

#### 4.2.2.2 Ingestion Exposure Parameters

##### Ingestion Rate

US EPA’s *Exposure Factors Handbook* (1997) discusses three adult soil ingestion studies with results ranging from 10 mg/day to 480 mg/day. Hawley’s (1985) value of 480 mg/day (as recommended by the MDEQ) was “derived from assumptions about soil/dust levels on hands and mouthing behavior” (US EPA, 1997). Since no supporting measurements were made for Hawley’s study, the US EPA states that Hawley’s estimate “must be considered conjectural” (1997). As such, the US EPA goes on to suggest adult soil ingestion rates of 50 mg/day for industrial settings and 100 mg/day for residential and agricultural settings, although “50 mg/day still represents a reasonable central estimate of adult soil ingestion and is the recommended value...” (1997). Accordingly, a value of 100 mg/day for the maintenance worker, construction worker, and adult off-Site resident is amply conservative and was used in this assessment.

The ingestion rate used for the adolescent visitor and off-Site child resident scenarios was 100 mg/day. The US EPA *Exposure Factors Handbook* (1997) recommends a value of 100 mg/day as a mean ingestion rate for children under six years of age. This value was conservatively used in this assessment to estimate soil and sediment ingestion exposures for an adolescent visitor aged 7 to 16 years and an off-Site resident aged one to six years.

### Gastrointestinal Matrix Effects of Soil

Incidental ingestion incorporates the matrix effect (ME; sometimes called the absorption adjustment factor [AAF]) into the general intake equation. When constituents are administered in solid vehicles such as food and soil, only a fraction of the ingested dose is extracted from the vehicle and subsequently absorbed through the gastrointestinal tract (US EPA *Estimated Exposure to Dioxin-like Compounds*, 1992). Gastrointestinal absorption of constituents sorbed onto such a medium is inhibited by physical-constituent bonding to the matrix (Hawley, 1985). This phenomenon is referred to as the gastrointestinal matrix effect of soil. Several studies referenced in the US EPA's *Estimated Exposure to Dioxin-like Compounds* (1992) have been performed to estimate the oral absorption factors of constituents from soil.

For PAHs, a gastrointestinal matrix effect value of 0.29 was utilized as recommended by Magee *et al.*, 1996. This value is the average of 12 oral soil AAF studies performed by Rozett *et al.*, Weyand *et al.*, and Goonet (Magee, 1996).

#### 4.2.2.3 Inhalation Exposure Parameters and Paradigms

##### Inhalation Rate

The inhalation rate used for the construction worker scenario was 20 m<sup>3</sup>/day. This is a common US EPA default value and was recommended by US EPA Region 4 (1995).

##### Retention Factor

According to the International Commission on Radiological Protection (ICRP), 75 percent of respirable dust particles (PM<sub>10</sub>, or particles less than 10 microns in aerodynamic diameter) are retained when inhaled, the vast majority of which is potentially subsequently swallowed (ICRP, 1968). This 75% was included in the inhalation intake equation as the retention factor parameter (RF). This parameter applies only to non-VOC constituents entrained onto dust particles.

### Concentration in Air

To estimate airborne dust levels during hypothetical construction activities, an emission rate of suspendible particles of less than 15 microns in aerodynamic diameter ( $PM_{15}$ ) was calculated (grams/second); particles less than 10 microns were considered to be respirable. Considering particles of 15 microns or less in diameter in the emission rate calculation is a conservative assumption, inasmuch as only particles with an aerodynamic diameter of less than five to seven microns are inhaled into the lung.

The two types of construction activities at the Site that have the potential to emit fugitive dusts are vehicular movement over bare (unpaved or unvegetated) surfaces and the excavation of soil. Estimation of fugitive dust emissions caused by each activity were examined separately, as follows, and were derived from existing estimates of general construction exposure. The sum of the emissions from these two activities was multiplied by the concentration of constituent in the soil ( $C_s$ ) in order to derive the total emission rate ( $E_i$ ) for non-VOCs as follows:

$$E_i = C_s \times (PERv + PERe) \quad [Equation 2]$$

where:

- $E_i$  = Emission rate (mg/sec);  
 $C_s$  = Concentration in soil (mg/kg);  
 $PERv$  = Particulate emission rate for vehicular movement (lb/vehicle mile);  
and  
 $PERe$  = Particulate emission rate for excavation (lb/vehicle mile).

The following empirical expression (US EPA, 1988) was used to estimate the fugitive dust generated by vehicles during construction activities:

$$PERv (\text{lbs/vehicle mile}) = k \times 5.9 \times (s/12)(S/30) \times (mvw/3)^{0.7} \times (ww/4)^{0.5} \times ((365 - p)/365) \quad [Equation 3]$$

where:

- $PERv$  = Vehicle particle emission rate (lb/vehicle mile traveled);

s = Percent silt content (unitless);  
 k = Particle size multiplier (unitless);  
 S = Mean vehicle speed (mph);  
 mvw = Mean vehicle weight (ton);  
 ww = Mean number of wheels per vehicle (unitless); and  
 p = Mean number of days with  $\geq$  0.01 inches of precipitation per year (unitless).

It was assumed that the vehicle travels during 40% of the 80-day exposure duration and 0.5 miles per day. The result is a value of 16 miles per construction event. Percent silt content was estimated to have a mean value of 50%, based on geotechnical data provided in the *Remedial Investigation Report* (Pisani & Assoc., 1997). US EPA default values were utilized and referenced for all other parameters. The particle size multiplier was assumed to be 0.50, corresponding to particles less than 15 microns (US EPA, 1996). Vehicle characteristics consist of the following: mean vehicle speed was assumed to be 15 mph, with mean vehicle weight assumed to be approximately 12.5 tons, for 8-wheeled vehicles (US EPA, 1988). The estimated mean number of days with precipitation equal to or greater than 0.01 inches per year is 110 (US EPA, 1988). Total resultant dust emissions for constituents during vehicular movement activities were estimated to be approximately 16.5 lbs/vehicle mile traveled, or 0.0001 kg/sec. Calculations are summarized in Table 20.

Future excavation may be performed by bulldozers, a backhoe, or other heavy construction equipment. The following estimate of particulate emissions, less than  $15 \mu\text{m}$  in diameter resulting from bulldozing activity, was based on the approach described in the US EPA *Compilation of Air Pollution Emission Factors* (1996), as developed from studies of emissions from uncontrolled open dust sources resulting from bulldozing at western surface coal mines.

$$\text{PERe (lb/hour)} = \frac{1.0 \times s^{1.5}}{M^{1.4}} \quad [\text{Equation 4}]$$

where:

PERe = Excavation particle emission rate (lb/hr);  
s = Percent silt content (unitless); and  
M = Soil moisture content (unitless).

Percent soil moisture content was assumed to be 15.1%, an average of Site-specific soil moisture data and percent silt content 50%, as described above.

The resultant fugitive dust emission rate during excavation activities was 7.9 lbs/hr or 0.001 kg/sec. Table 20 summarizes these calculations.

Once the emission rate (Ei in Equation 2) was calculated, it was converted to a concentration in ambient air. Gaussian models are conventionally used to determine downwind ambient air concentrations, Ca, from the emission rate, Ei, estimated. However, in this scenario, such models have limited applicability when the receptor(s) is at or very near the source of emission. In this case, a bulldozer operator, for example, is situated directly within the area of ground emissions of vapors and dusts. Average ambient air concentrations in this circumstance are best estimated by use of a near-field box model (US EPA, 1988).

The near-field box model assumes uniform wind speed and uniform mixing throughout the box. The release and mixing of VOCs or respirable dusts in ambient air is estimated as follows:

$$Ca \text{ (mg/m}^3\text{)} = \frac{Ei}{W_b \times H_b \times V} \quad [\text{Equation 5}]$$

where:

Ca = Concentration of constituent in ambient air (mg/m<sup>3</sup>);  
Ei = Emission rate of constituent (mg/sec);  
W<sub>b</sub> = Width of box in crosswind dimension within the area of residual constituent in soil (m);  
H<sub>b</sub> = Downwind height of box (m); and

V = Average wind speed through the box (m/sec).

The value of  $H_b$  in this calculation is determined by the downwind distance and the atmospheric turbulence at ground level, which determines the trajectory of a release from the upwind edge of the source of vapor or dust emissions. For neutral atmospheric conditions, the height at the downwind boundary ( $H_b$ ) may be expressed by the following function (Pasquill 1975, Horst 1979):

$$z = 6.25 r [H_b/r \times \ln(H_b/r) - 1.58 H_b/r + 1.58] \quad [\text{Equation 6}]$$

where:

$H_b$  = Downwind height of box (m);  
 $z$  = Downwind distance to boundary (m); and  
 $r$  = A terrain-dependent roughness height (m)

$H_b$  (defined in Equation 5) is adjusted until the  $z$  parameter is equal to  $W_b$  (defined in Equation 5). The resulting  $H_b$  value is the height of the box. On any given workday, it is estimated that grading or excavation activities occur over the entire "workable" Site area (exposure unit) from which dusts are generated. This area is estimated to be 2,500 m<sup>2</sup>, with length of the box estimated to be 50 meters (downwind distance) and the width of the box ( $W$ ) estimated to be 50 meters. The greater the roughness height, the greater the wind turbulence and constituent dilution (*i.e.*, the height of the box increases). For the purposes of this risk assessment, it is conservatively assumed that the roughness height is 0.20 meters, which corresponds to a terrain with grass, some small bushes, and occasional trees (US EPA *Rapid Assessment of Exposure to Particulate Emission from Surface Contamination Sites*, 1985). This assumption is appropriate for the actual Site conditions. An annual average wind speed (4.69 m/sec) is obtained from the STAR data set, accessed through the Personal Computer Graphical Exposure Modeling System (PCGEMS), for STAR station 03940, Jackson/Thompson, MS for the period 1974-1978 (Table 21).

## **5.0      Toxicity Assessment**

The toxicity assessment involves the evaluation of available toxicity information to be utilized in the risk assessment process. Toxicity values derived from a dose-response relationship can be used to estimate the potential for the occurrence of adverse effects in individuals exposed to various constituent levels.

Exposure to a constituent does not necessarily result in adverse effects. The relationship between dose and response defines the quantitative indices of toxicity required to evaluate the potential health risks associated with a given level of exposure. If the nature of the dose-response relationship is such that no effects can be demonstrated below a certain level of exposure, a threshold can be defined and an acceptable exposure level derived. Humans are routinely exposed to naturally-occurring constituents and man-made constituents through the typical diet, air, and water, with no apparent adverse effects. However, the potential for adverse effects may occur if the exposure level exceeds the threshold in a variably sensitive population. This threshold applies primarily to constituents which produce non-carcinogenic (systemic) effects, although there is a growing body of scientific evidence which suggests that exposure thresholds may exist for certain carcinogenic constituents as well.

Adverse effects can be caused by acute exposure, which is a single or short-term exposure to a toxic substance, or by chronic exposure on a continuous or repeated basis over an extended period of time. "Acceptable" acute or chronic levels of exposure are considered to be without any anticipated adverse effects. Such exposure levels are commonly expressed as reference doses (RfDs), health advisories, etc. An acceptable exposure level is calculated to provide an "adequate margin of safety."

Chronic RfDs, which have been derived by the US EPA for a large number of constituents, were utilized to evaluate exposures lasting seven to 70 years (US EPA, 1989). Activities involving exposures of shorter duration to COPCs at the Site are anticipated to result in hazard and risk estimates that are lower than those associated with the long-term exposures. Identification of

subchronic toxicity values corresponding to shorter-term exposure scenarios (*i.e.*, less than seven years) are included in the risk assessment to ensure that both short-term and long-term risks can be addressed.

Currently, the US EPA has not developed toxicity values to be utilized in dermal exposure scenarios; however, the US EPA does provide the following guidance for dermal exposure:

No RfDs or slope factors are available for the dermal route of exposure. In some cases, however, non-carcinogenic or carcinogenic risks associated with dermal exposure can be evaluated using an oral RfD or oral slope factor, respectively. (US EPA, 1989).

Provisional dermal toxicity values were developed and utilized in the dermal exposure pathways considered in the human health risk assessment to provide a more accurate Site-specific risk assessment. These dermal RfD values were developed by multiplying the published oral RfD for a given constituent by the fraction of that constituent that can be absorbed through the gastrointestinal tract (stomach/intestine lining). The absorption fraction utilized was 50% for semivolatiles as extracted from US EPA Region 4 guidance (1995).

A number of sources of toxicity information exists, and these sources vary with regard to the availability and strength of supporting evidence. The following protocol has been established for the determination of toxicity indices; it defines a hierarchy of sources to be consulted and the methodology for the determination of toxicity values. This protocol has been developed in accordance with current US EPA methodology. Toxicity values for the COPCs at the Site were obtained with reference to the following hierarchy of sources developed in accordance with MCEQ guidance (1999):

- 1) Toxicity values were obtained from the *Integrated Risk Information System* (IRIS, 1999) database. This database contains the RfDs and Cancer Slope Factors (CSFs), which have been verified by the US EPA's RfD and Carcinogen Risk Assessment Verification Endeavor (CRAVE) workgroups, and is, thus, the

agency's preferred source for toxicity values. IRIS supersedes all other information sources.

- 2) For toxicity values which are unavailable on IRIS, the most current source of information is the Health Effects Assessment Summary Tables (HEAST, US EPA, 1997), published by the US EPA. HEAST contains interim, as well as verified RfDs and CSFs. Supporting toxicity information for verified values is provided in an extensive reference section of HEAST.
- 3) In cases where IRIS or HEAST could not provide toxicity values, US EPA Region III's Risk-Based Concentration (RBC) Tables were visited. These tables often provide toxicity values generated by reliable sources other than IRIS or HEAST. For example, in response to specific requests from risk assessors, the US EPA National Center for Environmental Assessment (NCEA) develops provisional RfDs or CSFs for chemicals not listed in IRIS or HEAST. Region III's RBC tables will list such provisional values. Also, RfDs or CSFs that have since been withdrawn from IRIS or HEAST may still be listed on the Region III RBC tables, although they are flagged with a "W." These toxicity values were no longer agreed upon by US EPA scientists; however, the Region III RBC tables continue to publish such values because risk assessors still need to quantify exposures to these chemicals. Lastly, the Region III RBC tables will list toxicity indices found in "other" US EPA documents. These values are flagged with an "O" on the tables.

The US EPA has derived carcinogenic slope factors for both oral and inhalation pathways, and these are utilized to quantitatively estimate risks. In the first step of the US EPA's evaluation, the available data are analyzed to determine the likelihood that the agent is a human carcinogen. The evidence is characterized separately for human studies and animal studies as sufficient, limited, inadequate, no data, or evidence of no effect. The characterizations of these two types of data are combined, and based on the extent to which the agent has been shown to be a carcinogen in experimental animals or humans, or both, the agent is given a provisional weight-of-evidence classification. The US EPA scientists then adjust the provisional classification upward or downward, based on other supporting evidence of carcinogenicity (see Section 7.1.3, US EPA, 1989). For a further description of the role of supporting evidence, see the US EPA guidelines (US EPA, 1986).

The US EPA classification system for weight of evidence is shown in the table below. This system is adapted from the approach taken by the International Agency for Research on Cancer.

| US EPA WEIGHT-OF-EVIDENCE<br>CLASSIFICATION SYSTEM FOR<br>CARCINOGENICITY |  |
|---|--|
| Group   | Description  |
| A   | Human carcinogen   |
| B1 or<br>B2   | Probable human carcinogen<br><br>B1 indicates that limited human data are available<br><br>B2 indicates sufficient evidence in animals and inadequate or no evidence in humans |
| C   | Possible human carcinogen  |
| D   | Not classifiable as to human carcinogenicity   |
| E   | Evidence of non-carcinogenicity for humans   |

(US EPA, 1989)

Table 22 summarizes the available toxicity values for the identified COPCs. COPCs lacking published toxicity values were not able to be quantitatively evaluated in this assessment in accordance with MCEQ guidance (1999). The MCEQ limits the use of toxicity values to those that have been published in IRIS, HEAST, ATSDR toxicity profiles, or other peer-reviewed reference sources or literature approved by the MCEQ (1999).

## **6.0 Risk Characterization**

The objective of the risk characterization is to determine potential risk to receptors by combining the results of the exposure and toxicity assessments. Non-carcinogenic effects and carcinogenic risks are summarized in Table 23. Tables 24 through 64 provide algorithms and parameters for each pathway.

The estimated intakes calculated for each exposure pathway considered and each COPC were compared to RfDs for non-carcinogenic effects. The following formula was used to estimate the potential for non-carcinogenic health effects for each COPC.

$$HQ = ADI/RfD \quad [Equation 6]$$

where:

HQ = Hazard quotient - potential for noncancer health effects (unitless);  
ADI = Average daily intake of COPC (mg/kg-day); and  
RfD = Reference dose (mg/kg-day).

RfDs have been developed by the US EPA for chronic (*e.g.*, lifetime) and/or subchronic exposure to constituents based on the most sensitive non-carcinogenic effects. The chronic RfD for a constituent is an estimate of a lifetime daily exposure level for the human population, including sensitive subpopulations, that is likely to be without an appreciable risk of deleterious effects. The potential for noncancer health effects was evaluated by comparing the Site-specific exposure level with the RfD derived by the US EPA for a similar exposure period. This ratio of exposure to toxicity is called the hazard quotient (HQ). If the Site-specific exposure level exceeds the threshold (*i.e.*, the HQ exceeds a value greater than 1.0), there may be concern for potential noncancer effects.

To assess the overall potential for noncancer effects posed by multiple constituents, a hazard index (HI) is derived by summing the individual HQs. This approach assumes additivity of critical effects of multiple constituents. This is appropriate only for compounds that induce the

same effect by the same mechanism of action. This conservative approach significantly overestimates the actual potential for adverse health impacts.

In cancer risk assessment, the US EPA has required the use of the upper limit which produces an estimate of potential risk that has a 95% probability of exceeding the actual risk, which may, in fact, be zero. The following formula was utilized to estimate the upper bound excess cancer risk for each carcinogen (note that not all COPCs are carcinogens):

$$TR = CLDI \times SF \quad [Equation 7]$$

where:

- TR = Target risk - excess probability of an individual developing cancer (unitless);  
CLDI = Calculated lifetime average daily intake of carcinogenic COPC (mg/kg-day); and  
SF = Cancer slope factor (mg/kg-day)<sup>-1</sup>.

For exposures to multiple carcinogens, the upper limits of cancer risks are summed to derive a total cancer risk. The US EPA recognizes that it is not technically appropriate to sum upper confidence limits of the risk to produce a realistic total probability, but requires this approach be used.

Carcinogenic risk refers to the probability of developing cancer as a result of exposure to known or suspected carcinogens. The National Contingency Plan (NCP) endorses an acceptable risk range of  $10^{-4}$  to  $10^{-6}$  for exposure to multiple carcinogens. This range represents an incremental increase of 1 in 10,000 to 1 in 1,000,000 in the chance of developing cancer over a lifetime. The MCEQ (1999) indicates that the target risk level is  $1 \times 10^{-6}$  per individual carcinogen and a cumulative risk level of  $1 \times 10^{-4}$ . As such, risk levels totaled across oral, dermal, and inhalation pathways may exceed  $1 \times 10^{-6}$  and still be in compliance with MCEQ requirements (1999) as long as no single carcinogen exceeds  $1 \times 10^{-6}$  and the cumulative risk for a single receptor does not exceed  $1 \times 10^{-4}$ .

Table 23 provides a summary of the non-carcinogenic effects and carcinogenic risks associated with each of the pathways evaluated in this assessment.

The overall hazard index across the assessed pathways and EUs was 0.02 for the Site visitor scenario. This value is below the acceptable benchmark of 1.0. The highest hazard index associated with the Site visitor scenario was 0.08 corresponding to oral exposure to sediment in EU4. The overall cancer risk for exposures to Site visitors was estimated to be  $3 \times 10^{-6}$  however, this value is a summation of the risks of several carcinogens as no single carcinogen exceeded  $1 \times 10^{-6}$ . The highest cancer risk associated with the infrequent Site visitor scenario was  $9 \times 10^{-7}$  and is attributable to dermal exposure to benzo(a)pyrene in EU4 surface soils. This value is below the *de minimis* benchmark of  $1 \times 10^{-6}$ .

The overall hazard index for the maintenance worker scenarios was  $2 \times 10^{-2}$  and is well below the acceptable benchmark of 1.0. The highest hazard index associated with the maintenance worker scenario was 0.01 corresponding to oral exposure to sediment in EU4 and oral exposure to surface soil in EU 4. The overall cancer risk for the maintenance worker scenario was  $1 \times 10^{-5}$ , within US EPA's acceptable range of  $1 \times 10^{-4}$  to  $1 \times 10^{-6}$ . The cancer risk level for the maintenance worker was primarily attributable to dermal exposure to benzo(a)pyrene in surface soils in EU 4 and EU 5 and sediments in EU 4 . Additional discussion regarding remediation goals for this scenario has been provided in section 8.0.

The overall hazard index for the hypothetical future construction worker was  $3 \times 10^{-4}$  and is over three orders of magnitude below the acceptable benchmark of 1.0. The highest hazard index associated with the construction worker scenario was  $2 \times 10^{-4}$  corresponding to dermal exposure to surface and subsurface soils in EU 4. The overall cancer risk for the hypothetical future construction worker scenario was  $9 \times 10^{-7}$ . This value is below the *de minimis* benchmark of  $1 \times 10^{-6}$ .

The off-Site resident scenario revealed a hazard index of  $8 \times 10^{-5}$ . This value is considerably below the acceptable benchmark of 1.0. The overall cancer risk for the resident exposure scenario was estimated to be  $2 \times 10^{-6}$  however, this value is based on the summation of the levels of several carcinogens as no single carcinogen was estimated to exceed  $1 \times 10^{-6}$ . The highest cancer risk associated with the Site visitor scenario was  $9 \times 10^{-7}$ .

The scenario that drove the cancer risk level at the Site was maintenance worker dermal exposure to benzo(a)pyrene in EU 4 and EU 5. These scenarios exceeded a *de minimis* risk level of  $1 \times 10^{-6}$ , but were within US EPA's acceptable risk range of  $1 \times 10^{-4}$  to  $1 \times 10^{-6}$ .

## 6.1 Assessment of the Ingestion Path of Carcinogenic PAHs

### 6.1.1 Background

Under Subpart II of the Risk Evaluation Procedures, MCEQ (1999) states that US EPA's published cancer slope factor utilized to quantitatively estimate risks due to dermal absorption of carcinogenic PAHs cannot be used to quantify carcinogenic risks from incidental ingestion of PAHs. Rather, the guidance states that such potential ingestion risk shall be described qualitatively. This section addresses the qualitative added risk from this exposure route and also provides useful perspective concerning typical daily exposures to PAH compounds by the general populace.

PAHs are ubiquitous in the environment and occur from both natural sources and human-related activities. Microbial synthesis, higher plant synthesis, volcanic activity, and prairie and forest fires are all major contributors to the natural background levels of PAHs (ATSDR, 1999; Suess, 1976; Andelman *et al.*, 1980; and Youngblood *et al.*, 1975). However, the majority of PAHs in the environment is formed by anthropogenic sources. PAH mixtures are found in tar, soot, petroleum, tobacco smoke, automotive exhaust, engine lubricant wastes, asphalt, fried and broiled foods, smoked meats, etc., and are generally formed by the incomplete combustion of organic material. PAHs are present in ambient air, water, sediments, soils, and foods (ATSDR, 1999).

In the environment, PAHs almost always occur as complex mixtures of many compounds made up of three or more fused benzene rings in linear, angular, or cluster arrangements and contain only carbon and hydrogen (Edwards, 1983). The most studied member of the class is B(a)P. The use of B(a)P as a model PAH is so prevalent that PAHs and B(a)P are often thought of synonymously. B(a)P may be as much as three orders of magnitude more potent than certain other carcinogenic PAHs (Chu and Chen, 1984, as cited in ATSDR, 1999). It should be emphasized that B(a)P constitutes only 1-20% of the total PAHs usually found in a complex environmental mixture (Blumer, 1961). Menzie *et al.* (1992), report that B(a)P represents about 10 to 30% of carcinogenic PAHs in a mixture. Nevertheless, because B(a)P is well-characterized and many PAHs in a mixture may not be identified in typical gas chromatography/mass spectrometry (GC/MS) analysis, some authors still report only B(a)P concentrations as representative of a complex mixture. Typically, total PAHs (*i.e.*, the five to 20 PAHs quantified in published reports) in soils are about 8 to 10 times the value of B(a)P alone (Edwards, 1983). Typical concentrations of B(a)P in various materials or media are summarized in the following table:

#### **Polycyclic Aromatic Hydrocarbon Content Of Various Materials And Media**

| Material                                | Benzo(a)pyrene ( $\mu\text{g}/\text{kg}$ ) | Reference                   |
|---|--|-----------------------------|
| Coal                                    | 4,000                                      | Coomes – 1981               |
| Asphalt                                 | 10,000 – 100,000                           | Coomes – 1981               |
| Raw shale oil                           | 3,200                                      | Coomes – 1981               |
| Crude oil                               | 400 – 2,800                                | Pancirov & Brown – 1975     |
| Coal tar pitch                          | up to 12,500,000                           | Wallcave – 1971             |
| Eastern US soils<br>(rural)             | 40 – 1,300                                 | Blumer – 1961               |
| Soil near Swiss<br>highway              | 2,100 – 30,000                             | Blumer <i>et al.</i> – 1977 |
| Fresh waters ( $\mu\text{g}/\text{L}$ ) | 0.001 – 0.100                              | Andelman & Suess – 1980     |
| Urban air (ng/m)                        | 0.03 – 104                                 | IARC – 1973                 |

| Material                     | Benzo(a)pyrene ( $\mu\text{g}/\text{kg}$ ) | Reference                  |
|------------------------------|--|----------------------------|
| Marine sediments             | nil – 5,000                                | Zobell – 1971              |
| Margarine                    | 0.2 – 6.8                                  | Swallow – 1976             |
| Sunflower oil                | 8.0  | Swallow – 1976             |
| Barbecued ribs               | 10.5                                       | Libinsky & Shubik – 1965   |
| Liquid smoke flavoring       | 9,400,000                                  | Youngblood & Blumer – 1975 |
| Barbecued beef               | 3.5  | Fretheim – 1976            |
| Tea leaves                   | 3.9 – 21.3                                 | IARC – 1973                |
| Lettuce<br>(industrial area) | 8.6 – 150                                  | Edwards – 1983             |

The main source of contamination of the soil by PAHs is by deposition from air. It has been estimated that global emissions of B(a)P during the time period between 1966 and 1969 were about 4.6 million kg (about 10 million pounds) per year, mostly from fossil fuel burning, with motor vehicles contributing only about 1% of the total (Suess, 1976, as cited in ATSDR, 1988). However, background concentrations of PAHs in surface soils near major highways tend to be relatively high. Blumer and coworkers (1975) showed that the major source of PAHs found in soils of a Swiss mountain town located in a deep valley with frequent temperature inversions was automobile exhaust (Blumer *et al.*, 1977); concentrations of PAHs in soils close to a major highway were found to be much greater than those in soils of the same locality but nearer to industries.

The PAHs emitted from jet aircraft exhaust provide clear illustration of the influence of emission source locations to background soil levels of PAHs. Shabad reported that a jet airplane engine emits 2000 to 4000  $\mu\text{g}$  B(a)P/min when cruising at high altitude, but emits amazingly up to 40,000,000  $\mu\text{g}$  B(a)P/min during take-off (Shabad, 1980, as cited in ATSDR, 1988). Not surprisingly then, soil, snow, and vegetation near airports were found to contain B(a)P (and PAHs) at levels an order of magnitude or more higher than at a control site (Smirnov, 1970 and Audere *et al.* 1973, as cited in ATSDR, 1988). On a global scale, however, amounts of PAHs

emitted from jet engine exhausts are very small compared to emissions from fossil fuel combustion and other sources (Edwards, 1983).

Carcinogenic PAHs are found in all surface soils. Evidence of the global distribution of PAHs was given by Thomas (1986, as cited in ASTDR, 1999) who detected carcinogenic PAHs in arctic soils above 150 µg/kg.

Although background levels of PAHs will vary depending on locality and proximity to emission sources, some generalizations regarding typical levels in soils, sediments, water, air, etc., have been reported. Blumer and coworkers examined surface soils in many areas of the eastern United States and noted that levels between 40 and 1,300 µg/kg in soils from relatively rural areas (Blumer, 1961; 1976). Typical concentrations of B(a)P in soils of the world according to Edwards range from about 100 µg/kg to 1000 µg/kg, with total PAHs about 10 times the value of B(a)P alone (Edwards, 1983).

In source dominated areas soils may contain much higher concentrations. Butler (1984) examined soils in Switzerland and demonstrated that much higher levels of PAHs, up to 300,000 µg/kg, are found near complex road interchanges than in areas more distant (Blumer *et al.*, 1977; Butler *et al.*, 1984). B(a)P concentrations as high as 650,000 µg/kg have been reported in soil samples collected 10 m from a German soot plant and as high as 120,000 µg/kg 500 m from a tar plant (Freitz, 1971).

As anticipated, urban areas have higher soil concentrations than do more remote areas because of the proximity to sources of fossil fuel combustion. The majority of urban soil concentrations of carcinogenic PAH concentrations falls in the 600 to 3,000 µg/kg range (Menzie *et al.*, 1992). The highest ambient concentrations of carcinogenic PAHs in soils have been reported for road dust, which can contain levels of 8,000 to 336,000 µg/kg (Menzie *et al.*, 1992).

Sediments are major sinks for PAHs, primarily because of the low solubility of these compounds and their strong affinity for organic carbon in particulate matter. For example, sediment samples taken from the Detroit River in 1982 contained B(a)P at concentrations ranging from 120 µg/kg to 17,640 µg/kg (Health & Welfare Canada, 1979).

### 6.1.2 Qualitative Assessment of Incidental Ingestion of PAHs in Soil

#### 6.1.2.1 Bioavailability

Incidental ingestion of soils containing PAHs at the Site by maintenance workers, construction workers, and trespassers is possible. Therefore, additional risk from this route of exposure is probable. However, risks calculated due to potential dermal exposures to soils and sediments contaminated with PAHs, while within US EPA's acceptable risk range of  $1\times10^{-4}$  to  $\times10^{-6}$ , is likely overestimated. This is because the bioavailability of organic pollutants in soil declines with time. However, current analytical methods measure total and not bioavailable concentrations of these constituents in soil and sediments. Research has provided ample evidence that the availability of certain chemicals that have been in soil for some time is less than freshly added compounds, and hence the term aging or weathering has been applied to the phenomenon. Awareness is growing among environmental toxicologists, risk assessors, and regulatory agencies that the total concentration of a toxicant in a contaminated environment frequently overestimates the risk of pollutants to humans, animals and plants.

This phenomenon has been recognized for some time. Early toxicological studies demonstrated the time-dependent diminution in bioavailability. For example, Edwards *et al.* in 1957 (as cited in Alexander, 2000) reported that simultaneous biological assays of acute toxicity to the fruit fly, *Drosophila melanogaster*, and chemical concentration measurements gave similar results shortly after lindane was applied to soil, but much of the insecticide remaining in the soil after 22 months did not detectably affect the fruit flies. More recent research supports these observations. Typically, the bioremediation of soils containing PAHs, although reducing the concentration of many individual compounds, does not completely rid the treated site of PAHs that are known to be degraded in microbial cultures. The microorganisms are present, the environmental

conditions are conducive to their activity, but somehow the compounds become less and less accessible with time. The view that the contaminants became sequestered as they reside in the field gains credence in light of the finding that biodegradation of the seemingly resistant PAHs takes place if they are extracted and then added back to soil (Weissenfels *et al.*, 1992, as cited in Alexander, 2000).

Aging of PAHs and many other chemicals is toxicologically significant because the toxicity of harmful compounds declines substantially as they persist and become increasingly sequestered with time (Alexander, 2000). Recently, it has been demonstrated that the genotoxicity of the carcinogens benzo(*a*)pyrene and dimethyl-benzanthracene diminished rapidly and to a great extent within a 15-day period, even though standard analysis following vigorous extraction of soil showed only a very slight decline in concentration after about 2 months (Alexander and Alexander, 1999). Additional experimental evidence for the progressive decline in bioavailability with time has been obtained in tests with the earthworm, *E. fetida*. With the progress of aging, decreasing quantities of the PAHs naphthalene, phenanthrene, atrazine, anthracene, fluoranthene and pyrene were assimilated by the animals (Kelsey and Alexander, 1997 as cited in Alexander, 2000; Tang *et al.*, 1998, as cited in Alexander, 2000). The decline in availability of DDT and DDE to worms may even reach values of 90% (Morrison *et al.*, 2000). Yet the types of chemical analysis used for assessing risk and making regulatory decisions fail to show the parallel diminution in bioavailability and attendant toxicity.

Chemicals that leach through soil may pose a risk to humans who obtain drinking water from impacted downgradient wells (though this is not the case downgradient of the Hattiesburg Site). Because aging reduces the amounts of various compounds that are desorbed and may then pass through soil, it also indirectly affects potential human exposure (White *et al.*, 1997). The observation that the time-dependent decline in bioavailability is accompanied by a time-dependent decline in the quantity of compounds extracted from soil by a mild procedure suggests that a mild-extraction technique might serve as a better basis for a surrogate assay for

bioavailability (Chung and Alexander, 1998; Kelsey *et al.*, 1997 as cited in Alexander, 2000; Tang and Alexander, 1999 as cited in Alexander, 2000).

The analytical protocols used at the Hattiesburg Site rely on vigorous extraction techniques that fail to predict the bioavailability of the PAH compounds that have been aged in Site soils for many years. For the purposes of the risk assessment (e.g., direct dermal contact and dermal absorption with soils containing PAHs), it was assumed that virtually all the PAH compounds present in the soil matrix were available for absorption. This is clearly not likely to be the case and, accordingly, risks have been overestimated for this route of exposure.

Because PAHs are ubiquitous, humans are exposed to these chemicals as part of everyday living. PAHs have been detected in many food products, particularly leafy vegetables, especially those grown in urban areas or near roadways. Even smoking "one pack of cigarettes a day," which has been estimated to result in a B(a)P exposure of 0.4 µg/kg (ATSDR, 1988), is less than the typical intake via the diet. Santodonato *et al.* (1981) reviewed the literature on human exposure including dietary doses of PAHs and summarized the findings as follows.

| Source | B(a)P Concentration | Carcinogenic PAH Concentration | Total PAH Concentration |
|--------|---------------------|--------------------------------|-------------------------|
| Air    | 0.0095 – 0.0435 µg  | 0.038 µg                       | 0.207 µg                |
| Water  | 0.0011 µg           | 0.0042 µg                      | 0.0270 µg               |
| Food   | 0.16 – 1.6 µg       | no data                        | 1.6 – 16 µg             |

Statistics available from the United States Department of Agriculture (Pao *et al.*, and United States Department of Agriculture, 1986) were used to determine an average diet for US males. For the average American diet, the intake of carcinogenic PAHs was estimated to be between 1 and 5 µg/day. Menzie *et al.* (1992) reported comparable results for total potential dose of carcinogenic PAHs for nonsmoking adult males, and estimated a median of 3 µg/day and a

maximum of 15 µg/day. According to these authors, this would translate to a daily dietary dose of B(a)P of 0.6 to 3.0 µg.

It should be noted that the available monitoring data are very limited with regard to dairy products, such as cheese and milk (Santodonato *et al.*, 1981), which make up approximately one-third of the daily diet of 1600 g. In one study conducted in Canada, B(a)P concentrations of 7.6 to 387 µg/L, with a mean value of 129.5 µg/L were found in samples of milk from ten nursing mothers (HSDB, 1999). Assuming ingestion of one liter per day of mother's milk by a nursing baby, the intake of B(a)P alone would, on average, be about 130 µg.

To provide qualitative perspective with regard to the incidental ingestion of soil at the Hattiesburg Site, if a maintenance worker were to ingest 100 mg of soil per day, 150 days per year, with an average B(a)P concentration of 3.17 mg/kg in the former Process Area, the average daily intake of B(a)P would be about 0.13 µg. This appears to be comparable to or less than the typical dietary intake of B(a)P from food alone. A charcoal-broiled T-bone steak reportedly contains about 50 µg/kg of B(a)P, as well as various levels of other PAHs (Lijinsky and Ross, 1977, as cited in HSDB, 1999). For a 10-ounce portion, this translates to a dose of 14 µg B(a)P from this food item alone. It may be noted here also that, with few exceptions, PAHs in soil in the Process Area are covered with asphalt, and therefore, not readily available for direct contact and incidental ingestion.

## **7.0 Uncertainty Analysis**

Risk assessment uses a wide array of information sources and techniques. Even in those rare circumstances where constituent intake for an exposed individual may be measured relatively precisely, assumptions will still be required to evaluate the associated risk. Generally, data are not available for critical aspects of the risk assessment, and the use of professional judgment, inferences based on analogy, the use of default values, model estimation techniques, etc., result in uncertainty of varying degrees.

The expressions of risk in this assessment are not probabilistic; the expressions of risk are conditional, based on the conditions represented by the single-point values selected for the analysis. This section is intended to identify and qualitatively evaluate the more salient Site-specific uncertainties and their potential influence on the credibility of the estimated Site risks.

### **7.1 Uncertainty of Data Evaluation Factors**

Uncertainties in data analysis include analytical error, selection of COPCs, adequacy of sampling design, etc. Generally, there is far less uncertainty in this phase of the risk assessment process than other aspects contribute.

Laboratory analysis is extremely accurate relative to the potential error of "professional judgment" in exposure assessments. The uncertainty of analytical data is likely to be less than 25 percent, most of the time.

The adequacy of the sampling strategies to characterize Site conditions is a potentially large source of uncertainty. Because of the limited availability of resources, sample collection is generally limited. However, sampling (especially in multiple surveys) is not random, but is designed to locate the areas with the highest levels of constituents. Thus, test data are biased toward overestimation of average constituent levels. In addition, in most instances, the upper 95-percent confidence limit of the average concentration is utilized as an exposure-point

concentration in the risk assessment. The use of this value likely will result in an overestimation of risk, as the 95% UCL represents a value that will be greater than the true average 95% of the time.

Oftentimes, only a portion of detected constituents are carried through the risk assessment process because constituents are eliminated through COPC screening procedures (US EPA, 1989). This could result in an underestimation of risk, although the COPC selection process is intended to identify those constituents which account for the vast majority of potential risk. COPCs lacking published RfD values were not quantitatively evaluated and this may result in an underestimation of potential hazards (non-carcinogenic effects).

## 7.2     Uncertainty of Toxicity Values

The US EPA's IRIS states that the uncertainty associated with RfD values for non-carcinogenic endpoints of toxicity "span perhaps an order of magnitude." In fact, the uncertainty of extrapolating dose-response data from animals to humans with the application of multiple safety factors (100 to 10,000 or more) is likely to be several orders of magnitude. Current policies for deriving RfD values will often result in an overestimation of risk.

The uncertainty associated with the estimation of cancer risk contributes, by far, the major source of potential error and uncertainty. It is beyond the scope of this analysis to explore this toxicity assessment factor in any detail. However, a few salient points are noted below.

Some constituents classified as carcinogens have been shown to produce an increased incidence of cancer in mice but not rats, for example. If the mouse is not an adequate model for the rat, it may be wondered how reliable a model it is for human beings. The assumption of linearity and a non-threshold phenomenon in the dose versus risk relationship may not be valid and could result in a very large overestimation of actual cancer risk, if any even exist at low doses in humans.

The US EPA evaluated the uncertainty of cancer risk estimates from exposures to trichloroethene and several other related VOCs in public drinking water supplies (Cothern *et al.*, 1984). These US EPA scientists concluded the following:

- The largest uncertainty in the calculations is due to the choice of the model [Multistage, Weibull, Logit, Probit, etc.] used in extrapolating risk to low doses in humans, and is 5 to 6 orders of magnitude;
- If a single model were chosen [assumed to be valid], the overall uncertainty in risk estimates would be 2 to 3 orders of magnitude;
- The exposure estimates contribute, at most, an order of magnitude to the uncertainty; and
- It would appear that until a particular compound's mechanisms of cancer are better known, it is likely that the uncertainty in the toxicity will not be improved.

### 7.3 Uncertainties in Assessing Potential Exposure

Ideally, Site-specific exposure values should be used when assessing potential intakes of chemicals at a Site. Oftentimes, however, Site-specific data are not available; therefore, the risk assessor must estimate values that most accurately reflect Site conditions. In doing so, US EPA or other regulatory default values were utilized in place of Site-specific data. These values may over- or under-estimate risks, depending on Site conditions and the percentile range in which the default values fall (*e.g.*, 50<sup>th</sup>, 95<sup>th</sup>).

Although a considerable amount of published data is available on the most common exposure parameters (*e.g.*, body weight, skin surface area), even these data contain uncertainties. Studies conducted by different scientists often provide differing levels of detail, statistics, and accuracy based on sample size, study design, geographic area, etc. Such discrepancies can increase uncertainty when the data are combined to derive a single-point default value. These data may be the best available; however, the reflection of reality may still be imprecise.

Where published exposure parameters were not available, best professional judgment had to be used, thereby increasing uncertainty. The default or estimated exposure parameters used in this assessment likely resulted in a moderate over-estimation of risk.

MCEQ (1999) guidance recommends a qualitative evaluation of carcinogenic risks resulting from ingestion of carcinogenic PAH compounds. By not including carcinogenic PAH ingestion in the quantitative evaluation, potential risks are clearly underestimated.

Carcinogenic PAHs are thought to increase the risk of cancer by all routes of exposure. Accordingly, the risks estimated by dermal absorption and inhalation exposure may underestimate the upper bound of cancer risk from additional exposure via incidental ingestion of soil at the Site. However, the risks estimated for dermal absorption of PAHs adsorbed into soils adhering to skin may overestimate risks for a host of reasons.

Early studies conducted by Falk and coworkers indicated that the carcinogenic effect of B(a)P on subcutaneous injection in mice could be markedly inhibited by the simultaneous administration of various non-carcinogenic PAHs (Falk *et al.*, 1964, as cited in ATSDR, 1988. In other subcutaneous injection and skin-painting studies with mice, it was shown that a combination of several non-carcinogenic PAH compounds, mixed according to the proportion occurring in auto exhaust, did not enhance or inhibit the action of two potent PAH carcinogens, B(a)P and dibenz(a,h)anthracene.(ATSDR, 1988).

The carcinogenic potency of B(a)P and other carcinogenic PAHs is generally determined by injecting solutions under the skin, painting the skin with the carcinogenic PAH dissolved in a solvent, or dissolved in corn oil in feeding studies. This vehicle or matrix affords a high level of bioavailability of the carcinogenic PAH compound. Recently, Krueger *et al.* (1999) conducted *in vitro* percutaneous absorption studies with contaminated soils and organic solvent extracts of contaminated soils collected at former manufactured gas plant (MGP) sites. The MGP tar-contaminated soils contained PAHs at levels ranging from 10 to 2400 mg/kg. The dermal

penetration rates of PAH from the MGP tar-contaminated soils and soil solvent extracts were determined experimentally through human skin using tritium-labelled B(a)P as a surrogate. Results showed reductions of two to three orders of magnitude in PAH absorption through human skin from the most contaminated soils in comparison to the soil extracts. Reduction in PAH penetration was attributed to soil matrix properties. That is, PAH compounds adsorbed to organic carbon in a soil matrix are far less bioavailable for dermal flux than PAH compounds dissolved in a solvent. [No correction for such a profound soil matrix effect was applied in quantitatively estimating cancer risks due to dermal absorption of B(a)P and other carcinogenic PAHs in this assessment.]

## **8.0 Summary of Findings**

The results of the baseline human health risk assessment indicate potentially unacceptable risk values for the maintenance workers dermally exposed to sediment and surface soils in EU 4 and surface soils in EU5 as a result of the presence of benzo(a)pyrene. To determine the extent of remediation necessary to reduce these risks to acceptable levels, Site data in EU 4 and EU 5 were closely examined.

The benzo(a)pyrene exposure-point concentration used to evaluate maintenance worker exposures to sediment in EU 4 was 130 mg/kg (sample location SD02). This was the maximum benzo(a)pyrene concentration found in sediment in EU 4. The next highest concentration of benzo(a)pyrene in sediment was found at SD12 (71 mg/kg). Implementing a remedy to remove, treat, or preclude contact with sediment at sample location SD02 would leave a concentration of 71 mg/kg (sample location SD12) as the maximum concentration in sediment that could be potentially contacted by a maintenance worker in EU 4. Excluding sample SD02 and using 71 mg/kg as the exposure-point concentration drops the risk level for dermal contact with sediment by a maintenance worker to within acceptable levels ( $8 \times 10^{-7}$ , Table 65). Thus, implementing a remedy to address sediment at sample location SD02 would result in acceptable levels of risk for maintenance workers exposed to sediments in EU4.

In EU 4, the maximum concentration of benzo(a)pyrene was used as the exposure-point concentration for maintenance worker surface soil exposures. The maximum concentration was 500 mg/kg at sample location GEO-48/0-1' (Figure 2). The next two highest concentrations of benzo(a)pyrene in surface soil were 230 mg/kg (GEO-21/0-1') and 190 mg/kg (GEO-21/0-1'). These sample locations are presented on Figure 2. Implementing a remedy to remove, treat, or preclude contact with soils at sample locations GEO-48 and GEO-21 would leave a concentration of 56 mg/kg (sample location GEO-19/0-1') as the maximum concentration that could be potentially contacted by a maintenance worker in EU 4. Excluding samples GEO-48 and GEO-21 and using 56 mg/kg as an exposure point concentration, drops the risk level for dermal contact with surface soil by a maintenance worker to within acceptable levels ( $6 \times 10^{-7}$ ,

Table 66). Thus, implementing a remedy to address soils at sample locations GEO 48/0-1', GEO-21/0-1', and GEO-21/2-3' would result in acceptable levels of risk for maintenance workers exposed to surface soils in EU4.

In EU5, the surface sample location contributing most to elevated risk levels for the maintenance worker scenario was GEO-33/0-1' (see Figure 2). The benzo(a)pyrene concentration at this location (52.5 mg/kg) contributed most to the maintenance worker risk estimates. Sample location GEO-33/0-1' is located within a paved area in a parcel of land southwest of Courtesy Ford (Figure 2). Pavement in this area precludes direct contact with surface soils; therefore, it is not anticipated that current or future maintenance workers will have access to surface soils in or around sample location GEO-33/0-1'.

For the maintenance worker scenario, elimination of contact with the maximum concentration of benzo(a)pyrene in surface soil in EU 5 would reduce risk levels in EU5 to below the *de minimis* benchmark of  $1 \times 10^{-6}$ . Existing pavement already precludes direct contact with surface soils at location GEO-33/0-1'. The risk level associated with dermal exposure to benzo(a)pyrene was recalculated excluding the maximum benzo(a)pyrene concentration sample location in EU5 (52.5 mg/kg at GEO 33/0-1'). A reevaluation of risks is presented in Table 67. Results of this evaluation indicated risk levels below the *de minimis*  $1 \times 10^{-6}$  benchmark.

## Bibliography

- Andelman, J. B., and M. J. Suess. 1980. Polynuclear aromatic hydrocarbons in the water environment. Bull. WHO 43:479-508.
- Alexander, M. Aging, Bioavailability, and Overestimation of Risk from Environmental Pollutants. Environ. Sci. Technol, 2000, 34(20):4259.
- ATSDR (Agency for Toxic Substances and Disease Registry). Toxicological Profile for Benzo(a)Pyrene. Oak Ridge National Laboratory. 1988.
- ATSDR (Agency for Toxic Substances and Disease Registry). ATSDR's Toxicological Profiles on CD-ROM. Polycyclic Aromatic Hydrocarbons (PAHs), Update. CRC Press, 1999.
- Audere, A. K., Z. Y. Lindberg, G. A. Smirnov, and L. M. Shabad. 1973. Experiment in studying the influence of an airport located within the limits of a city on the level of environmental pollution by benzo(a)pyrene. Gig. Sanit. 38(9): 90-92.
- Bartek, M.J. and J.A. LaBudde. Percutaneous Absorption *in vitro*, in Animal Models in Dermatology. Ed. H.I. Maibach. New York: Churchill Livingstone, 1975. p. 103.
- Blumer, M. 1961. Benzopyrenes in soil. Science 134, 474-475.
- Blumer, M., W. Blumer, and T. Relch. 1977. Polycyclic aromatic hydrocarbons in soils of a mountain valley; correlation with highway traffic and cancer incidence. Environ. Sci. Technol. 11 (12), 1082-1084.
- Butler, J. D., V. Butterworth, C. Kellow, and H. G. Robinson. 1984. Some observations on the polycyclic aromatic hydrocarbon (PAH) content of surface soils in urban areas. Sci. Total Environ. 38, 75-85.
- Chu, M. M. L. and G. W. Chem. 1984. Evaluation and estimation of potential carcinogenic risks of polynuclear aromatic hydrocarbons. Paper presented at the Pacific Rim Risk Conference.
- Chung, N. and Alexander, M. Differences in Sequestration and Bioavailability of Organic Compounds Aged in Dissimilar Soils. Environ Sci. Technol. 32: 855.
- Coomes, R. M. 1981. Carcinogenic testing of oil shale materials. Twelfth Oil Shale Symposium Proceedings. Colorado School of Mines Pres.

Cothen, C. R., W. Conniglio, W. Marcus. Techniques for the Assessment of the Carcinogenic Risk to the US Population due to Exposure from Selected Volatile Organic Compounds from Drinking Water via the Ingestion, Inhalation and Dermal Routes. NTIS PB84-213941. Office of Drinking Water. Washington DC: Environmental Protection Agency, 1984.

Edwards, C.A., Beck, S.D. and Lichtenstein, E.P., J. Econ. Entomol. 1957, 50: 622.

Edwards, N. T. 1983. Polycyclic aromatic hydrocarbons (PAHs) in the terrestrial environment – a review. J. Environ. Qual. 12 (4), 427-441.

Falk, H. L., and P. T. S. Kotin. Inhibition of carcinogenesis: The effects of polycyclic hydrocarbons and related compounds. Arch. Environ. Health Vol. 9 (1964):169-179.

Fritz, W. 1971. Extent and sources of contamination of our food with carcinogenic hydrocarbons. Ernaehrungsforschung 16(4), 547-557.

Health & Welfare Canada. Polycyclic Aromatic Hydrocarbons, Report No. 80-EHD-50, (1979) p. 38.

Horst, T. W. Langrangian Similarity Modeling of Vertical Diffusion for a Ground Level Source. Int. Applied Met, Vol. 18 (1979): 733-740.

HSDB (Hazardous Substances Data Bank), 1999. National Library of Medicine (NLM) On-Line Toxicological Network (TOXNET). Bethesda, MD.

ICRP (International Commission on Radiological Protection). Report of Committee IV on Evaluation of Radiation Doses to Body Tissues from Internal Contamination due to Occupational Exposure. ICRP Publication 10. New York: Pergamon Press, 1968.

Kao, J.K., F.K. Patterson, and J. Hall. Skin Penetration and Metabolism of Topically Applied Chemicals in Six Mammalian Species, Including Man: an in vitro Study with Benzo(a)pyrene and Testcaterone. Toxicol. Appl. Pharmacol, Vol. 81 (1985): 502-516.

Kelsey, J.W. and Alexander, M. Declining Bioavailability and Inappropriate Estimation of Risk of Persistent Compounds. Environ. Toxicol. Chem. 1997. 16(3): 582

Kelsey, J.W., Kottler, B.D. and Alexander, M. Environ. Sci. Technol. 1997, 31: 214.

Kissel, J., K. Richter, and R. Fenske. Field Measurements of Dermal Soil Loading Attributable to Various Activities: Implications for Exposure Assessment. Risk Analysis, Vol. 16, No. 1 (1996): 115-125.

Magee, B., P. Anderson, and D. Burmaster. Absorption Adjustment Factor (AAF) Distributions for Polycyclic Aromatic Hydrocarbons (PAHs). Human and Ecological Assessment: An International Journal. Vol. 2, No. 4 (December 1996): 841-873.

Menzi, C.A., B.B. Potocki and J. Santodonato. Exposure to Carcinogenic PAHs in the Environment. Environ. Sci. Technol. Vol. 26, No. 7, 1992.

Michael Pisani & Associates. Remedial Investigation Report, Former Gulf States Creosoting Site, Hattiesburg, Mississippi. New Orleans, Louisiana. 1997.

Michael Pisani & Associates. Phase II Remedial Investigation Report, Former Gulf States Creosoting Site, Hattiesburg, Mississippi. New Orleans, Louisiana. 1998.

Mississippi Commission on Environmental Quality (MCEQ). Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi. 1999.

Mississippi Department of Environmental Quality (MDEQ). Guidance for Remediation of Uncontrolled Hazardous Substance Sites in Mississippi. Office of Pollution Control. 1990.

Mississippi Department of Environmental Quality (MDEQ). Letter to Glen Pilie, Adams and Reese, from Tony Russell, MDEQ. August 2, 2000.

Morrison, DE., Robertson, B.K. and Alexander, M. Bioavailability to Earthworms of Aged DDT, DDE, DDD, and Dieldrin in Soil. Environ. Sci. Technol. 2000 34: 709.

Moore, M.R., P.A. Meredith, W.S. Watson, D.J. Sumner, M.K. Taylor, and A. Goldberg. "The Percutaneous Absorption of Lead-203 in Humans From Cosmetic Preparations Containing Lead Acetate, as Assessed by Whole-Body Counting and Other Techniques." Food Cosmet. Toxicol. 18. (1980): 399.

Pancirov, R. J. and R. A. Brown. 1975. Analytical methods for polynuclear aromatic hydrocarbons in crude oil, heating oils, and marine tissues. In: Conference on prevention and control of oil pollution, San Francisco, CA, March, 1975. American Petroleum Institute, Wash., DC. pp 103-13.

Pao, E. M. *et al.* Home Economics Research Report No. 44. United States Department of Agriculture, Washington, DC. 1982.

Pasquill, I.. The Dispersion of Material in the Atmospheric Boundary Layer - The Basis for Generalization. In: Lectures on Air Pollution and Environmental Impact Analysis. Boston, MA: American Meteorological Society, 1975.

Ryan, E.A., E.T. Hawkins *et al.* "Assessing Risk From Dermal Exposure at Hazardous Waste Sites. in Bennett." Ed. G. and J. Bennett. Superfund '87: Proceedings of the Eighth National Conference. Washington, DC, 16-18 November 1987. The Hazardous Material Control Research Institute. p.166-168.

Santodonato, J., P. Howard, and D. Basu. Health and Ecological Assessment of Polynuclear Aromatic Hydrocarbons. Pathotox Publishers, Inc., Park Forest South, IL. 1981.

Shabad, L. M. 1980. Circulation of carcinogenic polycyclic aromatic hydrocarbons in the human environment and cancer prevention. *J. Natl. Cancer Inst.* 64(3): 405-410.

Shu, H.P., P. Teitelbaum, A.S. Webb, L. Marple, B. Brunck, D. Dei Rossi, F.J. Murray, and D.J. Paustenbach. "Bioavailability of Soil Bound TCDD: Dermal Bioavailability in the Rat." *Fundam. Appl. Toxicol.*, Vol. 10 (1988): 648-654.

Smirnov, G. A. 1970. The study of benzo(a)pyrene content in soil and vegetation in the airfield region. *Vopr. Onkol.* 16(5): 83-86.

State of Mississippi. Mississippi Code 1972 Annotated. Title 29 Public Lands, Buildings, and Property, Chapter 3, Sixteenth Section and Lieu Lands in General. 1998.

Suess, M. J. 1976. The environmental load and cycle of polycyclic aromatic hydrocarbons. *Sci. Total Environ.* 6:239-250.

Ta, Roy *et al.*, Studies Estimating the Dermal Bioavailability of Polynuclear Aromatic Hydrocarbons from Manufactured Gas Plan Tar-Contaminated Soils. *Env. Sci. Tech.* 32(20). 1998. 3113-3117.

Tang, W.C., White, J.C., and Alexander, M. *Appl. Microbiol. Biotechnol.* 1998, 49: 117.

Tang, J and Alexander, M. *Environ. Toxicol. Chem.* 1999, 18: 2711.

Thomas, J.F., M. Mukai, and B.D. Teggens. Fate of airborne benzo(a)pyrene. *Environ. Sci. Technol.* 2:33-39, 1968.

United States Department of Agriculture (USDA). Nationwide Food Consumption Survey: Continuing Survey of Food Intakes by Individuals, Men 19-50 years, 1 Day, 1985; United States Department of Agriculture. Human Nutrition Information Service. Nutrition Monitoring Division; Washington, DC, Report No. 86-1. 1986.

US EPA (United States Environmental Protection Agency). Rapid Assessment of Exposure to Particulate Emission from Surface Contamination Sites. EPA/OHEA/EPA. 600/8-85/002. Cowherd, C., Jr., G.E. Muleski, P.J. Engelhart and D.A. Gillett, Ed. Washington DC: Midwest Research Inst.. 1985.

US EPA (United States Environmental Protection Agency). "Guidelines for Carcinogenic Risk Assessment." Federal Register 51:33992-34003. 1986.

US EPA (United States Environmental Protection Agency). Superfund Exposure Assessment Manual (SEAM). EPA 540/1-88/001. Office of Remedial Response. Washington, DC, 1988.

US EPA (United States Environmental Protection Agency). Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A). Interim Final. EPA/540/1-89/002. Office of Emergency and Remedial Response. Washington, DC, 1989.

US EPA (United States Environmental Protection Agency). Human Health Evaluation Manual, Supplemental Guidance: Standard Default Exposure Factors. OSWER Directive. 9285.6-03. Office of Solid Waste and Emergency Response. Washington, DC, 1991.

US EPA (United States Environmental Protection Agency). Dermal Exposure Assessment: Principles and Applications. Office of Research and Development. EPA/600/8-91/011B. Washington, DC, 1992.

US EPA (United States Environmental Protection Agency). Estimating Exposure to Dioxin-Like Compounds. Office of Research and Development. EPA/600/6-88/005B. Washington, DC, 1992.

US EPA (United States Environmental Protection Agency). Provisional Guidance for Quantitative Risk Assessment of Polynuclear Aromatic Hydrocarbons. Office of Solid Waste and Environmental Remediation. EPA/600/R-93/089, July 1993

US EPA (United States Environmental Protection Agency). Land Use Directive in the CERCLA Remedy Selection Process. OSWER Directive 9355.7-04. Office of Solid Waste and Emergency Response. Washington, DC, May 1995.

US EPA (United States Environmental Protection Agency) Region 3. Technical Guidance Manual: Risk Assessment, Assessing Dermal Exposure From Soil. EPA/903-K-95-003. Office of Superfund Programs, Hazardous Waste Management Division. Washington, DC, 1995.

US EPA (United States Environmental Protection Agency) Region 4. Technical Services Supplemental Guidance to RAGS: Region 4 Bulletins. Waste Management Division, Atlanta, GA. 1995.

US EPA (United States Environmental Protection Agency). Supplement B to Compilation of Air Pollutant Emission Factors, Volume I: Stationary Point and Area Sources. AP-42, Fifth Edition, Supplement B. Office of Air Quality Planning and Standards, Office of Air and Radiation. Research Triangle Park, NC, 1996

US EPA (United States Environmental Protection Agency). Exposure Factors Handbook., EPA/600/P-95/002F. Office of Research and Development, Washington, DC, August 1997.

US EPA (United States Environmental Protection Agency). Health Effects Assessment Summary Tables (HEAST). Office of Health and Environmental Assessment, Environmental Criteria and Assessment Office (ECAO). Cincinnati OH, 1997.

US EPA (United States Environmental Protection Agency). IRIS (Integrated Risk Information System). A Continuously Updated Electronic Database Maintained by the US Environmental Protection Agency. Bethesda, Maryland: National Library of Medicine, 1999.

US EPA (United States Environmental Protection Agency) Region 3. Updated Risk-Based Concentration Tables. Office of RCRA Technical & Program Support Branch. Philadelphia, PA, April, 1999.

Wallcave, L., H. Garcia, R. Fedlman, W. Linjinsky, and P. Shubik. 1971. Skin tumorigenesis in mice by petroleum asphalts and coal-tar pitches of known polynuclear aromatic hydrocarbon content. *Toxicol. Appl. Pharmacol.* 18, 41-52.

Weissenfels, W.D., Klewer, H.J. and Langhoff, J., *J. Appl. Microbiol. Biotechnol.* 1992, 36: 689.

White, J.C., Kelsey, J.W., Hatzinger, P.B and Alexander, M. Factors Affecting Sequestration and Bioavailability of Phenanthrene in Soils. *Environ. Toxicol. Chem.* 1997, 16(10): 2040.

Youngblood, W. W., and M. Blumer. 1975. Polycyclic aromatic hydrocarbons in the environment: homologous series in soils and recent marine sediments. *Geochim. Cosmochim. Acta* 39:1303-1315.

Figure 1

**Site Conceptual Model and Selection of Exposure Pathways**  
**Kerr McGee, Hattiesburg, MS**

|                    |                          |                    |                     |              |                        |                           |   |  |  | Rationale for Selection or Exclusion of Exposure Pathway |  |
|--------------------|--------------------------|--------------------|---------------------|--------------|------------------------|---------------------------|---|--|--|--|--|
| Scenario Timeframe | Medium                   | Exposure Point     | Receptor Population | Receptor Age | Exposure Route         | On-Site/Off-Site          | Type of Analysis  |  |  |  |  |
| Current            | Soil Surface Soil (0-1') | Exposure Unit 1    | Visitor             | Adolescent   | Dermal Oral Inhalation | On-Site                   | None  | EU1 includes only surface water and sediment. Soils in this area are included in EU2   |  |  |  |
|                    |                          | Exposure Unit 2    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | None  | EU1 includes only surface water and sediment. Soils in this area are included in EU2   |  |  |  |
|                    |                          | Exposure Unit 3    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | Quantitative Qualitative  | EU1 includes only surface water and sediment. Soils in this area are included in EU2   |  |  |  |
|                    |                          | Exposure Unit 4    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | Quantitative Qualitative  | Area potentially attractive for occasional recreational use<br>Only cPAHs were COPCs; addressed qualitatively according to MCEQ guidance |  |  |  |
|                    |                          | Exposure Unit 5    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | Quantitative Qualitative  | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                        |  |  |  |
|                    | Surface Soil (0-6')      | Maintenance Worker | Adult               | Dermal       | On-Site                | Quantitative Quantitative | Area potentially attractive for occasional recreational use<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                        |  |  |  |
|                    |                          | Maintenance Worker | Adult               | Dermal       | On-Site                | Quantitative Quantitative | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                       | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                        |  |  |  |
|                    |                          | Maintenance Worker | Adult               | Dermal       | On-Site                | Quantitative Quantitative | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                       | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                        |  |  |  |
|                    |                          | Maintenance Worker | Adult               | Dermal       | On-Site                | Quantitative Quantitative | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                       | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                        |  |  |  |
|                    |                          | Maintenance Worker | Adult               | Dermal       | On-Site                | Quantitative Quantitative | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                       | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                        |  |  |  |
| Sediment           | Sediment                 | Exposure Unit 1    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | Quantitative Qualitative  | EU1 includes only surface water and sediment. Soils in this area are included in EU2   |  |  |  |
|                    |                          | Exposure Unit 2    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | Quantitative Qualitative  | EU1 includes only surface water and sediment. Soils in this area are included in EU2   |  |  |  |
|                    |                          | Exposure Unit 3    | Resident            | Child/Adult  | Dermal Oral            | Off-Site                  | None  | EU1 includes only surface water and sediment. Soils in this area are included in EU2   |  |  |  |
|                    | Sediment                 | Exposure Unit 4    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | Quantitative Qualitative  | Area potentially attractive for occasional recreational use<br>Only cPAHs were COPCs; addressed qualitatively according to MCEQ guidance |  |  |  |
|                    |                          | Exposure Unit 5    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | Quantitative Qualitative  | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                        |  |  |  |
|                    |                          | Exposure Unit 6    | Visitor             | Adolescent   | Dermal Oral            | On-Site                   | Quantitative Qualitative  | VOCs not present at levels of concern<br>Only VOCs were COPCs; addressed qualitatively according to MCEQ guidance                        |  |  |  |



**Figure 1**

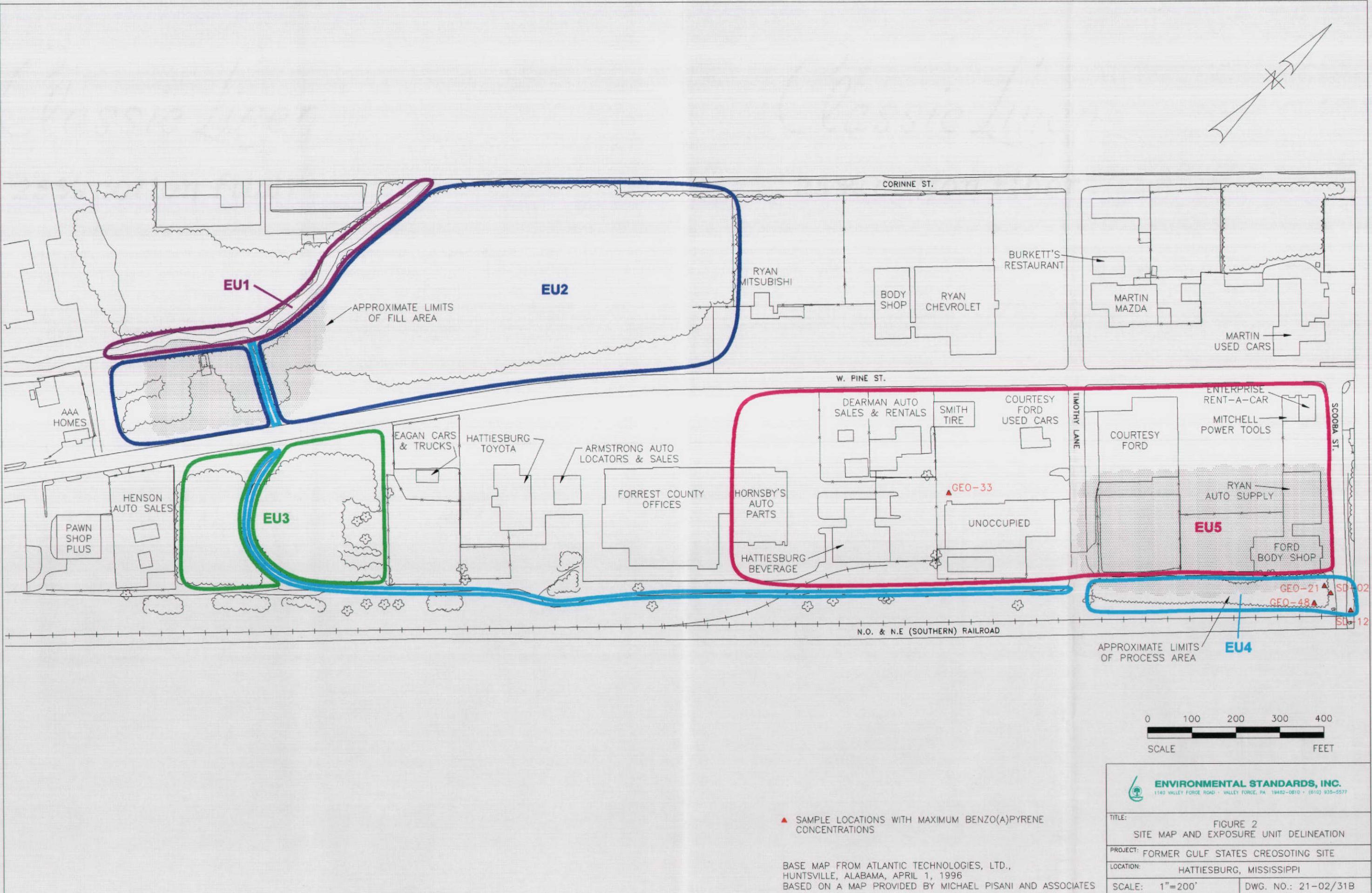
*Site Conceptual Model and Selection of Exposure Pathways*  
**Kerr McGee, Hattiesburg, MS**

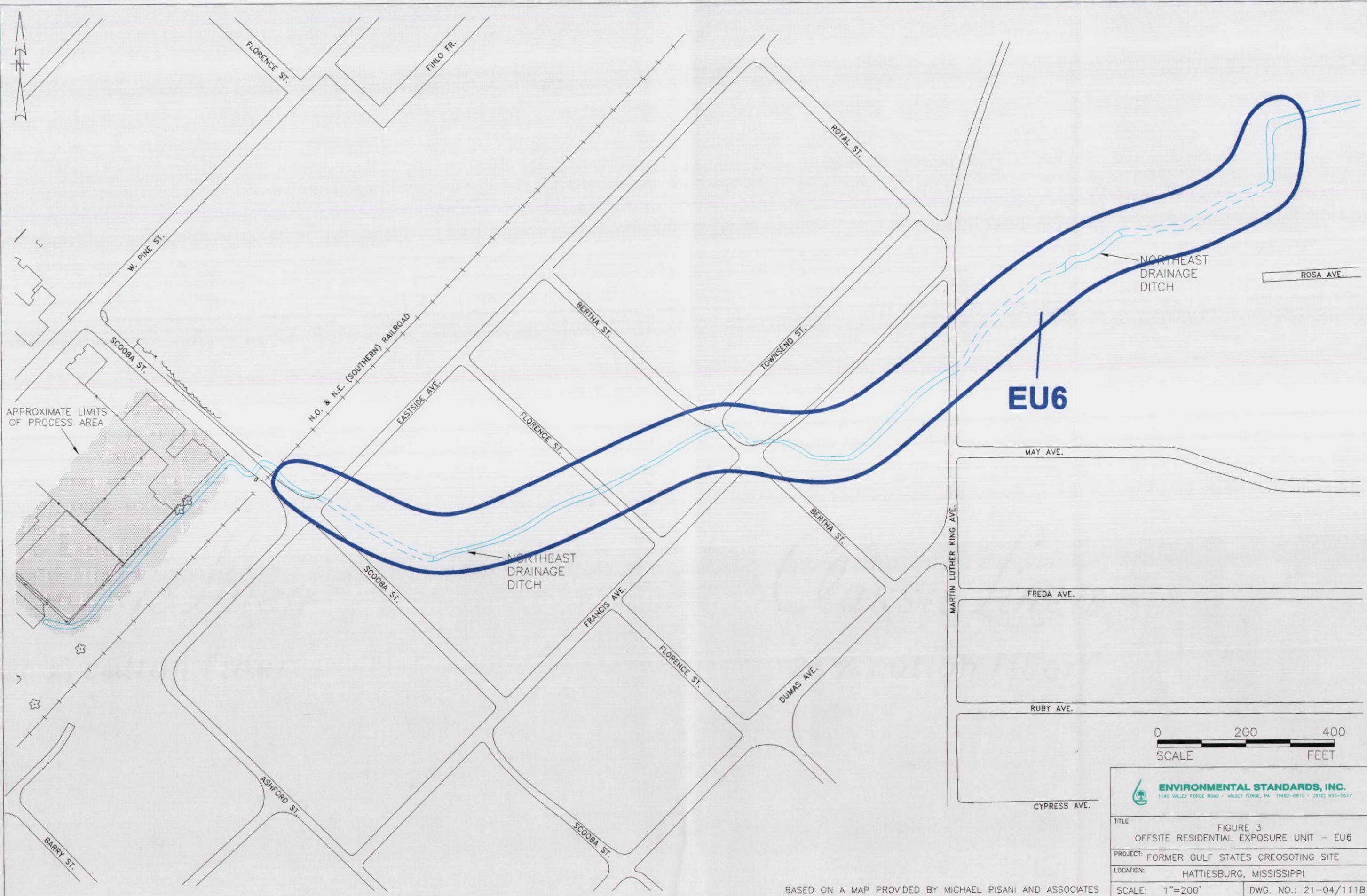
**Figure 1**

**Site Conceptual Model and Selection of Exposure Pathways  
Kerr McGee, Hattiesburg, MS**

| Scenario Timeframe |                 | Medium                              | Exposure Medium | Exposure Point      | Receptor Population | Receptor Age                           | Exposure Route  | On-Site/Off-Site                       | Type of Analysis  | Rationale for Selection or Exclusion of Exposure Pathway |
|--------------------|-----------------|-------------------------------------|-----------------|---------------------|---------------------|--|---|--|---|--|
|                    | Subsurface Soil | Subsurface Soil (0' to water table) | Exposure Unit 1 | Construction Worker | Adult               | Dermal Oral Inhalation                 | On-Site   | None                                   | EU1 includes only surface water and sediment. Soils in this area are included in EU2                                      |  |
|                    |                 |                                     | Exposure Unit 2 | Construction Worker | Adult               | Dermal Oral                            | On-Site   | None                                   | EU1 includes only surface water and sediment. Soils in this area are included in EU2                                      |  |
|                    |                 |                                     | Exposure Unit 3 | Construction Worker | Adult               | Dermal Oral                            | On-Site   | Quantitative Qualitative Quantitative  | Potentially constructable area in the future<br>Only cPAHs were COPCs; addressed qualitatively according to MCEQ guidance |  |
|                    |                 |                                     | Exposure Unit 4 | Construction Worker | Adult               | Dermal Oral                            | On-Site   | None                                   | COPCs eliminated during screening process   |  |
|                    |                 |                                     | Exposure Unit 5 | Construction Worker | Adult               | Dermal Oral                            | On-Site   | Quantitative Quantitative Quantitative | COPCs eliminated during screening process   |  |
| Sediment           | Exposure Unit 1 | Maintenance Worker                  | Adult           | Dermal Oral         | On-Site             | Quantitative Quantitative Quantitative | Infrequent construction activities may occur in the future<br>Only cPAHs were COPCs; addressed qualitatively according to MCEQ guidance |  |   |  |
|                    | Exposure Unit 4 | Maintenance Worker                  | Adult           | Dermal Oral         | On-Site             | Quantitative Quantitative Quantitative | Infrequent construction activities may occur in the future<br>Only cPAHs were COPCs; addressed qualitatively according to MCEQ guidance |  |   |  |
|                    | Exposure Unit 1 | Construction Worker                 | Adult           | Dermal Oral         | On-Site             | Quantitative Quantitative Quantitative | Potentially constructable area in the future<br>Only cPAHs were COPCs; addressed qualitatively according to MCEQ guidance               |  |   |  |
|                    | Exposure Unit 4 | Construction Worker                 | Adult           | Dermal Oral         | On-Site             | Quantitative Quantitative Quantitative | Infrequent construction activities may occur in the future<br>Only cPAHs were COPCs; addressed qualitatively according to MCEQ guidance |  |   |  |
| Surface Water      | Exposure Unit 1 | Maintenance Worker                  | Adult           | Dermal Oral         | On-Site             | Quantitative Quantitative Quantitative | Infrequent maintenance of Gordon's Creek<br>VOCs not present at levels of concern   |  |   |  |
|                    | Exposure Unit 4 | Maintenance Worker                  | Adult           | Dermal Oral         | On-Site             | Quantitative Quantitative Quantitative | Infrequent construction activities may occur in the future<br>VOCs not present at levels of concern                                     |  |   |  |
|                    | Exposure Unit 1 | Construction Worker                 | Adult           | Dermal Oral         | On-Site             | Quantitative Quantitative Quantitative | Infrequent construction activities may occur in the future<br>VOCs not present at levels of concern                                     |  |   |  |
|                    | Exposure Unit 4 | Construction Worker                 | Adult           | Dermal Oral         | On-Site             | Quantitative Quantitative Quantitative | Infrequent construction activities may occur in the future<br>VOCs not present at levels of concern                                     |  |   |  |







**Table I**  
**Statistical Summary and Selection of COPCs in EU1 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit | Detected | Minimum mg/kg | Maximum mg/kg | Mean mg/kg | Logarithmic Mean | Maximum mg/kg | Detected | Maximum mg/kg | Qualifier | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------|----------|---------------|---------------|------------|------------------|---------------|----------|---------------|-----------|-----------------------------------|--------------------------|
| Semi-volatiles         |            |               |                |                 |               |               |                 |          |               |               |            |                  |               |          |               |           |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 7.40E-02        | J        | 2.92E-01      | 5.10E-01      | SD-07      | 3.08E-01         |               |          |               |           |                                   |                          |
| Acenaphthene           | 83-32-9    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.80E-01        | J        | 3.15E-01      | 2.85E-01      | 4.50E-01   | SD-07            | 1.91E-01      |          |               |           |                                   |                          |
| Acenaphthylene         | 208-96-8   | 2             | 1              | 50              | 4.00E-02      | 4.00E-02      | 7.80E-02        | J        | 4.90E-02      | 3.95E-02      | 7.80E-02   | J                | 4.10E-02      |          |               |           |                                   |                          |
| Anthracene             | 120-12-7   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 2.60E-01        | J        | 3.60E-01      | 3.46E-01      | 4.60E-01   | SD-07            | 1.41E-01      |          |               |           |                                   |                          |
| Benz(a)anthracene      | 56-55-3    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.80E-01        | J        | 3.85E-01      | 3.26E-01      | 5.90E-01   | SD-07            | 2.90E-01      |          |               |           |                                   |                          |
| Benz(a)pyrene          | 50-32-8    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.20E-01        | J        | 2.55E-01      | 2.16E-01      | 3.90E-01   | J                | 1.91E-01      |          |               |           |                                   |                          |
| Benz(b)fluoranthene    | 205-99-2   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.70E-01        | J        | 3.75E-01      | 3.14E-01      | 5.80E-01   | SD-07            | 2.90E-01      |          |               |           |                                   |                          |
| Benz(ghi)perylene      | 191-24-2   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 6.50E-02        | J        | 1.23E-01      | 1.08E-01      | 1.80E-01   | J                | 8.13E-02      |          |               |           |                                   |                          |
| Benz(k)fluoranthene    | 207-08-9   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 6.40E-02        | J        | 1.27E-01      | 1.10E-01      | 1.90E-01   | J                | 8.91E-02      |          |               |           |                                   |                          |
| Carbazole              | 86-74-8    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.60E-01        | J        | 3.65E-01      | 3.02E-01      | 5.70E-01   | SD-07            | 2.90E-01      |          |               |           |                                   |                          |
| Chrysene               | 218-01-9   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.80E-01        | J        | 3.55E-01      | 3.09E-01      | 5.30E-01   | SD-07            | 2.47E-01      |          |               |           |                                   |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 2             | 1              | 50              | 4.00E-02      | 4.00E-02      | 6.20E-02        | J        | 4.10E-02      | 3.52E-02      | 6.20E-02   | J                | 2.97E-02      |          |               |           |                                   |                          |
| Dibenzofuran           | 132-64-9   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.50E-01        | J        | 2.80E-01      | 2.48E-01      | 4.10E-01   | SD-07            | 1.84E-01      |          |               |           |                                   |                          |
| Fluoranthene           | 206-44-0   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 6.80E-01        | J        | 1.19E+00      | 1.08E+00      | 1.70E+00   | SD-07            | 7.21E-01      |          |               |           |                                   |                          |
| Fluorene               | 86-73-7    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 2.30E-01        | J        | 4.25E-01      | 3.78E-01      | 6.20E-01   | SD-07            | 2.76E-01      |          |               |           |                                   |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 6.90E-02        | J        | 1.45E-01      | 1.23E-01      | 2.20E-01   | J                | 1.07E-01      |          |               |           |                                   |                          |
| Naphthalene            | 91-20-3    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.80E-01        | J        | 6.40E-01      | 4.45E-01      | 1.10E+00   | SD-07            | 6.51E-01      |          |               |           |                                   |                          |
| Phenanthrene           | 85-01-8    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 7.20E-01        | J        | 1.21E+00      | 1.11E+00      | 1.70E+00   | SD-07            | 6.93E-01      |          |               |           |                                   |                          |
| Pyrone                 | 129-00-0   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 4.80E-01        | J        | 9.40E-01      | 8.20E-01      | 1.40E+00   | SD-07            | 6.51E-01      |          |               |           |                                   |                          |



**Table 1**  
**Statistical Summary and Selection of COPCs in EU1 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Unrestricted<br>Soil TRG<br>mg/kg | Is Maximum<br>Detected ><br>TRG? | Is the 95%<br>UCL ><br>TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|--|----------------------------------|-----------------------------|
| <b>Semivolatiles</b>   |                  |                                 |                                   |  |  |                                  |                             |
| 2-methylnaphthalene    | 1.67E+00         | 1.60E+22                        | Unknown                           | 5.10E-01                                 | 3.13E+03                                 | no                               |                             |
| Acenaphthene           | 1.17E+00         | 3.23E+04                        | Unknown                           | 4.50E-01                                 | 4.69E+03                                 | no                               |                             |
| Acenaphthylene         | 2.32E-01         | 8.34E+09                        | Unknown                           | 7.80E-02                                 | 4.69E+03                                 | no                               |                             |
| Anthracene             | 9.91E-01         | 2.23E+01                        | Unknown                           | 4.60E-01                                 | 2.35E+04                                 | no                               |                             |
| Benz(a)anthracene      | 1.68E+00         | 1.25E+08                        | Unknown                           | 5.90E-01                                 | 8.75E-01                                 | no                               | YES*                        |
| Benz(a)pyrene          | 1.11E+00         | 6.25E+07                        | Unknown                           | 3.90E-01                                 | 8.75E-02                                 | YES                              | YES - COPC                  |
| Benz(b)fluoranthene    | 1.67E+00         | 4.79E+08                        | Unknown                           | 5.80E-01                                 | 8.75E-01                                 | no                               | YES*                        |
| Benz(ghi)perylene      | 4.86E-01         | 2.08E+05                        | Unknown                           | 1.80E-01                                 | 2.35E+03                                 | no                               | YES*                        |
| Benz(k)fluoranthene    | 5.25E-01         | 1.71E+06                        | Unknown                           | 1.90E-01                                 | 8.75E+00                                 | no                               | YES*                        |
| Carbazole              | 1.66E+00         | 2.15E+09                        | Unknown                           | 5.70E-01                                 | 3.19E+01                                 | no                               |                             |
| Chrysene               | 1.46E+00         | 3.73E+06                        | Unknown                           | 5.30E-01                                 | 8.75E+01                                 | no                               | YES*                        |
| Dibenz(a,h)anthracene  | 1.74E-01         | 2.15E+06                        | Unknown                           | 6.20E-02                                 | 8.75E-02                                 | no                               | YES*                        |
| Dibenzofuran           | 1.10E+00         | 3.27E+05                        | Unknown                           | 4.10E-01                                 | 3.13E+02                                 | no                               |                             |
| Fluoranthene           | 4.41E+00         | 1.22E+05                        | Unknown                           | 1.70E+00                                 | 3.13E+03                                 | no                               |                             |
| Fluorene               | 1.66E+00         | 3.35E+05                        | Unknown                           | 6.20E-01                                 | 3.13E+03                                 | no                               | YES*                        |
| Indeno(1,2,3-cd)pyrene | 6.21E-01         | 1.88E+07                        | Unknown                           | 2.20E-01                                 | 8.75E-01                                 | no                               | YES*                        |
| Naphthalene            | 3.54E+00         | 6.10E+19                        | Unknown                           | 1.10E+00                                 | 6.45E+02                                 | no                               |                             |
| Phenanthrene           | 4.30E+00         | 2.92E+04                        | Unknown                           | 1.70E+00                                 | 2.35E+03                                 | no                               |                             |
| Pyrene                 | 3.84E+00         | 7.40E+06                        | Unknown                           | 1.40E+00                                 | 2.35E+03                                 | no                               |                             |

\* Retained as a COPC, as per MDEQ Comments (8/2/2000): constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 2

*Statistical Summary and Selection of COPCs in EU1 Surface Water  
Kerr McGee, Hattiesburg, MS*

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/L | Maximum mg/L | Mean mg/L | Logarithmic Mean mg/L | Maximum mg/L | Detected Qualifier | Detected mg/L | Maximum mg/L | Location of Maximum Concentration | Standard Deviation mg/L |
|------------------------|------------|---------------|----------------|-----------------|--------------|--------------|-----------|-----------------------|--------------|--------------------|---------------|--------------|-----------------------------------|-------------------------|
| <b>Semivolatiles</b>   |            |               |                |                 |              |              |           |                       |              |                    |               |              |                                   |                         |
| Benzo(a)anthracene     | 56-55-3    | 2             | 1              | 50              | 1.00E-03     | 1.00E-03     | J         | 7.50E-04              | 7.07E-04     | 1.00E-03           | J             | SW-08        | 3.54E-04                          |                         |
| Benzo(a)pyrene         | 50-32-8    | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04     | 0.00E+00           | NA            | SW-08        | 0.00E+00                          |                         |
| Benzo(b)fluoranthene   | 205-99-2   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | 0.00E+00  | NA                    | 5.00E-04     | 0.00E+00           | 0.00E+00      | NA           | SW-08                             | 0.00E+00                |
| Benzo(k)fluoranthene   | 207-08-9   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | 0.00E+00  | NA                    | 5.00E-04     | 0.00E+00           | 0.00E+00      | NA           | SW-08                             | 0.00E+00                |
| Chrysene               | 218-01-9   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | 0.00E+00  | NA                    | 5.00E-04     | 0.00E+00           | 0.00E+00      | NA           | SW-08                             | 0.00E+00                |
| Dibenz(a,h)anthracene  | 53-70-3    | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | 0.00E+00  | NA                    | 5.00E-04     | 0.00E+00           | 0.00E+00      | NA           | SW-08                             | 0.00E+00                |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | 0.00E+00  | NA                    | 5.00E-04     | 5.00E-04           | 0.00E+00      | NA           | SW-08                             | 0.00E+00                |
| Fluoranthene           | 206-44-0   | 2             | 1              | 50              | 1.00E-03     | 1.00E-03     | 7.50E-03  | 4.00E-03              | 1.94E-03     | 7.50E-03           | 4.00E-03      | SW-08        | 4.95E-03                          |                         |
| Pyrene                 | 129-00-0   | 2             | 1              | 50              | 1.00E-03     | 1.00E-03     | J         | 7.50E-04              | 7.07E-04     | 1.00E-03           | J             | SW-08        | 3.54E-04                          |                         |

NA - Not applicable; constituent not detected in media.



**Table 2**

**Statistical Summary and Selection of COPCs in EU1 Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | Human Health    |                 |                   |   |  |                                   |
|------------------------|-----------------|-----------------|-------------------|---|--|-----------------------------------|
|                        | 95% UCL<br>mg/L | 95% UCL<br>mg/L | 99%<br>Confidence | Exposure Point<br>Concentration<br>mg/L | Consumption of Water<br>& Organisms AWQC<br>mg/L | Is Maximum<br>Detected ><br>AWQC? |
| <b>Semivolatile</b>    |                 |                 |                   |   |  |                                   |
| Benz(a)anthracene      | 2.33E-03        | 4.37E-01        | Unknown           | 1.00E-03                                | 4.40E-06   | YES - COPC                        |
| Benz(a)pyrene          | 5.00E-04        | 5.00E-04        | Unknown           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Benz(b)fluoranthene    | 5.00E-04        | 5.00E-04        | Unknown           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Benz(k)fluoranthene    | 5.00E-04        | 5.00E-04        | Unknown           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Chrysene               | 5.00E-04        | 5.00E-04        | Unknown           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Dibenz(a,h)anthracene  | 5.00E-04        | 5.00E-04        | Unknown           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Indeno(1,2,3-cd)pyrene | 5.00E-04        | 5.00E-04        | Unknown           | 5.00E-04                                | 4.40E-06   | YES**                             |
| Fluoranthene           | 2.61E-02        | 2.90E-42        | Unknown           | 7.50E-03                                | 3.00E-01   | no                                |
| Pyrene                 | 2.33E-03        | 4.37E-01        | Unknown           | 1.00E-03                                | 9.60E-01   | no                                |

NA - Not Available

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.



**Table 3**  
**Statistical Summary and Selection of COPCs in EU2 Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum Detection Limit mg/kg | Detected mg/kg | Minimum Detected mg/kg | Detected Qualifier mg/kg | Mean mg/kg | Logarithmic Mean | Maximum Detected mg/kg | Detected Qualifier mg/kg | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|-------------------------------|----------------|------------------------|--------------------------|------------|------------------|------------------------|--------------------------|-----------------------------|-----------------------------------|--------------------------|
| <b>Semi-volatiles</b>  |            |               |                |                 |               |                               |                |                        |                          |            |                  |                        |                          |                             |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 14            | 2              | 14.29           | 3.30E-02      | 7.00E-02                      | J              | 3.06E-02               | 2.15E-02                 | 1.60E-01   | J                | SS-10                  | 3.99E-02                 |                             |                                   |                          |
| Acenaphthene           | 83-32-9    | 14            | 1              | 7.14            | 3.30E-02      | 4.90E-02                      | J              | 1.88E-02               | 1.78E-02                 | 4.90E-02   | J                | GEO-13                 | 8.69E-03                 |                             |                                   |                          |
| Acenaphthylene         | 208-96-8   | 14            | 6              | 42.86           | 3.30E-02      | 3.70E-02                      | J              | 1.59E-01               | 4.29E-02                 | 1.30E+00   | J                | GEO-13                 | 3.52E-01                 |                             |                                   |                          |
| Anthracene             | 120-12-7   | 14            | 7              | 50              | 3.30E-02      | 3.70E-02                      | J              | 1.89E-01               | 5.00E-02                 | 1.60E+00   | J                | GEO-13                 | 4.28E-01                 |                             |                                   |                          |
| Benz(a)anthracene      | 56-55-3    | 14            | 12             | 85.71           | 3.30E-02      | 4.10E-02                      | J              | 8.98E-01               | 2.28E-01                 | 6.70E+00   | J                | GEO-13                 | 1.78E+00                 |                             |                                   |                          |
| Benz(a)pyrene          | 50-32-8    | 14            | 11             | 78.57           | 6.70E-02      | 8.40E-02                      | J              | 8.31E-01               | 2.82E-01                 | 5.20E+00   | J                | GEO-13                 | 1.42E+00                 |                             |                                   |                          |
| Benz(b)fluoranthene    | 205-99-2   | 14            | 12             | 85.71           | 6.70E-02      | 6.70E-02                      | J              | 1.10E-01               | 1.84E+00                 | 5.95E-01   | J                | GEO-13                 | 2.62E+00                 |                             |                                   |                          |
| Benzoguaiacol          | 191-24-2   | 14            | 10             | 71.43           | 6.70E-02      | 6.70E-02                      | J              | 1.70E-01               | 5.17E-01                 | 2.20E+00   | J                | GEO-13                 | 6.95E-01                 |                             |                                   |                          |
| Benz(k)fluoranthene    | 207-08-9   | 14            | 9              | 64.29           | 1.30E-01      | 1.30E-01                      | J              | 1.90E-01               | 7.01E-01                 | 2.88E+01   | J                | GEO-13                 | 1.04E+00                 |                             |                                   |                          |
| Carbazole              | 86-74-8    | 14            | 4              | 28.57           | 3.30E-02      | 3.30E-02                      | J              | 4.30E-02               | 6.28E-02                 | 2.94E-02   | J                | GEO-13                 | 1.05E-01                 |                             |                                   |                          |
| Chrysene               | 218-01-9   | 14            | 12             | 85.71           | 3.30E-02      | 3.30E-02                      | J              | 6.20E-02               | 1.19E-01                 | 3.11E-01   | J                | GEO-13                 | 2.16E+00                 |                             |                                   |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 14            | 7              | 50              | 6.70E-02      | 6.70E-02                      | J              | 1.85E-01               | 8.87E-02                 | 9.10E-01   | J                | GEO-13                 | 2.66E-01                 |                             |                                   |                          |
| Dibenzofuran           | 132-64-9   | 14            | 2              | 14.29           | 3.30E-02      | 3.30E-02                      | J              | 2.63E-02               | 2.08E-02                 | 9.80E-02   | J                | SS-10                  | 2.54E-02                 |                             |                                   |                          |
| Di-n-butylphthalate    | 84-74-2    | 14            | 9              | 64.29           | 3.30E-02      | 7.20E-02                      | J              | 3.60E-02               | 4.30E-02                 | 3.68E-02   | J                | SS-10                  | 2.50E-02                 |                             |                                   |                          |
| Fluoranthene           | 206-44-0   | 14            | 12             | 85.71           | 3.30E-02      | 6.60E-02                      | J              | 1.40E+00               | 3.00E-01                 | 1.20E+01   | J                | GEO-13                 | 3.16E+00                 |                             |                                   |                          |
| Fluorene               | 86-73-7    | 14            | 2              | 14.29           | 3.30E-02      | 4.50E-02                      | J              | 4.38E-02               | 2.21E-02                 | 3.70E-01   | J                | GEO-13                 | 9.42E-02                 |                             |                                   |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 14            | 10             | 71.43           | 6.70E-02      | 9.60E-02                      | J              | 6.59E-01               | 2.37E-01                 | 3.70E+00   | J                | GEO-13                 | 1.03E+00                 |                             |                                   |                          |
| Naphthalene            | 91-20-3    | 14            | 2              | 14.29           | 3.30E-02      | 8.80E-02                      | J              | 3.26E-02               | 2.20E-02                 | 1.70E-01   | J                | SS-10                  | 4.39E-02                 |                             |                                   |                          |
| Phenanthrene           | 85-01-8    | 14            | 8              | 57.14           | 3.30E-02      | 3.70E-02                      | J              | 1.28E-01               | 5.30E-02                 | 7.40E-01   | J                | GEO-13                 | 2.08E-01                 |                             |                                   |                          |
| Pyrene                 | 129-00-0   | 14            | 12             | 85.71           | 6.70E-02      | 9.80E-02                      | J              | 1.70E+00               | 4.60E-01                 | 1.40E+01   | J                | GEO-13                 | 3.66E+00                 |                             |                                   |                          |

**Table 3**  
**Statistical Summary and Selection of COPCs in EU2 Soil (0'-1' Egs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>Unrestricted<br>Soil TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? | Is the 95%<br>UCL > TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|---|---|--------------------------|
| <b>Semi-volatiles</b>  |                  |                                 |                                   |  |   |   |                          |
| 2-methylnaphthalene    | 4.95E-02         | 4.29E-02                        | Unknown                           | 4.29E-02                                 | 3.13E+03                                    | no                                      |                          |
| Acenaphthene           | 2.29E-02         | 2.17E-02                        | Unknown                           | 2.17E-02                                 | 4.69E+03                                    | no                                      |                          |
| Acenaphthylene         | 3.26E-01         | 4.99E-01                        | Unknown                           | 4.99E-01                                 | 4.69E+03                                    | no                                      |                          |
| Anthracene             | 3.91E-01         | 6.29E-01                        | Unknown                           | 6.29E-01                                 | 2.35E+04                                    | no                                      |                          |
| Benzo(a)anthracene     | 1.74E+00         | 9.91E+00                        | Lognormal                         | 6.70E+00                                 | 8.75E-01                                    | YES                                     | YES - COPC               |
| Benzo(a)pyrene         | 1.50E+00         | 5.08E+00                        | Lognormal                         | 5.08E+00                                 | 8.75E-02                                    | YES                                     | YES - COPC               |
| Benzo(b)fluoranthene   | 3.08E+00         | 2.53E+01                        | Lognormal                         | 9.20E+00                                 | 8.75E-01                                    | YES                                     | YES - COPC               |
| Benzo(g,h,i)perylene   | 8.46E-01         | 2.74E+00                        | Lognormal                         | 2.30E+00                                 | 2.35E+03                                    | no                                      |                          |
| Benzo(k)fluoranthene   | 1.19E+00         | 2.93E+00                        | Lognormal                         | 2.93E+00                                 | 8.75E+00                                    | no                                      |                          |
| Carbazole              | 1.12E-01         | 1.24E-01                        | Unknown                           | 1.24E-01                                 | 3.19E+01                                    | no                                      |                          |
| Chrysene               | 2.22E+00         | 1.68E+01                        | Lognormal                         | 8.00E+00                                 | 8.75E+01                                    | no                                      |                          |
| Dibenz(a,h)anthracene  | 3.11E-01         | 4.93E-01                        | Unknown                           | 4.93E-01                                 | 8.75E-02                                    | YES                                     | YES - COPC               |
| Dibenzofuran           | 3.83E-02         | 3.57E-02                        | Unknown                           | 3.57E-02                                 | 3.13E+02                                    | no                                      |                          |
| Di-n-butylphthalate    | 5.48E-02         | 6.30E-02                        | Normal/Lognormal                  | 6.30E-02                                 | 2.28E+03                                    | no                                      |                          |
| Fluoranthene           | 2.89E+00         | 1.66E+01                        | Lognormal                         | 1.20E+01                                 | 3.13E+03                                    | no                                      |                          |
| Fluorene               | 8.84E-02         | 5.84E-02                        | Unknown                           | 5.84E-02                                 | 3.13E+03                                    | no                                      |                          |
| Indeno(1,2,3-cd)pyrene | 1.15E+00         | 4.29E+00                        | Lognormal                         | 3.70E+00                                 | 8.75E-01                                    | YES                                     | YES - COPC               |
| Naphthalene            | 5.34E-02         | 4.71E-02                        | Unknown                           | 4.71E-02                                 | 6.45E+02                                    | no                                      |                          |
| Phenanthrene           | 2.26E-01         | 3.96E-01                        | Lognormal                         | 3.96E-01                                 | 2.35E+03                                    | no                                      |                          |
| Pyrene                 | 3.43E+00         | 1.25E+01                        | Lognormal                         | 1.25E+01                                 | 2.35E+03                                    | no                                      |                          |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 4

*Statistical Summary and Selection of COPCs in EU2 Soil (0-6' bgs)*  
*Kerr McGee, Hattiesburg, MS*

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit | Minimum mg/kg | Maximum mg/kg | Detected Qualifier | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected Qualifier | Mean mg/kg | Maximum mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------|---------------|---------------|--------------------|------------|------------------------|---------------|--------------------|------------|---------------|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>       |            |               |                |                 |               |               |                 |               |               |                    |            |                        |               |                    |            |               |                                   |                          |
| 2-methylphthalene          | 91-57-6    | 20            | 2              | 10              | 3.30E-02      | 7.00E-02      | J               | 2.71E-02      | 2.07E-02      | 1.60E-01           | J          | SS-10                  | 3.34E-02      |                    |            |               |                                   |                          |
| Acenaphthene               | 83-32-9    | 21            | 1              | 4.76            | 3.30E-02      | 3.00E-01      | 4.90E-02        | J             | 2.51E-02      | 2.01E-02           | 4.90E-02   | J                      | GEO-13        | 2.95E-02           |            |               |                                   |                          |
| Acenaphthylene             | 208-96-8   | 21            | 6              | 28.57           | 3.30E-02      | 3.00E-01      | 3.70E-02        | J             | 1.19E-01      | 3.60E-02           | 1.30E+00   | J                      | GEO-13        | 2.91E-01           |            |               |                                   |                          |
| Anthracene                 | 120-12-7   | 21            | 8              | 38.1            | 3.30E-02      | 3.90E-02      | 4.10E-02        | J             | 1.37E-01      | 3.94E-02           | 1.60E+00   | J                      | GEO-13        | 3.54E-01           |            |               |                                   |                          |
| Benz(a)anthracene          | 56-55-3    | 21            | 14             | 66.67           | 3.30E-02      | 3.80E-02      | 4.10E-02        | J             | 6.10E-01      | 1.12E-01           | 6.70E+00   | J                      | GEO-13        | 1.49E+00           |            |               |                                   |                          |
| Benz(a)pyrene              | 50-32-8    | 21            | 12             | 57.14           | 3.70E-02      | 6.70E-02      | 8.40E-02        | J             | 5.65E-01      | 1.25E-01           | 5.20E+00   | J                      | GEO-13        | 1.21E+00           |            |               |                                   |                          |
| Benz(b)fluoranthene        | 205-99-2   | 21            | 16             | 76.19           | 3.70E-02      | 6.70E-02      | 9.50E-02        | J             | 1.29E+00      | 3.16E-01           | 9.20E+00   | J                      | GEO-13        | 2.26E+00           |            |               |                                   |                          |
| Benz(g,h)perylene          | 191-24-2   | 21            | 11             | 52.38           | 3.70E-02      | 6.70E-02      | 8.50E-02        | J             | 3.54E-01      | 1.04E-01           | 2.30E+00   | J                      | GEO-13        | 6.08E-01           |            |               |                                   |                          |
| Benz(k)fluoranthene        | 207-08-9   | 21            | 13             | 61.9            | 3.70E-02      | 1.30E-01      | 5.10E-02        | J             | 5.21E-01      | 1.84E-01           | 3.60E+00   | J                      | GEO-13        | 8.79E-01           |            |               |                                   |                          |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 20            | 1              | 5               | 6.70E-02      | 7.80E-02      | 3.70E-01        | J             | 5.15E-02      | 3.91E-02           | 3.70E-01   | J                      | GEO-13        | 7.50E-02           |            |               |                                   |                          |
| Carbazole                  | 86-74-8    | 20            | 4              | 20              | 3.30E-02      | 3.90E-02      | 4.30E-02        | J             | 4.96E-02      | 2.57E-02           | 3.50E-01   | J                      | GEO-13        | 8.92E-02           |            |               |                                   |                          |
| Chrysene                   | 218-01-9   | 21            | 13             | 61.9            | 3.30E-02      | 7.40E-02      | 5.10E-02        | J             | 8.03E-01      | 1.32E-01           | 8.00E+00   | J                      | GEO-13        | 1.83E+00           |            |               |                                   |                          |
| Dibenz(a,h)anthracene      | 53-70-3    | 21            | 8              | 38.1            | 3.70E-02      | 6.70E-02      | 1.88E-02        | J             | 1.29E-01      | 5.30E-02           | 9.10E-01   | J                      | GEO-13        | 2.29E-01           |            |               |                                   |                          |
| Dibenzofuran               | 132-64-9   | 20            | 2              | 10              | 3.30E-02      | 3.90E-02      | 7.20E-02        | J             | 2.41E-02      | 2.02E-02           | 9.80E-02   | J                      | SS-10         | 2.13E-02           |            |               |                                   |                          |
| Di-n-butylphthalate        | 84-74-2    | 20            | 9              | 45              | 3.30E-02      | 7.80E-02      | 3.60E-02        | J             | 4.15E-02      | 3.71E-02           | 1.10E-01   | J                      | SS-10         | 2.08E-02           |            |               |                                   |                          |
| Fluoranthene               | 206-44-0   | 21            | 14             | 66.67           | 3.30E-02      | 3.80E-02      | 5.00E-02        | J             | 9.54E-01      | 1.42E-01           | 1.20E-01   | J                      | GEO-13        | 2.63E+00           |            |               |                                   |                          |
| Fluorene                   | 86-73-7    | 21            | 4              | 19.05           | 3.30E-02      | 3.80E-02      | 2.90E-02        | J             | 5.08E-02      | 2.45E-02           | 3.70E-01   | J                      | GEO-13        | 9.99E-02           |            |               |                                   |                          |
| Indeno(1,2,3-cd)pyrene     | 193-39-5   | 21            | 11             | 52.38           | 3.70E-02      | 6.70E-02      | 9.60E-02        | J             | 4.50E-01      | 1.11E-01           | 3.70E+00   | J                      | GEO-13        | 8.86E-01           |            |               |                                   |                          |
| Naphthalene                | 91-20-3    | 21            | 2              | 9.52            | 3.30E-02      | 3.00E-01      | 8.80E-02        | J             | 3.43E-02      | 2.31E-02           | 1.70E-01   | J                      | SS-10         | 4.47E-02           |            |               |                                   |                          |
| Phenanthrene               | 85-01-8    | 21            | 9              | 42.86           | 3.30E-02      | 3.90E-02      | 3.70E-02        | J             | 1.01E-01      | 4.22E-02           | 7.40E-01   | J                      | GEO-13        | 1.77E-01           |            |               |                                   |                          |
| Phenol                     | 108-95-2   | 20            | 2              | 10              | 3.30E-02      | 7.80E-02      | 1.10E-01        | J             | 3.51E-02      | 2.52E-02           | 1.90E-01   | J                      | GEO-13        | 4.24E-02           |            |               |                                   |                          |
| Pyrene                     | 129-00-0   | 21            | 14             | 66.67           | 3.70E-02      | 6.70E-02      | 6.80E-02        | J             | 1.16E+00      | 1.92E-01           | 1.40E+01   | J                      | GEO-13        | 3.05E+00           |            |               |                                   |                          |



Table 4

Statistical Summary and Selection of COPCs in EU2 Soil (0'-6' bg3)  
Kerr McGee, Hattiesburg, MS

| Constituent                | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG?<br>UCL > TRG? |            |  |
|----------------------------|------------------|---------------------------------|-----------------------------------|--|------------------------|---|------------|--|
|                            |                  |                                 |                                   |  |                        | Is the<br>95%<br>UCL > TRG?                           |            |  |
| Semi-volatiles             |                  |                                 |                                   |  |                        |   |            |  |
| 2-methylnaphthalene        | 4.00E-02         | 3.22E-02                        | Unknown                           | 3.22E-02                                 | 8.18E+04               | no  |            |  |
| Acenaphthene               | 3.62E-02         | 2.90E-02                        | Unknown                           | 2.90E-02                                 | 1.23E+05               | no  |            |  |
| Acenaphthylene             | 2.28E-01         | 1.83E-01                        | Unknown                           | 1.83E-01                                 | 1.23E+05               | no  |            |  |
| Anthracene                 | 2.70E-01         | 2.09E-01                        | Unknown                           | 2.09E-01                                 | 6.13E+05               | no  |            |  |
| Benzo(a)anthracene         | 1.17E+00         | 2.80E+00                        | Lognormal                         | 2.80E+00                                 | 7.84E+00               | no  | YES*       |  |
| Benzo(a)pyrene             | 1.02E+00         | 2.64E+00                        | Lognormal                         | 2.64E+00                                 | 7.84E-01               | YES   | YES - COPC |  |
| Benzo(b)fluoranthene       | 2.14E+00         | 1.09E+01                        | Lognormal                         | 9.20E+00                                 | 7.84E+00               | YES   | YES - COPC |  |
| Benzo(g,h)perylene         | 5.83E-01         | 1.41E+00                        | Lognormal                         | 1.41E+00                                 | 6.13E+04               | no  |            |  |
| Benzo(k)fluoranthene       | 8.52E-01         | 1.84E+00                        | Lognormal                         | 1.84E+00                                 | 7.84E+01               | no  | YES*       |  |
| Bis(2-ethylhexyl)phthalate | 8.05E-02         | 5.77E-02                        | Unknown                           | 5.77E-02                                 | 4.09E+02               | no  |            |  |
| Carbazole                  | 8.41E-02         | 6.51E-02                        | Unknown                           | 6.51E-02                                 | 2.86E+02               | no  |            |  |
| Chrysene                   | 1.49E+00         | 5.33E+00                        | Lognormal                         | 5.33E+00                                 | 7.84E+02               | no  | YES*       |  |
| Dibenz(a,h)anthracene      | 2.16E-01         | 2.39E-01                        | Unknown                           | 2.39E-01                                 | 7.84E+01               | YES   | YES**      |  |
| Dibenzofuran               | 3.23E-02         | 2.86E-02                        | Unknown                           | 2.86E-02                                 | 8.18E+03               | no  |            |  |
| Di-n-butylphthalate        | 4.95E-02         | 5.24E-02                        | Lognormal                         | 5.24E-02                                 | 2.28E+03               | no  |            |  |
| Fluoranthene               | 1.94E+00         | 5.34E+00                        | Lognormal                         | 5.34E+00                                 | 8.17E+04               | no  | YES*       |  |
| Fluorene                   | 8.84E-02         | 6.16E-02                        | Unknown                           | 6.16E-02                                 | 8.17E+04               | no  |            |  |
| Indeno(1,2,3-cd)pyrene     | 7.83E-01         | 1.97E+00                        | Lognormal                         | 1.97E+00                                 | 7.84E+00               | no  |            |  |
| Naphthalene                | 5.11E-02         | 4.37E-02                        | Unknown                           | 4.37E-02                                 | 8.24E+02               | no  |            |  |
| Phenanthrene               | 1.67E-01         | 1.88E-01                        | Unknown                           | 1.88E-01                                 | 6.13E+04               | no  |            |  |
| Phenol                     | 5.15E-02         | 4.60E-02                        | Unknown                           | 4.60E-02                                 | 1.23E+05               | no  |            |  |
| Pyrene                     | 2.31E+00         | 7.47E+00                        | Lognormal                         | 7.47E+00                                 | 6.13E+04               | no  |            |  |

**Table 5**  
**Statistical Summary and Selection of COPCs in EU2 Soil (0-10' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected Qualifier | Detected Qualifier | Detected Qualifier | Maximum mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|------------|------------------------|---------------|--------------------|--------------------|--------------------|---------------|-----------------------------------|--------------------------|
| <b>Semi-volatiles</b>      |            |               |                |                 |               |               |            |                        |               |                    |                    |                    |               |                                   |                          |
| 2-methylnaphthalene        | 91-57-6    | 20            | 2              | 10              | 3.30E-02      | 3.90E-02      | 7.00E-02   | 2.71E-02               | 2.07E-02      | J                  | 1.60E-01           | J                  | SS-10         | 3.34E-02                          |                          |
| Acenaphthene               | 83-32-9    | 21            | 1              | 4.76            | 3.30E-02      | 3.00E-01      | 4.90E-02   | 2.51E-02               | 2.01E-02      | J                  | 4.90E-02           | J                  | GEO-13        | 2.95E-02                          |                          |
| Acenaphthylene             | 208-96-8   | 21            | 6              | 28.57           | 3.30E-02      | 3.00E-01      | 3.70E-02   | J                      | 1.19E-01      | 3.60E-02           | 1.30E+00           | J                  | GEO-13        | 2.91E-01                          |                          |
| Anthracene                 | 120-12-7   | 21            | 8              | 38.1            | 3.30E-02      | 3.90E-02      | 4.10E-02   | J                      | 1.37E-01      | 3.94E-02           | 1.60E+00           | J                  | GEO-13        | 3.54E-01                          |                          |
| Benz(a)anthracene          | 56-55-3    | 21            | 14             | 66.67           | 3.30E-02      | 3.80E-02      | 4.10E-02   | J                      | 6.10E-01      | 1.12E-01           | 6.70E+00           | J                  | GEO-13        | 1.49E+00                          |                          |
| Benz(a)pyrene              | 50-32-8    | 21            | 12             | 57.14           | 3.70E-02      | 6.70E-02      | 8.40E-02   | J                      | 5.65E-01      | 1.25E-01           | 5.20E+00           | J                  | GEO-13        | 1.21E+00                          |                          |
| Benz(b)fluoranthene        | 205-99-2   | 21            | 16             | 76.19           | 3.70E-02      | 6.70E-02      | 9.50E-02   | J                      | 1.29E+00      | 3.16E-01           | 9.20E+00           | J                  | GEO-13        | 2.26E+00                          |                          |
| Benz(ghi)perylene          | 191-24-2   | 21            | 11             | 52.38           | 3.70E-02      | 6.70E-02      | 8.50E-02   | J                      | 3.54E-01      | 1.04E-01           | 2.30E+00           | J                  | GEO-13        | 6.08E-01                          |                          |
| Benz(k)fluoranthene        | 207-08-9   | 21            | 13             | 61.9            | 3.70E-02      | 1.30E-01      | 5.10E-02   | J                      | 5.21E-01      | 1.84E-01           | 3.60E+00           | J                  | GEO-13        | 8.79E-01                          |                          |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 20            | 1              | 5               | 6.70E-02      | 7.80E-02      | 3.70E-01   | J                      | 5.15E-02      | 3.91E-02           | 3.70E-01           | J                  | GEO-13        | 7.50E-02                          |                          |
| Carbazole                  | 86-74-8    | 20            | 4              | 20              | 3.30E-02      | 3.90E-02      | 4.30E-02   | J                      | 4.96E-02      | 2.57E-02           | 3.50E-01           | J                  | GEO-13        | 8.92E-02                          |                          |
| Chrysene                   | 218-01-9   | 21            | 13             | 61.9            | 3.30E-02      | 7.40E-02      | 5.10E-02   | J                      | 8.03E-01      | 1.32E-01           | 8.00E+00           | J                  | GEO-13        | 1.83E+00                          |                          |
| Dibenz(a,h)anthracene      | 53-70-3    | 21            | 8              | 38.1            | 3.70E-02      | 6.70E-02      | 1.88E-02   | J                      | 1.29E-01      | 5.30E-02           | 9.10E-01           | J                  | GEO-13        | 2.29E-01                          |                          |
| Dibenzofuran               | 132-64-9   | 20            | 2              | 10              | 3.30E-02      | 3.90E-02      | 7.20E-02   | J                      | 2.41E-02      | 2.02E-02           | 9.80E-02           | J                  | SS-10         | 2.13E-02                          |                          |
| Di-n-butylphthalate        | 84-74-2    | 20            | 9              | 45              | 3.30E-02      | 7.80E-02      | 3.60E-02   | J                      | 4.15E-02      | 3.71E-02           | 1.10E-01           | J                  | SS-10         | 2.08E-02                          |                          |
| Fluoranthene               | 206-44-0   | 21            | 14             | 66.67           | 3.30E-02      | 3.80E-02      | 5.00E-02   | J                      | 9.54E-01      | 1.42E-01           | 1.20E+01           | J                  | GEO-13        | 2.63E+00                          |                          |
| Fluorene                   | 86-73-7    | 21            | 4              | 19.05           | 3.30E-02      | 3.80E-02      | 2.90E-02   | J                      | 5.08E-02      | 2.45E-02           | 3.70E-01           | J                  | GEO-13        | 9.99E-02                          |                          |
| Indeno(1,2,3-cd)pyrene     | 193-39-5   | 21            | 11             | 52.38           | 3.70E-02      | 6.70E-02      | 9.60E-02   | J                      | 4.50E-01      | 1.11E-01           | 3.70E+00           | J                  | GEO-13        | 8.86E-01                          |                          |
| Naphthalene                | 91-20-3    | 21            | 2              | 9.52            | 3.30E-02      | 3.00E-01      | 8.80E-02   | J                      | 3.43E-02      | 2.31E-02           | 1.70E-01           | J                  | SS-10         | 4.47E-02                          |                          |
| Phenanthrene               | 85-01-8    | 21            | 9              | 42.86           | 3.30E-02      | 3.90E-02      | 3.70E-02   | J                      | 1.01E-01      | 4.22E-02           | 7.40E-01           | J                  | GEO-13        | 1.77E-01                          |                          |
| Phenol                     | 108-95-2   | 20            | 2              | 10              | 3.30E-02      | 7.80E-02      | 1.10E-01   | J                      | 3.51E-02      | 2.52E-02           | 1.90E-01           | J                  | GEO-13        | 4.24E-02                          |                          |
| Pyrene                     | 129-00-0   | 21            | 14             | 66.67           | 3.70E-02      | 6.70E-02      | 6.80E-02   | J                      | 1.16E+00      | 1.92E-01           | 1.40E+01           | J                  | GEO-13        | 3.05E+00                          |                          |

**Table 5**  
**Statistical Summary and Selection of COPCs in EU2 Soil (0-10' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? |                          |
|----------------------------|------------------|---------------------------------|-----------------------------------|--|------------------------|---|--------------------------|
|                            |                  |                                 |                                   |  |                        | UCL > TRG?                              | Is the 95%<br>UCL > TRG? |
| Semi-volatiles             |                  |                                 |                                   |  |                        |   |                          |
| 2-methylnaphthalene        | 4.00E-02         | 3.22E-02                        | Unknown                           | 3.22E-02                                 | 8.18E+04               | no                                      |                          |
| Acenaphthene               | 3.62E-02         | 2.90E-02                        | Unknown                           | 2.90E-02                                 | 1.23E+05               | no                                      |                          |
| Acenaphthylene             | 2.28E-01         | 1.83E-01                        | Unknown                           | 1.83E-01                                 | 1.23E+05               | no                                      |                          |
| Anthracene                 | 2.70E-01         | 2.09E-01                        | Unknown                           | 2.09E-01                                 | 6.13E+05               | no                                      |                          |
| Benz(a)anthracene          | 1.17E+00         | 2.80E+00                        | Lognormal                         | 2.80E+00                                 | 7.84E+00               | no                                      | YES*                     |
| Benz(a)pyrene              | 1.02E+00         | 2.64E+00                        | Lognormal                         | 2.64E+00                                 | 7.84E-01               | YES                                     | YES - COPC               |
| Benz(b)fluoranthene        | 2.14E+00         | 1.09E+01                        | Lognormal                         | 9.20E+00                                 | 7.84E+00               | YES                                     | YES - COPC               |
| Benz(g,h)perylene          | 5.83E-01         | 1.41E+00                        | Lognormal                         | 1.41E+00                                 | 6.13E+04               | no                                      |                          |
| Benz(k)fluoranthene        | 8.52E-01         | 1.84E+00                        | Lognormal                         | 1.84E+00                                 | 7.84E+01               | no                                      |                          |
| Bis(2-ethylhexyl)phthalate | 8.05E-02         | 5.77E-02                        | Unknown                           | 5.77E-02                                 | 4.09E+02               | no                                      |                          |
| Carbazole                  | 8.41E-02         | 6.51E-02                        | Unknown                           | 6.51E-02                                 | 2.86E+02               | no                                      |                          |
| Chrysene                   | 1.49E+00         | 5.33E+00                        | Lognormal                         | 5.33E+00                                 | 7.84E+02               | no                                      | YES*                     |
| Dibenz(a,h)anthracene      | 2.16E-01         | 2.39E-01                        | Unknown                           | 2.39E-01                                 | 7.84E-01               | YES                                     | YES**                    |
| Dibenzofuran               | 3.23E-02         | 2.86E-02                        | Unknown                           | 2.86E-02                                 | 8.18E+03               | no                                      |                          |
| Di-n-butylphthalate        | 4.95E-02         | 5.24E-02                        | Lognormal                         | 5.24E-02                                 | 2.28E+03               | no                                      |                          |
| Fluoranthene               | 1.94E+00         | 5.34E+00                        | Lognormal                         | 5.34E+00                                 | 8.17E+04               | no                                      |                          |
| Fluorene                   | 8.84E-02         | 6.16E-02                        | Unknown                           | 6.16E-02                                 | 8.17E+04               | no                                      |                          |
| Indeno(1,2,3-cd)pyrene     | 7.83E-01         | 1.97E+00                        | Lognormal                         | 1.97E+00                                 | 7.84E+00               | no                                      | YES*                     |
| Naphthalene                | 5.11E-02         | 4.37E-02                        | Unknown                           | 4.37E-02                                 | 8.24E+02               | no                                      |                          |
| Phenanthrene               | 1.67E-01         | 1.88E-01                        | Unknown                           | 1.88E-01                                 | 6.13E+04               | no                                      |                          |
| Phenol                     | 5.15E-02         | 4.60E-02                        | Unknown                           | 4.60E-02                                 | 1.23E+05               | no                                      |                          |
| Pyrene                     | 2.31E+00         | 7.47E+00                        | Lognormal                         | 7.47E+00                                 | 6.13E+04               | no                                      |                          |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

\*\*Logarithmic 95% UCL is less than benchmark but retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

**Table 6**  
*Statistical Summary and Selection of COPCs in EU3 Soil (0-1' bgs)*  
*Kerr McGee, Hattiesburg, MS*

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected Qualifier | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected Qualifier | Maximum mg/kg | Detected Qualifier | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|--------------------|------------|------------------------|---------------|--------------------|---------------|--------------------|-----------------------------------|--------------------------|
| Semivolatiles          |            |               |                |                 |               |               |                       |                    |            |                        |               |                    |               |                    |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 3             | 1              | 33.33           | 3.30E-02      | 3.30E-02      | 2.30E-01              | J                  | 8.77E-02   | 3.97E-02               | 2.30E-01      | J                  | SS-16         | 1.23E-01           |                                   |                          |
| Acenaphthylene         | 208-96-8   | 3             | 2              | 66.67           | 3.30E-02      | 3.30E-02      | 1.20E-01              | J                  | 1.02E-01   | 6.96E-02               | 1.70E-01      | J                  | SS-16         | 7.83E-02           |                                   |                          |
| Anthracene             | 120-12-7   | 3             | 2              | 66.67           | 3.30E-02      | 3.30E-02      | 1.20E-01              | J                  | 1.02E-01   | 6.96E-02               | 1.70E-01      | J                  | SS-16         | 7.83E-02           |                                   |                          |
| Benz[a]anthracene      | 56-55-3    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 5.60E-02              | J                  | 3.62E-01   | 2.46E-01               | 5.40E-01      | J                  | SS-15         | 2.66E-01           |                                   |                          |
| Benz[a]pyrene          | 50-32-8    | 3             | 2              | 66.67           | 6.70E-02      | 6.70E-02      | 5.60E-01              | J                  | 4.35E-01   | 2.37E-01               | 7.10E-01      | J                  | SS-16         | 3.55E-01           |                                   |                          |
| Benz[b]fluoranthene    | 205-99-2   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 1.90E-01              | J                  | 9.30E-01   | 6.83E-01               | 1.40E+00      | J                  | SS-16         | 6.49E-01           |                                   |                          |
| Benz[ghi]perylene      | 191-24-2   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 8.00E-02              | J                  | 6.53E-01   | 4.03E-01               | 1.20E+00      | J                  | SS-16         | 5.60E-01           |                                   |                          |
| Benzok(k)fluoranthene  | 207-08-9   | 3             | 2              | 66.67           | 1.30E-01      | 1.30E-01      | 4.70E-01              | J                  | 3.42E-01   | 2.46E-01               | 4.90E-01      | J                  | SS-16         | 2.40E-01           |                                   |                          |
| Carbazole              | 86-74-8    | 3             | 2              | 66.67           | 3.30E-02      | 3.30E-02      | 4.60E-02              | J                  | 5.75E-02   | 4.37E-02               | 1.10E-01      | J                  | SS-15         | 4.78E-02           |                                   |                          |
| Chrysene               | 218-01-9   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 1.10E-01              | J                  | 5.93E-01   | 4.25E-01               | 8.70E-01      | J                  | SS-16         | 4.20E-01           |                                   |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 3             | 2              | 66.67           | 6.70E-02      | 6.70E-02      | 1.40E-01              | J                  | 1.11E-01   | 9.09E-02               | 1.60E-01      | J                  | SS-16         | 6.80E-02           |                                   |                          |
| Dibenzofuran           | 132-64-9   | 3             | 2              | 66.67           | 3.30E-02      | 3.30E-02      | 3.60E-02              | J                  | 4.85E-02   | 3.81E-02               | 9.30E-02      | J                  | SS-16         | 3.98E-02           |                                   |                          |
| Di-n-butyl phthalate   | 84-74-2    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 4.00E-02              | J                  | 8.30E-02   | 7.58E-02               | 1.10E-01      | J                  | SS-16         | 3.76E-02           |                                   |                          |
| Fluoranthene           | 206-44-0   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 1.20E-01              | J                  | 5.27E-01   | 3.99E-01               | 7.80E-01      | J                  | SS-16         | 3.56E-01           |                                   |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 8.60E-02              | J                  | 3.85E-01   | 2.89E-01               | 6.00E-01      | J                  | SS-16         | 2.67E-01           |                                   |                          |
| Naphthalene            | 91-20-3    | 3             | 1              | 33.33           | 3.30E-02      | 3.30E-02      | 1.60E-01              | J                  | 6.43E-02   | 3.52E-02               | 1.60E-01      | J                  | SS-16         | 8.28E-02           |                                   |                          |
| Phenanthrene           | 85-01-8    | 3             | 2              | 66.67           | 3.30E-02      | 3.30E-02      | 1.30E-01              | J                  | 1.32E-01   | 8.12E-02               | 2.50E-01      | J                  | SS-16         | 1.17E-01           |                                   |                          |
| Pyrene                 | 129-00-0   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 1.20E-01              | J                  | 6.90E-01   | 4.85E-01               | 1.00E+00      | J                  | SS-17         | 4.94E-01           |                                   |                          |

**Table 6**  
**Statistical Summary and Selection of COPCs in EU3 Soil (0'-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>Unrestricted<br>Soil TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? | Is the<br>Logarithmic<br>95% UCL ><br>TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|---|---|--|
| Semivolatiles          |                  |                                 |                                   |  |   |   |  |
| 2-methylnaphthalene    | 2.95E-01         | 2.43E+08                        | Unknown                           | 2.30E-01                                 | 3.13E+03                                    | no                                      |  |
| Acenaphthylene         | 2.34E-01         | 3.45E+05                        | Normal/Lognormal                  | 1.70E-01                                 | 4.69E+03                                    | no                                      |  |
| Anthracene             | 2.34E-01         | 3.45E+05                        | Normal/Lognormal                  | 1.70E-01                                 | 2.35E+04                                    | no                                      |  |
| Benzo(a)anthracene     | 8.11E-01         | 2.15E+06                        | Normal/Lognormal                  | 5.40E-01                                 | 8.75E-01                                    | no                                      | YES*                                       |
| Benzo(a)pyrene         | 1.93E+00         | 3.82E+11                        | Normal/Lognormal                  | 7.10E-01                                 | 8.75E-02                                    | YES                                     | YES - COPC                                 |
| Benzo(b)fluoranthene   | 2.02E+00         | 1.13E+05                        | Normal/Lognormal                  | 1.40E+00                                 | 8.75E-01                                    | YES                                     | YES - COPC                                 |
| Benzo(ghi)perylene     | 1.60E+00         | 1.70E+08                        | Normal/Lognormal                  | 1.20E+00                                 | 2.35E+03                                    | no                                      |  |
| Benzo(k)fluoranthene   | 7.46E-01         | 1.06E+05                        | Normal/Lognormal                  | 4.90E-01                                 | 8.75E+00                                    | no                                      | YES*                                       |
| Carbazole              | 1.38E-01         | 2.81E+02                        | Normal/Lognormal                  | 1.10E-01                                 | 3.19E+01                                    | no                                      |  |
| Chrysene               | 1.30E+00         | 2.63E+05                        | Normal/Lognormal                  | 8.70E-01                                 | 8.75E+01                                    | no                                      | YES*                                       |
| Dibenz(a,h)anthracene  | 2.26E-01         | 1.35E+02                        | Normal/Lognormal                  | 1.60E-01                                 | 8.75E+02                                    | YES                                     | YES - COPC                                 |
| Dibenzofuran           | 1.16E-01         | 5.59E-01                        | Normal/Lognormal                  | 9.30E-02                                 | 3.13E+02                                    | no                                      |  |
| Di-t-butyl phthalate   | 1.46E-01         | 1.52E+00                        | Normal/Lognormal                  | 1.10E-01                                 | 2.28E+03                                    | no                                      |  |
| Fluoranthene           | 1.13E+00         | 1.59E+04                        | Normal/Lognormal                  | 7.80E-01                                 | 3.13E+03                                    | no                                      |  |
| Indeno(1,2,3-cd)pyrene | 8.36E-01         | 1.56E+04                        | Normal/Lognormal                  | 6.00E-01                                 | 8.75E-01                                    | no                                      | YES  |
| Naphthalene            | 2.04E-01         | 6.64E-05                        | Unknown                           | 1.60E-01                                 | 6.45E-02                                    | no                                      |  |
| Phenanthrene           | 3.29E-01         | 2.65E-07                        | Normal/Lognormal                  | 2.50E-01                                 | 2.35E+03                                    | no                                      |  |
| Pyrene                 | 1.52E+00         | 7.45E-05                        | Normal/Lognormal                  | 1.00E+00                                 | 2.35E+03                                    | no                                      |  |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 7

*Statistical Summary and Selection of COPCs in EU3 Soil (0-6' bgs)*  
*Kerr McGee, Hattiesburg, MS*

| Constituent            | CAS Number | Total Samples | Hlt Hits | Hlt Frequency % | Minimum mg/kg  | Maximum mg/kg  | Logarithmic Mean mg/kg | Maximum mg/kg | Detected Qualifier | Detected Qualifier | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------|-----------------|----------------|----------------|------------------------|---------------|--------------------|--------------------|-----------------------------|-----------------------------------|--------------------------|
|                        |            |               |          |                 | Detected Limit | Detected Limit | Mean                   | Detected      | Detected           | Detected           |                             |                                   |                          |
| Semivolatiles          |            |               |          |                 |                |                |                        |               |                    |                    |                             |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 7             | 1        | 14.29           | 3.30E-02       | 4.00E-02       | 2.30E-01               | J             | 4.85E-02           | 2.62E-02           | 2.30E-01                    | J                                 | SS-16                    |
| Acenaphthylene         | 208-96-8   | 7             | 2        | 28.57           | 3.30E-02       | 4.00E-02       | 1.20E-01               | J             | 5.47E-02           | 3.33E-02           | 1.70E-01                    | J                                 | SS-16                    |
| Anthracene             | 120-12-7   | 7             | 2        | 28.57           | 3.30E-02       | 4.00E-02       | 1.20E-01               | J             | 5.47E-02           | 3.33E-02           | 1.70E-01                    | J                                 | SS-16                    |
| Benz(a)anthracene      | 56-55-3    | 7             | 3        | 42.86           | 3.70E-02       | 4.00E-02       | 5.60E-02               | J             | 1.66E-01           | 5.71E-02           | 5.40E-01                    | J                                 | SS-15                    |
| Benz(a)pyrene          | 50-32-8    | 7             | 2        | 28.57           | 3.70E-02       | 6.70E-02       | 5.60E-01               | J             | 1.97E-01           | 5.62E-02           | 7.10E-01                    | J                                 | SS-16                    |
| Benz(b)fluoranthene    | 205-99-2   | 7             | 3        | 42.86           | 3.70E-02       | 4.00E-02       | 1.90E-01               | J             | 4.10E-01           | 8.85E-02           | 1.40E+00                    | J                                 | SS-16                    |
| Benz(g)iperylene       | 191-24-2   | 7             | 3        | 42.86           | 3.70E-02       | 4.00E-02       | 8.00E-02               | J             | 2.91E-01           | 7.06E-02           | 1.20E+00                    | J                                 | SS-16                    |
| Benz(k)fluoranthene    | 207-08-9   | 7             | 2        | 28.57           | 3.70E-02       | 1.30E-01       | 4.70E-01               | J             | 1.57E-01           | 5.72E-02           | 4.90E-01                    | J                                 | SS-16                    |
| Carbazole              | 86-74-8    | 7             | 2        | 28.57           | 3.30E-02       | 4.00E-02       | 4.60E-02               | J             | 3.56E-02           | 2.72E-02           | 1.10E-01                    | J                                 | SS-15                    |
| Chrysene               | 218-01-9   | 7             | 3        | 42.86           | 3.70E-02       | 4.00E-02       | 1.10E-01               | J             | 2.65E-01           | 7.22E-02           | 8.70E-01                    | J                                 | SS-16                    |
| Dibenz(a,h)anthracene  | 53-70-3    | 7             | 2        | 28.57           | 3.70E-02       | 6.70E-02       | 1.49E-01               | J             | 5.86E-02           | 3.73E-02           | 1.60E-01                    | J                                 | SS-16                    |
| Dibenzofuran           | 132-64-9   | 7             | 2        | 28.57           | 3.30E-02       | 4.00E-02       | 3.60E-02               | J             | 3.17E-02           | 2.57E-02           | 9.30E-02                    | J                                 | SS-16                    |
| Di-n-butylphthalate    | 84-74-2    | 7             | 3        | 42.86           | 7.50E-02       | 7.90E-02       | 4.00E-02               | J             | 5.74E-02           | 5.13E-02           | 1.10E-01                    | J                                 | SS-16                    |
| Fluoranthene           | 206-44-0   | 7             | 3        | 42.86           | 3.70E-02       | 4.00E-02       | 1.20E-01               | J             | 2.37E-01           | 7.03E-02           | 7.80E-01                    | J                                 | SS-16                    |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 7             | 3        | 42.86           | 3.70E-02       | 4.00E-02       | 8.60E-02               | J             | 1.76E-01           | 6.13E-02           | 6.00E-01                    | J                                 | SS-16                    |
| Naphthalene            | 91-20-3    | 7             | 1        | 14.29           | 3.30E-02       | 4.00E-02       | 1.60E-01               | J             | 3.85E-02           | 2.48E-02           | 1.60E-01                    | J                                 | SS-16                    |
| Phenanthrene           | 85-01-8    | 7             | 2        | 28.57           | 3.30E-02       | 4.00E-02       | 1.30E-01               | J             | 6.76E-02           | 3.55E-02           | 2.50E-01                    | J                                 | SS-16                    |
| Phenol                 | 108-95-2   | 7             | 2        | 28.57           | 3.30E-02       | 7.90E-02       | 9.60E-02               | J             | 6.47E-02           | 3.95E-02           | 2.30E-01                    | J                                 | GEO-17                   |
| Pyrene                 | 129-00-0   | 7             | 3        | 42.86           | 3.70E-02       | 4.00E-02       | 1.20E-01               | J             | 3.07E-01           | 7.64E-02           | 1.00E+00                    | J                                 | 4.58E-01                 |

**Table 7**  
**Statistical Summary and Selection of COPCs in EU3 Soil (0-6' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | Logarithmic      |                  | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? |
|------------------------|------------------|------------------|-----------------------------------|--|------------------------|---|
|                        | 95% UCL<br>mg/kg | 95% LCL<br>mg/kg |                                   |  |                        |   |
| <b>Semivolatiles</b>   |                  |                  |                                   |  |                        |   |
| 2-methylnaphthalene    | 1.07E-01         | 1.70E-01         | Unknown                           | 1.70E-01                                 | 8.18E+04               | no                                      |
| Acenaphthylene         | 1.01E-01         | 2.50E-01         | Unknown                           | 1.70E-01                                 | 1.23E+05               | no                                      |
| Anthracene             | 1.01E-01         | 2.50E-01         | Unknown                           | 1.70E-01                                 | 6.13E+05               | no                                      |
| Benz(a)anthracene      | 3.42E-01         | 5.64E+00         | Unknown                           | 5.40E-01                                 | 7.84E+00               | no                                      |
| Benz(a)pyrene          | 4.19E-01         | 1.08E+01         | Unknown                           | 7.10E-01                                 | 7.84E-01               | no                                      |
| Benz(b)fluoranthene    | 8.61E-01         | 1.83E+02         | Lognormal                         | 1.40E+00                                 | 7.84E+00               | no                                      |
| Benz(ghi)perylene      | 6.35E-01         | 3.79E+01         | Lognormal                         | 1.20E+00                                 | 6.13E+04               | no                                      |
| Benz(k)fluoranthene    | 3.20E-01         | 4.72E+00         | Lognormal                         | 4.90E-01                                 | 7.84E+01               | no                                      |
| Carbazole              | 6.08E-02         | 8.08E-02         | Unknown                           | 8.08E-02                                 | 2.86E+02               | no                                      |
| Chrysene               | 5.52E-01         | 3.06E+01         | Lognormal                         | 8.70E-01                                 | 7.84E+02               | no                                      |
| Dibenz(a,h)anthracene  | 1.05E-01         | 2.52E-01         | Unknown                           | 1.60E-01                                 | 7.84E-01               | no                                      |
| Dibenzofuran           | 5.21E-02         | 6.21E-02         | Unknown                           | 6.21E-02                                 | 8.18E+03               | no                                      |
| Di-n-butylphthalate    | 8.12E-02         | 9.38E-02         | Unknown                           | 9.38E-02                                 | 2.28E+03               | no                                      |
| Fluoranthene           | 4.87E-01         | 2.06E+01         | Lognormal                         | 7.80E-01                                 | 8.17E+04               | no                                      |
| Indeno(1,2,3-cd)pyrene | 3.59E-01         | 6.92E+00         | Lognormal                         | 6.00E-01                                 | 7.84E+00               | no                                      |
| Naphthalene            | 7.79E-02         | 1.03E-01         | Unknown                           | 1.03E-01                                 | 8.24E+02               | no                                      |
| Phenanthrene           | 1.34E-01         | 4.34E-01         | Unknown                           | 2.50E-01                                 | 6.13E+04               | no                                      |
| Phenol                 | 1.22E-01         | 3.11E-01         | Lognormal                         | 2.30E-01                                 | 1.23E+05               | no                                      |
| Pyrene                 | 6.43E-01         | 5.32E+01         | Lognormal                         | 1.00E+00                                 | 6.13E+04               | no                                      |

Table 8

Statistical Summary and Selection of COPCs in EU3 Soil (0-20' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg   | Maximum mg/kg   | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected Concentration | Location of Maximum Concentration | Standard Deviation mg/kg |        |          |
|------------------------|------------|---------------|----------------|-----------------|-----------------|-----------------|------------|------------------------|---------------|------------------------|-----------------------------------|--------------------------|--------|----------|
|                        |            |               |                |                 | Detection Limit | Detection Limit | Detected   | Detected               | Detected      | Detected               |                                   |                          |        |          |
|                        |            |               |                |                 | mg/kg           | mg/kg           | Qualifier  | mg/kg                  | mg/kg         | Qualifier              |                                   |                          |        |          |
| Semivolatiles          |            |               |                |                 |                 |                 |            |                        |               |                        |                                   |                          |        |          |
| 2-methylnaphthalene    | 91-57-6    | 7             | 1              | 14.29           | 3.30E-02        | 4.00E-02        | 2.30E-01   | J                      | 4.85E-02      | 2.62E-02               | 2.30E-01                          | J                        | SS-16  | 8.00E-02 |
| Acenaphthylene         | 208-96-8   | 7             | 2              | 28.57           | 3.30E-02        | 4.00E-02        | 1.20E-01   | J                      | 5.47E-02      | 3.33E-02               | 1.70E-01                          | J                        | SS-16  | 6.34E-02 |
| Anthracene             | 120-12-7   | 7             | 2              | 28.57           | 3.30E-02        | 4.00E-02        | 1.20E-01   | J                      | 5.47E-02      | 3.33E-02               | 1.70E-01                          | J                        | SS-16  | 6.34E-02 |
| Benz(a)anthracene      | 56-55-3    | 7             | 3              | 42.86           | 3.70E-02        | 4.00E-02        | 5.60E-02   | J                      | 1.66E-01      | 5.71E-02               | 5.40E-01                          | J                        | SS-15  | 2.39E-01 |
| Benz(a)pyrene          | 50-32-8    | 7             | 2              | 28.57           | 3.70E-02        | 6.70E-02        | 5.60E-01   | J                      | 1.97E-01      | 5.62E-02               | 7.10E-01                          | J                        | SS-16  | 3.02E-01 |
| Benz(b)fluoranthene    | 205-99-2   | 7             | 3              | 42.86           | 3.70E-02        | 4.00E-02        | 1.90E-01   | J                      | 4.10E-01      | 8.85E-02               | 1.40E+00                          | J                        | SS-16  | 6.14E-01 |
| Benz(ghi)perylene      | 191-24-2   | 7             | 3              | 42.86           | 3.70E-02        | 4.00E-02        | 8.00E-02   | J                      | 2.91E-01      | 7.06E-02               | 1.20E+00                          | J                        | SS-16  | 4.69E-01 |
| Benz(k)fluoranthene    | 207-08-9   | 7             | 2              | 28.57           | 3.70E-02        | 1.30E-01        | 4.70E-01   | J                      | 1.57E-01      | 5.72E-02               | 4.90E-01                          | J                        | SS-16  | 2.21E-01 |
| Carbazole              | 86-74-8    | 7             | 2              | 28.57           | 3.30E-02        | 4.00E-02        | 4.60E-02   | J                      | 3.56E-02      | 2.72E-02               | 1.10E-01                          | J                        | SS-15  | 3.44E-02 |
| Chrysene               | 218-01-9   | 7             | 3              | 42.86           | 3.70E-02        | 4.00E-02        | 1.10E-01   | J                      | 2.65E-01      | 7.22E-02               | 8.70E-01                          | J                        | SS-16  | 3.91E-01 |
| Dibenz(a,b)anthracene  | 53-70-3    | 7             | 2              | 28.57           | 3.70E-02        | 6.70E-02        | 1.40E-01   | J                      | 5.86E-02      | 3.73E-02               | 1.60E-01                          | J                        | SS-16  | 6.29E-02 |
| Dibenzofuran           | 132-64-9   | 7             | 2              | 28.57           | 3.30E-02        | 4.00E-02        | 3.60E-02   | J                      | 3.17E-02      | 2.57E-02               | 9.30E-02                          | J                        | SS-16  | 2.78E-02 |
| Di-n-butylphthalate    | 84-74-2    | 7             | 3              | 42.86           | 7.50E-02        | 7.90E-02        | 4.90E-02   | J                      | 5.74E-02      | 5.13E-02               | 1.10E-01                          | J                        | SS-16  | 3.23E-02 |
| Fluoranthene           | 206-44-0   | 7             | 3              | 42.86           | 3.70E-02        | 4.00E-02        | 1.20E-01   | J                      | 2.37E-01      | 7.03E-02               | 7.80E-01                          | J                        | SS-16  | 3.40E-01 |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 7             | 3              | 42.86           | 3.70E-02        | 4.00E-02        | 8.60E-02   | J                      | 1.76E-01      | 6.13E-02               | 6.00E-01                          | J                        | SS-16  | 2.49E-01 |
| Naphthalene            | 91-20-3    | 7             | 1              | 14.29           | 3.30E-02        | 4.00E-02        | 1.60E-01   | J                      | 3.85E-02      | 2.48E-02               | 1.60E-01                          | J                        | SS-16  | 5.36E-02 |
| Phenanthrene           | 85-01-8    | 7             | 2              | 28.57           | 3.30E-02        | 4.00E-02        | 1.30E-01   | J                      | 6.76E-02      | 3.55E-02               | 2.50E-01                          | J                        | SS-16  | 9.05E-02 |
| Phenol                 | 108-93-2   | 7             | 2              | 28.57           | 3.30E-02        | 7.90E-02        | 9.60E-02   | J                      | 6.47E-02      | 3.95E-02               | 2.30E-01                          | J                        | GEO-17 | 7.81E-02 |
| Pyrene                 | 129-00-0   | 7             | 3              | 42.86           | 3.70E-02        | 4.00E-02        | 1.20E-01   | J                      | 3.07E-01      | 7.64E-02               | 1.00E+00                          | J                        | SS-17  | 4.58E-01 |



**Table 8**  
**Statistical Summary and Selection of COPCs in EU3 Soil (0-20' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL,<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? |   |
|------------------------|------------------|----------------------------------|-----------------------------------|--|------------------------|---|---|
|                        |                  |                                  |                                   |  |                        | Tier I<br>Soil<br>TRG<br>mg/kg          | Tier I<br>Is the<br>Maximum<br>Detected ><br>TRG? |
| <b>Semi-volatiles</b>  |                  |                                  |                                   |  |                        |   |   |
| 2-methylnaphthalene    | 1.07E-01         | 1.70E-01                         | Unknown                           | 1.70E-01                                 | 8.18E+04               | no                                      |   |
| Acenaphthylene         | 1.01E-01         | 2.50E-01                         | Unknown                           | 1.70E-01                                 | 1.23E+05               | no                                      |   |
| Anthracene             | 1.01E-01         | 2.50E-01                         | Unknown                           | 1.70E-01                                 | 6.13E+05               | no                                      |   |
| Benz(a)anthracene      | 3.42E-01         | 5.64E+00                         | Unknown                           | 5.40E-01                                 | 7.84E+00               | no                                      |   |
| Benz(a)pyrene          | 4.19E-01         | 1.08E+01                         | Unknown                           | 7.10E-01                                 | 7.84E-01               | no                                      |   |
| Benz(b)fluoranthene    | 8.61E-01         | 1.83E+02                         | Lognormal                         | 1.40E+00                                 | 7.84E+00               | no                                      |   |
| Benz(g,h)perylene      | 6.35E-01         | 3.79E+01                         | Lognormal                         | 1.20E+00                                 | 6.13E+04               | no                                      |   |
| Benz(k)fluoranthene    | 3.20E-01         | 4.72E+00                         | Lognormal                         | 4.90E-01                                 | 7.84E+01               | no                                      |   |
| Carbazole              | 6.08E-02         | 8.08E-02                         | Unknown                           | 8.08E-02                                 | 2.86E+02               | no                                      |   |
| Chrysene               | 5.52E-01         | 3.06E+01                         | Lognormal                         | 8.70E-01                                 | 7.84E+02               | no                                      |   |
| Dibenz(a,h)anthracene  | 1.05E-01         | 2.52E-01                         | Unknown                           | 1.60E-01                                 | 7.84E-01               | no                                      |   |
| Dibenzofuran           | 5.21E-02         | 6.21E-02                         | Unknown                           | 6.21E-02                                 | 8.18E+03               | no                                      |   |
| Di-n-butylphthalate    | 8.12E-02         | 9.38E-02                         | Unknown                           | 9.38E-02                                 | 2.28E+03               | no                                      |   |
| Fluoranthene           | 4.87E-01         | 2.06E+01                         | Lognormal                         | 7.80E-01                                 | 8.17E+04               | no                                      |   |
| Indeno(1,2,3-cd)pyrene | 3.59E-01         | 6.92E+00                         | Lognormal                         | 6.00E-01                                 | 7.84E+00               | no                                      |   |
| Naphthalene            | 7.79E-02         | 1.03E-01                         | Unknown                           | 1.03E-01                                 | 8.24E+02               | no                                      |   |
| Phenanthrene           | 1.34E-01         | 4.34E-01                         | Unknown                           | 2.50E-01                                 | 6.13E+04               | no                                      |   |
| Phenol                 | 1.22E-01         | 3.11E-01                         | Lognormal                         | 2.30E-01                                 | 1.23E+05               | no                                      |   |
| Pyrene                 | 6.43E-01         | 5.32E+01                         | Lognormal                         | 1.00E+00                                 | 6.13E+04               | no                                      |   |

Table 9

Statistical Summary and Selection of COPCs in EU4 Sediment  
Kerr McGee, Hattiesburg, MS

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Maximum mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected | Maximum mg/kg | Mean mg/kg | Qualifer | Concentration | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|---------------|------------|------------------------|---------------|----------|---------------|------------|----------|---------------|-----------------------------------|--------------------------|
|                        |            |               |                |                 |               |               |                       |          |               |               |            |                        |               |          |               |            |          |               |                                   |                          |
| <b>Semivolatiles</b>   |            |               |                |                 |               |               |                       |          |               |               |            |                        |               |          |               |            |          |               |                                   |                          |
| 2,4-dimethylphenol     | 105-67-9   | 1             | 1              | 100             | 0.00E+00      | 1.50E+00      | J                     | 1.50E+00 | 1.50E+00      | 1.50E+00      | 1.50E+00   | 1.50E+00               | J             | 1.50E+00 | 1.50E+00      | 1.50E+00   | J        | SD-02         | NA                                |                          |
| 2-methylnaphthalene    | 91-57-6    | 1             | 1              | 100             | 0.00E+00      | 1.50E+00      | J                     | 1.50E+03 | 1.50E+03      | 1.50E+03      | 1.50E+03   | 1.50E+03               | J             | 1.50E+03 | 1.50E+03      | 1.50E+03   | J        | SD-02         | NA                                |                          |
| Acenaphthene           | 83-32-9    | 8             | 2              | 25              | 5.70E-01      | 1.00E+03      | 6.50E+00              | J        | 1.08E+02      | 6.96E+00      | 3.45E+02   | 3.45E+02               | 3.45E+02      | J        | 3.45E+02      | 3.45E+02   | 3.45E+02 | J             | SD-12                             | 1.99E+02                 |
| Acenaphthylene         | 208-96-8   | 8             | 1              | 12.5            | 5.70E-01      | 1.85E+02      | 3.50E+01              | J        | 1.72E+01      | 3.57E+00      | 3.50E+01   | 3.50E+01               | 3.50E+01      | J        | 3.50E+01      | 3.50E+01   | 3.50E+01 | J             | SD-02                             | 3.26E+01                 |
| Anthracene             | 120-12-7   | 8             | 4              | 50              | 3.80E-02      | 5.56E-01      | 1.80E+00              | Z        | 4.01E+02      | 1.78E+00      | 1.90E+03   | 1.90E+03               | 1.90E+03      | Z        | 1.90E+03      | 1.90E+03   | 1.90E+03 | Z             | SD-02                             | 7.57E+02                 |
| Benzo(a)anthracene     | 56-55-3    | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.29E-01              | Z        | 6.52E+01      | 4.98E+00      | 3.30E+02   | 3.30E+02               | 3.30E+02      | Z        | 3.30E+02      | 3.30E+02   | 3.30E+02 | Z             | SD-02                             | 1.23E+02                 |
| Benzo(a)pyrene         | 50-32-8    | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.80E-01              | Z        | 2.69E+01      | 4.43E+00      | 1.30E+02   | 1.30E+02               | 1.30E+02      | Z        | 1.30E+02      | 1.30E+02   | 1.30E+02 | Z             | SD-02                             | 4.81E+01                 |
| Benzo(b)fluoranthene   | 205-99-2   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 4.25E-01              | Z        | 3.29E+01      | 6.00E+00      | 1.80E+02   | 1.80E+02               | 1.80E+02      | Z        | 1.80E+02      | 1.80E+02   | 1.80E+02 | Z             | SD-02                             | 6.31E+01                 |
| Benzo(g,h)perylene     | 191-24-2   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 1.73E-01              | J        | 7.24E+00      | 2.05E+00      | 3.60E+01   | 3.60E+01               | 3.60E+01      | J        | 3.60E+01      | 3.60E+01   | 3.60E+01 | J             | SD-02                             | 1.25E+01                 |
| Benzo(k)fluoranthene   | 207-08-9   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.13E-01              | Z        | 1.38E+01      | 2.88E+00      | 6.40E+01   | 6.40E+01               | 6.40E+01      | Z        | 6.40E+01      | 6.40E+01   | 6.40E+01 | Z             | SD-02                             | 2.38E+01                 |
| Carbazole              | 86-74-8    | 1             | 1              | 100             | 0.00E+00      | 0.00E+00      | 5.90E+02              | Z        | 5.90E+02      | 5.90E+02      | 5.90E+02   | 5.90E+02               | 5.90E+02      | Z        | 5.90E+02      | 5.90E+02   | 5.90E+02 | Z             | SD-02                             | NA                       |
| Chrysene               | 218-01-9   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.50E-01              | Z        | 5.44E+01      | 4.85E+00      | 2.90E+02   | 2.90E+02               | 2.90E+02      | Z        | 2.90E+02      | 2.90E+02   | 2.90E+02 | Z             | SD-02                             | 1.05E+02                 |
| Dibenz(s,h)anthracene  | 53-70-3    | 8             | 7              | 87.5            | 6.00E-02      | 6.00E-02      | 5.70E-02              | J        | 2.93E+00      | 5.99E-01      | 1.20E+01   | 1.20E+01               | 1.20E+01      | J        | 1.20E+01      | 1.20E+01   | 1.20E+01 | J             | SD-02                             | 4.76E+00                 |
| Dibenzofuran           | 132-64-9   | 1             | 1              | 100             | 0.00E+00      | 0.00E+00      | 9.40E+02              | Z        | 9.40E+02      | 9.40E+02      | 9.40E+02   | 9.40E+02               | 9.40E+02      | Z        | 9.40E+02      | 9.40E+02   | 9.40E+02 | Z             | SD-02                             | NA                       |
| Fluoranthene           | 206-44-0   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.60E-01              | Z        | 3.27E+02      | 1.42E+01      | 1.60E+03   | 1.60E+03               | 1.60E+03      | Z        | 1.60E+03      | 1.60E+03   | 1.60E+03 | Z             | SD-02                             | 6.16E+02                 |
| Fluorene               | 86-73-7    | 8             | 3              | 37.5            | 5.30E-02      | 4.50E-01      | 7.40E+00              | Z        | 2.32E+02      | 1.93E+00      | 1.20E+03   | 1.20E+03               | 1.20E+03      | Z        | 1.20E+03      | 1.20E+03   | 1.20E+03 | Z             | SD-02                             | 4.51E+02                 |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 2.23E-01              | J        | 1.08E+01      | 2.98E+00      | 4.70E+01   | 4.70E+01               | 4.70E+01      | J        | 4.70E+01      | 4.70E+01   | 4.70E+01 | J             | SD-02                             | 1.75E+01                 |
| Naphthalene            | 91-20-3    | 8             | 2              | 25              | 5.70E-01      | 1.85E+02      | 8.20E+00              | J        | 3.89E+02      | 7.60E+00      | 3.00E+03   | 3.00E+03               | 3.00E+03      | J        | 3.00E+03      | 3.00E+03   | 3.00E+03 | J             | SD-02                             | 1.06E+03                 |
| Phenanthrene           | 85-01-8    | 8             | 3              | 37.5            | 3.10E-02      | 1.05E+00      | 2.38E+01              | Z        | 6.49E+02      | 3.49E+00      | 3.20E+03   | 3.20E+03               | 3.20E+03      | Z        | 3.20E+03      | 3.20E+03   | 3.20E+03 | Z             | SD-02                             | 1.24E+03                 |
| Pyrene                 | 129-00-0   | 8             | 8              | 100             | 0.00E+00      | 0.00E+00      | 4.59E-01              | J        | 2.48E+02      | 1.67E+01      | 1.00E+03   | 1.00E+03               | 1.00E+03      | J        | 1.00E+03      | 1.00E+03   | 1.00E+03 | J             | SD-02                             | 4.44E+02                 |

**Table 9**  
**Statistical Summary and Selection of COPCs in EU4 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent             | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg | Logarithmic<br>99%<br>Confidence | Distribution<br>99%<br>Concentration | Exposure Point<br>Unrestricted<br>Concentration<br>mg/kg | Tier I<br>Soil TRG<br>mg/kg | Is the<br>95%<br>UCL ><br>TRG? |                               |
|-------------------------|------------------|------------------|----------------------------------|--------------------------------------|--|-----------------------------|--------------------------------|-------------------------------|
|                         |                  |                  |                                  |                                      |  |                             | Detected ><br>Soil TRG         | Maximum<br>Detected ><br>TRG? |
| <b>Semi-volatiles</b>   |                  |                  |                                  |                                      |  |                             |                                |                               |
| 2,4-dimethylphenol      | NA               | NA               | Unknown                          | 1.50E+00                             | 1.56E+03   | no                          |                                |                               |
| 2-methylnaphthalene     | NA               | NA               | Unknown                          | 1.50E+03                             | 3.13E+03   | no                          |                                |                               |
| Acenaphthene            | 2.41E+02         | 8.24E+05         | Lognormal                        | 3.45E+02                             | 4.69E+03   | no                          |                                |                               |
| Acenaphthylene          | 3.90E+01         | 1.12E+03         | Lognormal                        | 3.50E+01                             | 4.69E+03   | no                          |                                |                               |
| Anthracene              | 9.08E+02         | 1.74E+15         | Lognormal                        | 1.90E+03                             | 2.35E+04   | no                          |                                |                               |
| Benz(a)anthracene       | 1.48E+02         | 3.91E+05         | Lognormal                        | 3.30E+02                             | 8.75E-01   | YES                         | YES - COPC                     |                               |
| Benzo(a)pyrene          | 5.91E+01         | 6.94E+03         | Lognormal                        | 1.30E+02                             | 8.75E-02   | YES                         | YES - COPC                     |                               |
| Benzo(b)fluoranthene    | 7.52E+01         | 4.80E+03         | Lognormal                        | 1.80E+02                             | 8.75E-01   | YES                         | YES - COPC                     |                               |
| Benzo(ghi)perylene      | 1.56E+01         | 2.97E+02         | Lognormal                        | 3.00E+01                             | 2.35E+03   | no                          |                                |                               |
| Benzo(k)fluoranthene    | 2.98E+01         | 1.74E+03         | Lognormal                        | 6.40E+01                             | 8.75E+00   | YES                         | YES - COPC                     |                               |
| Carbazole               | NA               | NA               | Unknown                          | 5.90E+02                             | 3.19E+01   | YES                         | YES - COPC                     |                               |
| Chrysene                | 1.25E+02         | 1.18E+05         | Lognormal                        | 2.90E+02                             | 8.75E+01   | YES                         | YES - COPC                     |                               |
| Dibenz(a,h)anthracene   | 6.12E+00         | 9.77E+02         | Lognormal                        | 1.20E+01                             | 8.75E-02   | YES                         | YES - COPC                     |                               |
| Dibenzofuran            | NA               | NA               | Unknown                          | 9.40E+02                             | 3.13E+02   | YES                         | YES - COPC                     |                               |
| Fluoranthene            | 7.40E+02         | 6.03E+07         | Lognormal                        | 1.60E+03                             | 3.13E+03   | no                          |                                |                               |
| Fluorene                | 5.34E+02         | 1.26E+12         | Lognormal                        | 1.20E+03                             | 3.13E+03   | no                          |                                |                               |
| Indeno(1,2,3-c,d)pyrene | 2.26E+01         | 5.97E+02         | Lognormal                        | 4.70E+01                             | 8.75E-01   | YES                         | YES - COPC                     |                               |
| Naphthalene             | 1.10E+03         | 8.93E+06         | Lognormal                        | 3.00E+03                             | 6.45E+02   | YES                         | YES - COPC                     |                               |
| Phenanthrene            | 1.48E+03         | 1.30E+15         | Lognormal                        | 3.20E+03                             | 2.35E+03   | YES                         | YES - COPC                     |                               |
| Pyrene                  | 5.46E+02         | 5.33E+06         | Lognormal                        | 1.00E+03                             | 2.35E+03   | no                          |                                |                               |



Table 10

*Statistical Summary and Selection of COPCs in EU4 Surface Water  
Kerr McGee, Hattiesburg, MS*

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum Detection Limit mg/L | Maximum Detection Limit mg/L | Detected | Minimum Detected Qualifier mg/L | Maximum Detected Qualifier mg/L | Detected | Maximum Detected Qualifier mg/L | Detected | Maximum Detected Qualifier mg/L | Location of Maximum Concentration | Standard Deviation mg/L |  |
|----------------------------|------------|---------------|----------------|-----------------|------------------------------|------------------------------|----------|---------------------------------|---------------------------------|----------|---------------------------------|----------|---------------------------------|-----------------------------------|-------------------------|--|
|                            |            |               |                |                 |                              |                              |          |                                 |                                 |          |                                 |          |                                 |                                   |                         |  |
| Semivolatiles              |            |               |                |                 |                              |                              |          |                                 |                                 |          |                                 |          |                                 |                                   |                         |  |
| Acenaphthene               | 83-32-9    | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 1.40E-02 | 1.40E-02                        | 1.40E-02                        | 1.40E-02 | 1.40E-02                        | 1.40E-02 | 1.40E-02                        | SW-02                             | 0.00E+00                |  |
| Anthracene                 | 120-12-7   | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 1.30E-02 | 1.30E-02                        | 1.30E-02                        | 1.30E-02 | 1.30E-02                        | 1.30E-02 | 1.30E-02                        | SW-02                             | 0.00E+00                |  |
| Benz(a)anthracene          | 56-55-3    | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 5.00E-03 | 5.00E-03                        | 5.00E-03                        | 5.00E-03 | 5.00E-03                        | 5.00E-03 | 5.00E-03                        | SW-02                             | 0.00E+00                |  |
| Benz(a)pyrene              | 50-32-8    | 1             | 0              | 0               | 1.00E-03                     | 1.00E-03                     | 0.00E+00 | NA                              | 5.00E-04                        | 5.00E-04 | 0.00E+00                        | 5.00E-04 | 0.00E+00                        | NA                                | 0.00E+00                |  |
| Benz(b)fluoranthene        | 205-99-2   | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 1.20E-02 | 1.20E-02                        | 1.20E-02                        | 1.20E-02 | 1.20E-02                        | 1.20E-02 | 1.20E-02                        | SW-02                             | 0.00E+00                |  |
| Benz(k)fluoranthene        | 207-08-9   | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 2.00E-03 | 2.00E-03                        | 2.00E-03                        | 2.00E-03 | 2.00E-03                        | 2.00E-03 | 2.00E-03                        | J                                 | 0.00E+00                |  |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 3.00E-03 | 3.00E-03                        | 3.00E-03                        | 3.00E-03 | 3.00E-03                        | 3.00E-03 | 3.00E-03                        | J                                 | 0.00E+00                |  |
| Carbazole                  | 86-74-8    | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 1.00E-02 | 1.00E-02                        | 1.00E-02                        | 1.00E-02 | 1.00E-02                        | 1.00E-02 | 1.00E-02                        | J                                 | 0.00E+00                |  |
| Chrysene                   | 218-01-9   | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 6.00E-03 | 6.00E-03                        | 6.00E-03                        | 6.00E-03 | 6.00E-03                        | 6.00E-03 | 6.00E-03                        | J                                 | 0.00E+00                |  |
| Dibenz(a,h)anthracene      | 53-70-3    | 1             | 0              | 0               | 1.00E-03                     | 1.00E-03                     | 0.00E+00 | NA                              | 5.00E-04                        | 5.00E-04 | 0.00E+00                        | 5.00E-04 | 0.00E+00                        | NA                                | 0.00E+00                |  |
| Dibenzofuran               | 132-56-9   | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 1.10E-02 | 1.10E-02                        | 1.10E-02                        | 1.10E-02 | 1.10E-02                        | 1.10E-02 | 1.10E-02                        | SW-02                             | 0.00E+00                |  |
| Fluoranthene               | 206-44-0   | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 3.90E-02 | 3.90E-02                        | 3.90E-02                        | 3.90E-02 | 3.90E-02                        | 3.90E-02 | 3.90E-02                        | SW-02                             | 0.00E+00                |  |
| Fluorene                   | 86-73-7    | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 1.20E-02 | 1.20E-02                        | 1.20E-02                        | 1.20E-02 | 1.20E-02                        | 1.20E-02 | 1.20E-02                        | SW-02                             | 0.00E+00                |  |
| Indeno[1,2,3-cd]pyrene     | 193-39-5   | 1             | 0              | 0               | 1.00E-03                     | 1.00E-03                     | 0.00E+00 | NA                              | 5.00E-04                        | 5.00E-04 | 0.00E+00                        | 5.00E-04 | 0.00E+00                        | NA                                | 0.00E+00                |  |
| Phenanthrene               | 85-01-8    | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 1.70E-02 | 1.70E-02                        | 1.70E-02                        | 1.70E-02 | 1.70E-02                        | 1.70E-02 | 1.70E-02                        | SW-02                             | 0.00E+00                |  |
| Pyrene                     | 129-00-0   | 1             | 1              | 100             | 0.00E+00                     | 0.00E+00                     | 2.10E-02 | 2.10E-02                        | 2.10E-02                        | 2.10E-02 | 2.10E-02                        | 2.10E-02 | 2.10E-02                        | SW-02                             | 0.00E+00                |  |

NA - Not applicable: constituent not detected in media.

**Table 10**  
**Statistical Summary and Selection of COPCs in EU4 Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | 95% UCL<br>mg/L | Logarithmic<br>95% UCL<br>mg/L | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/L | Human Health                            |                                   |
|----------------------------|-----------------|--------------------------------|--------------------------------|---|---|-----------------------------------|
|                            |                 |                                |                                |   | Consumption<br>& Organisms AWQC<br>mg/L | Is Maximum<br>Detected ><br>AWQC? |
| <b>Semi挥发物</b>             |                 |                                |                                |   |   |                                   |
| Acenaphthene               | NA              | NA                             | Unknown                        | 1.40E-02                                | 1.20E+00                                | no                                |
| Anthracene                 | NA              | NA                             | Unknown                        | 1.30E-02                                | 9.60E+00                                | no                                |
| Benzo(a)anthracene         | NA              | NA                             | Unknown                        | 5.00E-03                                | 4.40E-06                                | YES - COPC                        |
| Benzo(a)pyrene             | NA              | NA                             | Unknown                        | 5.00E-04                                | 4.40E-06                                | YES**                             |
| Benzo(b)fluoranthene       | NA              | NA                             | Unknown                        | 1.20E-02                                | 4.40E-06                                | YES - COPC                        |
| Benzo(k)fluoranthene       | NA              | NA                             | Unknown                        | 2.00E-03                                | 4.40E-06                                | YES - COPC                        |
| Bis(2-ethylhexyl)phthalate | NA              | NA                             | Unknown                        | 3.00E-03                                | 1.80E-03                                | YES - COPC                        |
| Carbazole*                 | NA              | NA                             | Unknown                        | 1.00E-02                                | NA                                      | no                                |
| Chrysene                   | NA              | NA                             | Unknown                        | 6.00E-03                                | 4.40E-06                                | YES - COPC                        |
| Dibenz(a,h)anthracene      | NA              | NA                             | Unknown                        | 5.00E-04                                | 4.40E-06                                | YES**                             |
| Dibenzofuran*              | NA              | NA                             | Unknown                        | 1.10E-02                                | NA                                      | no                                |
| Fluoranthene               | NA              | NA                             | Unknown                        | 3.90E-02                                | 3.00E-01                                | no                                |
| Fluorene                   | NA              | NA                             | Unknown                        | 1.20E-02                                | 1.30E+00                                | no                                |
| Indeno[1,2,3-cd]pyrene     | NA              | NA                             | Unknown                        | 5.00E-04                                | 4.40E-06                                | YES**                             |
| Phenanthrene*              | NA              | NA                             | Unknown                        | 1.70E-02                                | NA                                      | no                                |
| Pyrene                     | NA              | NA                             | Unknown                        | 2.10E-02                                | 9.60E-01                                | no                                |

NA - Not Available

\* Constituent will be retained as a COPC due to lack of published screening criteria.

\*\*Retained as a COPC, as per MDEQ Comments (8/2/2000): constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 11

*Statistical Summary and Selection of COPCs in EU4 Soil (0-1' bgs)*  
*Kerr McGee, Hattiesburg, MS*

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected | Detected | Maximum mg/kg | Location of Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|------------|------------------------|---------------|----------|----------|---------------|---------------------------|--------------------------|
| <b>Semi-volatiles</b>  |            |               |                |                 |               |               |                       |          |               |            |                        |               |          |          |               |                           |                          |
| 2,4-dimethylphenol     | 105-67-9   | 3             | 1              | 33.33           | 4.10E-01      | 9.90E+00      | 2.50E-01              | J        | 1.80E+00      | 6.33E-01   | 2.50E-01               | J             | GEO-19   | 2.73E+00 |               |                           |                          |
| 2-methylnaphthalene    | 91-57-6    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 2.70E-01              | J        | 9.36E-01      | 3.42E+00   | 2.80E+02               | J             | GEO-21   | 1.61E+02 |               |                           |                          |
| 2-methylphenol         | 95-48-7    | 3             | 1              | 33.33           | 2.00E-01      | 5.00E+00      | 7.30E-02              | J        | 8.91E-01      | 2.63E+01   | 7.30E-02               | J             | GEO-19   | 1.39E+00 |               |                           |                          |
| 3- and 4-methylphenol  | 106-44-5   | 3             | 1              | 33.33           | 4.10E-01      | 9.90E+00      | 2.10E-01              | J        | 1.79E+00      | 5.97E-01   | 2.10E-01               | J             | GEO-19   | 2.74E+00 |               |                           |                          |
| Acenaphthene           | 83-32-9    | 6             | 2              | 33.33           | 2.00E-01      | 1.50E+03      | 1.00E+00              | J        | 1.58E+02      | 7.02E+00   | 1.90E+02               | J             | GEO-21   | 3.00E+02 |               |                           |                          |
| Acenaphthylene         | 208-96-8   | 6             | 3              | 50              | 2.80E+00      | 1.50E+03      | 1.40E+00              | J        | 1.37E+02      | 1.34E+01   | 4.70E+01               | J             | GEO-21   | 3.01E+02 |               |                           |                          |
| Anthracene             | 120-12-7   | 6             | 5              | 83.33           | 5.30E-02      | 5.30E+02      | 2.10E+00              | Z        | 6.35E+02      | 2.06E+01   | 3.00E+03               | Z             | GEO-48   | 1.20E+03 |               |                           |                          |
| Benz(a)anthracene      | 56-55-3    | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 2.10E+00              | Z        | 2.16E+02      | 3.61E+01   | 9.30E+02               | Z             | GEO-48   | 3.65E+02 |               |                           |                          |
| Benz(a)pyrene          | 50-32-8    | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 3.00E+00              | Z        | 1.35E+02      | 3.16E+01   | 5.00E+02               | Z             | GEO-48   | 1.99E+02 |               |                           |                          |
| Benz(b)fluoranthene    | 205-99-2   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 3.50E+00              | Z        | 1.86E+02      | 4.59E+01   | 5.30E+02               | Z             | GEO-48   | 2.43E+02 |               |                           |                          |
| Benz(g)phenylacetylene | 191-24-2   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 1.60E+00              | J        | 4.27E+01      | 1.45E+01   | 1.30E+02               | J             | GEO-48   | 5.43E+01 |               |                           |                          |
| Benz(k)fluoranthene    | 207-08-9   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 1.80E+00              | Z        | 8.28E+01      | 1.94E+01   | 2.90E+02               | Z             | GEO-48   | 1.18E+02 |               |                           |                          |
| Carbazole              | 86-74-8    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 6.00E-01              | J        | 7.88E+01      | 9.34E+00   | 2.30E+02               | Z             | GEO-21   | 1.31E+02 |               |                           |                          |
| Chrysene               | 218-01-9   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 2.70E+00              | Z        | 1.79E+02      | 3.72E+01   | 6.90E+02               | Z             | GEO-48   | 2.73E+02 |               |                           |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 4.80E-01              | J        | 1.83E+01      | 4.98E+00   | 6.40E+01               | J             | GEO-48   | 2.57E+01 |               |                           |                          |
| Dibenzofuran           | 132-64-9   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 3.40E-01              | J        | 6.37E+01      | 3.65E+00   | 1.90E+02               | Z             | GEO-21   | 1.09E+02 |               |                           |                          |
| Fluoranthene           | 206-44-0   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 2.80E+00              | Z        | 9.04E+02      | 7.65E+01   | 4.60E+03               | Z             | GEO-48   | 1.83E+03 |               |                           |                          |
| Fluorene               | 86-73-7    | 6             | 4              | 66.67           | 2.00E-01      | 2.60E-01      | 1.40E+00              | Z        | 3.44E+02      | 4.84E+00   | 1.80E+03               | Z             | GEO-48   | 7.21E+02 |               |                           |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 2.00E+00              | Z        | 6.94E+01      | 1.94E+01   | 2.50E+02               | J             | GEO-48   | 9.91E+01 |               |                           |                          |
| Naphthalene            | 91-20-3    | 6             | 4              | 66.67           | 2.80E+00      | 1.20E+01      | 6.80E-01              | J        | 4.50E+02      | 1.30E+01   | 2.20E+03               | J             | GEO-48   | 8.79E+02 |               |                           |                          |
| N-nitrosodiphenylamine | 86-30-6    | 3             | 1              | 33.33           | 3.70E-02      | 5.00E+00      | 2.00E+01              | J        | 9.06E-01      | 2.10E-01   | 2.00E+01               | Z             | GEO-20   | 1.38E+00 |               |                           |                          |
| Phenanthrene           | 85-01-8    | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 3.10E-01              | J        | 1.19E+03      | 2.33E+01   | 6.40E+03               | Z             | GEO-48   | 2.57E+03 |               |                           |                          |
| Pyrene                 | 129-00-0   | 6             | 6              | 100             | 0.00E+00      | 0.00E+00      | 5.10E+00              | Z        | 8.76E+02      | 9.14E+01   | 4.40E+03               | Z             | GEO-20   | 1.74E+03 |               |                           |                          |

**Table 11**  
**Statistical Summary and Selection of COPCs in EU4 Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL  |          | Logarithmic<br>95% UCL |          | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I<br>Unrestricted Soil<br>TRG<br>mg/kg | Is the<br>Maximum<br>Detected ><br>TRG? | Is the 95%<br>UCL > TRG? |
|------------------------|----------|----------|------------------------|----------|-----------------------------------|--|---|---|--------------------------|
|                        | mg/kg    | mg/kg    | mg/kg                  | mg/kg    |                                   |  |   |   |                          |
| <b>Semivolatiles</b>   |          |          |                        |          |                                   |  |   |   |                          |
| 2,4-dimethylphenol     | 6.40E+00 | 1.83E+13 | Normal/Lognormal       | 2.50E-01 | 1.56E+03                          | no                                       |   |   |                          |
| 2-methylnaphthalene    | 3.66E+02 | 4.06E+62 | Lognormal              | 2.80E+02 | 3.13E+03                          | no                                       |   |   |                          |
| 2-methylphenol         | 3.24E+00 | 3.99E+15 | Normal/Lognormal       | 7.30E-02 | 3.91E+03                          | no                                       |   |   |                          |
| 3- and 4-methylphenol  | 6.40E+00 | 9.25E+13 | Lognormal              | 2.10E-01 | 3.91E+02                          | no                                       |   |   |                          |
| Acenaphthene           | 4.05E+02 | 5.72E+11 | Lognormal              | 1.90E+02 | 4.69E+03                          | no                                       |   |   |                          |
| Acenaphthylene         | 3.84E+02 | 4.19E+06 | Lognormal              | 4.70E+01 | 4.69E+03                          | no                                       |   |   |                          |
| Anthracene             | 1.62E+03 | 8.99E+17 | Lognormal              | 3.00E+03 | 2.35E+04                          | no                                       |   |   |                          |
| Benzof(a)anthracene    | 5.17E+02 | 1.56E+07 | Lognormal              | 9.30E+02 | 8.75E+01                          | YES                                      | YES - COPC                                  |   |                          |
| Benz(a)pyrene          | 2.98E+02 | 9.02E+05 | Normal/Lognormal       | 5.00E+02 | 8.75E+02                          | YES                                      | YES - COPC                                  |   |                          |
| Benz(b)fluoranthene    | 3.85E+02 | 1.50E+06 | Normal/Lognormal       | 5.30E+02 | 8.75E+01                          | YES                                      | YES - COPC                                  |   |                          |
| Benz(ghi)perylene      | 8.73E+01 | 2.25E+04 | Normal/Lognormal       | 1.30E+02 | 2.35E+03                          | no                                       |   |   |                          |
| Benz(k)fluoranthene    | 1.80E+02 | 6.89E+05 | Normal/Lognormal       | 2.90E+02 | 8.75E+00                          | YES                                      | YES - COPC                                  |   |                          |
| Carbazole              | 3.00E+02 | 1.23E+39 | Normal/Lognormal       | 2.30E+02 | 3.19E+01                          | YES                                      | YES - COPC                                  |   |                          |
| Chrysene               | 4.04E+02 | 3.19E+06 | Normal/Lognormal       | 6.90E+02 | 8.75E+01                          | YES                                      | YES - COPC                                  |   |                          |
| Dibenz(a,h)anthracene  | 3.95E+01 | 3.75E+04 | Normal/Lognormal       | 6.40E+01 | 8.75E+02                          | YES                                      | YES - COPC                                  |   |                          |
| Dibenzofuran           | 2.48E+02 | 7.02E+50 | Lognormal              | 1.90E+02 | 3.13E+02                          | no                                       |   |   |                          |
| Fluoranthene           | 2.41E+03 | 3.17E+09 | Lognormal              | 4.60E+03 | 3.13E+03                          | YES                                      | YES - COPC                                  |   |                          |
| Fluorene               | 9.37E+02 | 1.95E+16 | Lognormal              | 1.80E+03 | 3.13E+03                          | no                                       |   |   |                          |
| Indeno(1,2,3-cd)pyrene | 1.51E+02 | 1.07E+05 | Normal/Lognormal       | 2.50E+02 | 8.75E+01                          | YES                                      | YES - COPC                                  |   |                          |
| Naphthalene            | 1.17E+03 | 7.77E+12 | Lognormal              | 2.20E+03 | 6.45E+02                          | YES                                      | YES - COPC                                  |   |                          |
| N-nitrosodiphenylamine | 3.24E+00 | 6.32E+24 | Normal/Lognormal       | 2.00E-01 | 1.30E+02                          | no                                       |   |   |                          |
| Phenanthrene           | 3.31E+03 | 9.46E+14 | Lognormal              | 6.40E+03 | 2.35E+03                          | YES                                      | YES - COPC                                  |   |                          |
| Pyrene                 | 2.31E+03 | 6.40E-08 | Lognormal              | 4.40E+03 | 2.35E+03                          | YES                                      | YES - COPC                                  |   |                          |

Table 12

**Statistical Summary and Selection of COPCs in EU4 Soil (0-6' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Maximum mg/kg | Logarithmic Mean | Detected | Maximum mg/kg | Detected | Qualifier | Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|---------------|------------------|----------|---------------|----------|-----------|---------------------|-----------------------------------|--------------------------|
| Semivolatiles          |            |               |                |                 |               |               |                       |          |               |               |                  |          |               |          |           |                     |                                   |                          |
| 2,4-dimethylphenol     | 105-67-9   | 9             | 3              | 33.33           | 7.30E-02      | 9.90E+00      | 7.90E-02              | J        | 1.63E+00      | 2.30E-01      | 8.90E+00         | J        | GEO-21        | 3.16E+00 |           |                     |                                   |                          |
| 2-methylnaphthalene    | 91-57-6    | 9             | 8              | 88.89           | 4.00E-02      | 4.00E-02      | 6.20E-02              | J        | 2.11E+02      | 1.91E+00      | 1.50E+03         | J        | GEO-21        | 4.93E+02 |           |                     |                                   |                          |
| 2-methylphenol         | 95-48-7    | 9             | 1              | 11.11           | 3.70E-02      | 5.00E+00      | 7.30E-02              | J        | 4.61E-01      | 8.75E-02      | 7.30E-02         | J        | GEO-19        | 8.70E-01 |           |                     |                                   |                          |
| 3- and 4-methylphenol  | 106-44-5   | 9             | 1              | 11.11           | 7.30E-02      | 9.90E+00      | 2.10E-01              | J        | 9.17E-01      | 1.82E-01      | 2.10E-01         | J        | GEO-19        | 1.72E+00 |           |                     |                                   |                          |
| Acenaphthene           | 83-32-9    | 18            | 9              | 50              | 4.00E-02      | 1.50E+03      | 9.70E-02              | J        | 1.30E+02      | 2.05E+00      | 1.20E+03         | J        | GEO-21        | 3.21E+02 |           |                     |                                   |                          |
| Acenaphthylene         | 208-96-8   | 18            | 7              | 38.89           | 3.80E-02      | 1.50E+03      | 8.30E-02              | J        | 4.90E+01      | 1.14E+00      | 5.00E+01         | J        | GEO-21        | 1.76E+02 |           |                     |                                   |                          |
| Anthracene             | 120-12-7   | 18            | 14             | 77.78           | 2.60E-03      | 5.30E-02      | 1.20E-01              | Z        | 3.15E+02      | 1.93E+00      | 3.00E+03         | Z        | GEO-48        | 8.06E+02 |           |                     |                                   |                          |
| Benz(a)anthracene      | 56-55-3    | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 4.90E-03              | J        | 9.72E+01      | 2.57E+00      | 9.30E+02         | Z        | GEO-48        | 2.34E+02 |           |                     |                                   |                          |
| Benz(a)pyrene          | 50-32-8    | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 1.10E-02              | J        | 5.71E+01      | 2.04E+00      | 5.00E+02         | Z        | GEO-48        | 1.29E+02 |           |                     |                                   |                          |
| Benz(b)fluoranthene    | 205-99-2   | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 1.10E-02              | J        | 7.89E+01      | 2.53E+00      | 5.30E+02         | Z        | GEO-48        | 1.65E+02 |           |                     |                                   |                          |
| Benz(ghi)perylene      | 191-24-2   | 18            | 17             | 94.44           | 3.80E-02      | 3.80E-02      | 8.90E-03              | J        | 1.86E+01      | 8.49E-01      | 1.30E+02         | J        | GEO-48        | 3.79E+01 |           |                     |                                   |                          |
| Benz(k)fluoranthene    | 207-08-9   | 18            | 16             | 88.89           | 3.80E-02      | 4.00E-02      | 5.60E-03              | J        | 3.27E+01      | 1.03E+00      | 2.90E+02         | Z        | GEO-48        | 7.57E+01 |           |                     |                                   |                          |
| Carbazole              | 86-74-8    | 9             | 8              | 88.89           | 4.00E-02      | 4.00E-02      | 2.10E-01              | J        | 9.80E+01      | 2.78E+00      | 6.20E+02         | Z        | GEO-21        | 2.10E+02 |           |                     |                                   |                          |
| Chrysene               | 218-01-9   | 18            | 17             | 94.44           | 5.10E-03      | 5.10E-03      | 9.90E-03              | J        | 8.58E+01      | 2.42E+00      | 6.90E+02         | Z        | GEO-48        | 1.88E+02 |           |                     |                                   |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 18            | 14             | 77.78           | 2.60E-03      | 4.00E-02      | 3.80E-03              | J        | 7.32E+00      | 2.96E-01      | 6.40E+01         | J        | GEO-48        | 1.68E+01 |           |                     |                                   |                          |
| Dibenzofuran           | 132-64-9   | 9             | 8              | 88.89           | 4.00E-02      | 4.00E-02      | 7.80E-02              | J        | 1.54E+02      | 2.43E+00      | 1.10E+03         | J        | GEO-21        | 3.61E+02 |           |                     |                                   |                          |
| Fluoranthene           | 206-44-0   | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 1.00E-02              | J        | 4.29E+02      | 7.56E+00      | 4.60E+03         | Z        | GEO-48        | 1.15E+03 |           |                     |                                   |                          |
| Fluorene               | 86-73-7    | 18            | 12             | 66.67           | 1.30E-02      | 2.60E-01      | 1.40E-01              | J        | 2.09E+02      | 1.40E+00      | 1.80E+03         | Z        | GEO-48        | 5.31E+02 |           |                     |                                   |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 18            | 17             | 94.44           | 6.00E-03      | 6.00E-03      | 9.00E-03              | J        | 2.83E+01      | 1.05E+00      | 2.50E+02         | J        | GEO-48        | 6.42E+01 |           |                     |                                   |                          |
| Naphthalene            | 91-20-3    | 18            | 12             | 66.67           | 4.00E-02      | 1.20E+01      | 7.60E-02              | J        | 3.73E+02      | 2.99E+00      | 3.50E+03         | Z        | GEO-21        | 9.37E+02 |           |                     |                                   |                          |
| N-nitrosodiphenylamine | 86-30-6    | 9             | 1              | 11.11           | 3.70E-02      | 5.00E+00      | 2.00E-01              | J        | 4.66E-01      | 8.11E-02      | 2.00E-01         | Z        | GEO-20        | 8.69E-01 |           |                     |                                   |                          |
| Phenanthrene           | 85-01-8    | 18            | 17             | 94.44           | 4.00E-02      | 4.00E-02      | 5.30E-03              | J        | 6.51E+02      | 4.31E+00      | 6.40E+03         | Z        | GEO-48        | 1.72E+03 |           |                     |                                   |                          |
| Pyrene                 | 129-00-0   | 18            | 18             | 100             | 0.00E+00      | 0.00E+00      | 1.60E-02              | J        | 3.77E+02      | 7.72E+00      | 4.40E+03         | Z        | GEO-20        | 1.06E+03 |           |                     |                                   |                          |



Table 12

*Statistical Summary and Selection of COPCs in EU4 Soil (0-6' bgs)*  
*Kerr McGee, Hattiesburg, MS*

| Constituent            | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg | Logarithmic<br>99%<br>Confidence | Distribution<br>Concentration<br>mg/kg | Exposure Point<br>TRG | Restricted Soil<br>mg/kg | Tier 1                        |  | Is the 95%<br>UCL ><br>TRG? |
|------------------------|------------------|------------------|----------------------------------|--|-----------------------|--------------------------|-------------------------------|--|-----------------------------|
|                        |                  |                  |                                  |  |                       |                          | Maximum<br>Detected ><br>TRG? |  |                             |
| <b>Semivolatiles</b>   |                  |                  |                                  |  |                       |                          |                               |  |                             |
| 2,4-dimethylphenol     | 3.59E+00         | 1.33E+02         | Lognormal                        | 8.90E+00                               | 4.08E+04              | no                       |                               |  |                             |
| 2-methylnaphthalene    | 5.17E+02         | 1.40E+11         | Lognormal                        | 1.50E+03                               | 8.18E+04              | no                       |                               |  |                             |
| 2-methylphenol         | 1.00E+00         | 1.71E+01         | Lognormal                        | 7.30E-02                               | 1.02E+05              | no                       |                               |  |                             |
| 3- and 4-methylphenol  | 1.98E+00         | 3.49E+01         | Lognormal                        | 2.10E-01                               | 1.02E+04              | no                       |                               |  |                             |
| Acenaphthene           | 2.61E+02         | 7.82E+05         | Lognormal                        | 1.20E+03                               | 1.23E+05              | no                       |                               |  |                             |
| Acenaphthylene         | 1.21E+02         | 8.90E+03         | Lognormal                        | 5.00E+01                               | 1.23E+05              | no                       |                               |  |                             |
| Anthracene             | 6.46E+02         | 3.29E+08         | Lognormal                        | 3.00E+03                               | 6.13E+05              | no                       |                               |  |                             |
| Benzo(a)anthracene     | 1.93E+02         | 1.43E+06         | Lognormal                        | 9.30E+02                               | 7.84E+00              | YES - COPC               |                               |  |                             |
| Benzo(a)pyrene         | 1.10E+02         | 1.87E+05         | Lognormal                        | 5.00E+02                               | 7.84E-01              | YES - COPC               |                               |  |                             |
| Benzo(b)fluoranthene   | 1.47E+02         | 5.68E+05         | Lognormal                        | 5.30E+02                               | 7.84E+00              | YES - COPC               |                               |  |                             |
| Benzo(ghi)perylene     | 3.42E+01         | 2.74E+04         | Lognormal                        | 1.30E+02                               | 6.13E+04              | no                       |                               |  |                             |
| Benzo(k)fluoranthene   | 6.37E+01         | 2.16E+05         | Lognormal                        | 2.90E+02                               | 7.84E+01              | YES - COPC               |                               |  |                             |
| Carbazole              | 2.28E+02         | 1.00E+08         | Lognormal                        | 6.20E+02                               | 2.86E+02              | YES - COPC               |                               |  |                             |
| Chrysene               | 1.63E+02         | 3.02E+06         | Lognormal                        | 6.90E+02                               | 7.84E+02              | no                       | YES*                          |  |                             |
| Dibenz(a,h)anthracene  | 1.44E+01         | 1.67E+04         | Lognormal                        | 6.40E+01                               | 7.84E-01              | YES                      |                               |  |                             |
| Dibenzofuran           | 3.78E+02         | 4.19E+09         | Lognormal                        | 1.10E+03                               | 8.18E+03              | no                       |                               |  |                             |
| Fluoranthene           | 8.99E+02         | 1.07E+07         | Lognormal                        | 4.60E+03                               | 8.17E+04              | no                       |                               |  |                             |
| Fluorene               | 4.27E+02         | 4.71E+07         | Lognormal                        | 1.80E+03                               | 8.17E+04              | no                       |                               |  |                             |
| Indeno(1,2,3-cd)pyrene | 5.47E+01         | 1.10E+05         | Lognormal                        | 2.50E+02                               | 7.84E+00              | YES - COPC               |                               |  |                             |
| Naphthalene            | 7.57E+02         | 2.43E+07         | Lognormal                        | 3.50E+03                               | 8.24E+02              | YES - COPC               |                               |  |                             |
| N-nitrosodiphenylamine | 1.00E+00         | 2.74E+01         | Lognormal                        | 2.00E-01                               | 1.17E+03              | no                       |                               |  |                             |
| Phenanthrene           | 1.35E+03         | 6.44E+08         | Lognormal                        | 6.40E+03                               | 1.26E+02              | YES                      | YES - COPC                    |  |                             |
| Pyrene                 | 8.10E+02         | 5.56E+06         | Lognormal                        | 4.40E+03                               | 2.60E+02              | YES                      | YES - COPC                    |  |                             |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 13

Statistical Summary and Selection of COPCs in EU4 Soil (0-20' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Mean mg/kg | Maximum mg/kg | Detected | Maximum mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|------------|---------------|----------|---------------|-----------------------------------|--------------------------|
|                        |            |               |                |                 |               |               |                       |          |               |            |               |          |               |                                   |                          |
| Semivolatiles          |            |               |                |                 |               |               |                       |          |               |            |               |          |               |                                   |                          |
| 2,4-dimethylphenol     | 105-67-9   | 13            | 4              | 30.77           | 7.30E-02      | 9.90E+00      | 7.90E-02              | J        | 1.21E+00      | 2.11E-01   | 8.90E+00      | J        | GEO-21        | 2.67E+00                          |                          |
| 2-methylnaphthalene    | 91-57-6    | 13            | 12             | 92.31           | 4.00E-02      | 6.20E-02      | J                     | 1.91E+02 | 4.52E+00      | 1.50E+03   | J             | GEO-21   | 4.16E+02      |                                   |                          |
| 2-methylphenol         | 95-48-7    | 13            | 2              | 15.38           | 3.70E-02      | 5.00E+00      | 5.10E-02              | J        | 3.87E-01      | 9.13E-02   | 7.30E-02      | J        | GEO-19        | 7.34E-01                          |                          |
| 3- and 4-methylphenol  | 106-44-5   | 13            | 1              | 7.69            | 7.30E-02      | 9.90E+00      | 2.10E-01              | J        | 7.37E-01      | 1.67E-01   | 2.10E-01      | J        | GEO-19        | 1.45E+00                          |                          |
| Acenaphthene           | 83-32-9    | 22            | 13             | 59.09           | 4.00E-02      | 1.50E+03      | 9.70E-02              | J        | 1.23E+02      | 3.12E+00   | 1.20E+03      | J        | GEO-21        | 2.93E+02                          |                          |
| Acenaphthylene         | 208-96-8   | 22            | 10             | 45.45           | 3.80E-02      | 1.50E+03      | 8.30E-02              | J        | 4.07E+01      | 1.05E+00   | 5.00E+01      | J        | GEO-21        | 1.59E+02                          |                          |
| Anthracene             | 120-12-7   | 22            | 18             | 81.82           | 2.60E-03      | 5.30E-02      | 1.20E-01              | Z        | 2.64E+02      | 2.65E+00   | 3.00E+03      | Z        | GEO-48        | 7.33E+02                          |                          |
| Benz(a)anthracene      | 56-55-3    | 22            | 22             | 100             | 0.00E+00      | 0.00E+00      | 4.90E-03              | J        | 8.38E+01      | 3.08E+00   | 9.30E+02      | Z        | GEO-48        | 2.13E+02                          |                          |
| Benz(a)pyrene          | 50-32-8    | 22            | 22             | 100             | 0.00E+00      | 0.00E+00      | 1.10E-02              | J        | 4.83E+01      | 2.14E+00   | 5.00E+02      | Z        | GEO-48        | 1.18E+02                          |                          |
| Benz(b)fluoranthene    | 205-99-2   | 22            | 22             | 100             | 0.00E+00      | 0.00E+00      | 1.10E-02              | Z        | 6.68E+01      | 2.74E+00   | 5.30E+02      | Z        | GEO-48        | 1.51E+02                          |                          |
| Benz(ghi)perylene      | 191-24-2   | 22            | 20             | 90.91           | 3.80E-02      | 4.00E-02      | 8.90E-03              | J        | 1.57E+01      | 8.12E-01   | 1.30E+02      | J        | GEO-48        | 3.47E+01                          |                          |
| Benz(k)fluoranthene    | 207-08-9   | 22            | 20             | 90.91           | 3.80E-02      | 4.00E-02      | 5.60E-03              | J        | 2.76E+01      | 1.10E+00   | 2.90E+02      | Z        | GEO-48        | 6.91E+01                          |                          |
| Carbazole              | 86-74-8    | 13            | 12             | 92.31           | 4.00E-02      | 4.00E-02      | 2.10E-01              | J        | 7.47E+01      | 3.60E+00   | 6.20E+02      | J        | GEO-21        | 1.76E+02                          |                          |
| Chrysene               | 218-01-9   | 22            | 21             | 95.45           | 5.10E-03      | 5.10E-03      | 9.90E-03              | J        | 7.39E+01      | 2.85E+00   | 6.90E+02      | Z        | GEO-48        | 1.72E+02                          |                          |
| Dibenz(a,h)anthracene  | 53-70-3    | 22            | 16             | 72.73           | 2.60E-03      | 3.30E-01      | 3.80E-03              | J        | 6.30E+00      | 2.88E-01   | 6.40E+01      | J        | GEO-48        | 1.54E+01                          |                          |
| Dibenzofuran           | 132-64-9   | 13            | 12             | 92.31           | 4.00E-02      | 4.00E-02      | 7.80E-02              | J        | 1.33E+02      | 4.64E+00   | 1.10E+03      | J        | GEO-21        | 3.03E+02                          |                          |
| Fluoranthene           | 206-44-0   | 22            | 22             | 100             | 0.00E+00      | 0.00E+00      | 1.00E-02              | J        | 3.76E+02      | 1.01E+01   | 4.60E+03      | Z        | GEO-48        | 1.04E+03                          |                          |
| Fluorene               | 86-73-7    | 22            | 16             | 72.73           | 1.30E-02      | 2.60E-01      | 1.40E-01              | J        | 1.91E+02      | 2.40E+00   | 1.80E+03      | Z        | GEO-48        | 4.82E+02                          |                          |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 22            | 21             | 95.45           | 6.00E-03      | 6.00E-03      | 9.00E-03              | J        | 2.37E+01      | 1.04E+00   | 2.50E+02      | J        | GEO-48        | 5.87E+01                          |                          |
| Naphthalene            | 91-20-3    | 22            | 16             | 72.73           | 4.00E-02      | 1.20E+01      | 7.60E-02              | J        | 3.61E+02      | 5.28E+00   | 3.50E+03      | J        | GEO-21        | 8.57E+02                          |                          |
| N-nitrosodiphenylamine | 86-30-6    | 13            | 1              | 7.69            | 3.70E-02      | 5.00E+00      | 2.00E-01              | J        | 3.88E-01      | 8.06E-02   | 2.00E-01      | J        | GEO-20        | 7.35E-01                          |                          |
| Phenanthrene           | 85-01-8    | 22            | 21             | 95.45           | 4.00E-02      | 4.00E-02      | 5.30E-03              | J        | 5.78E+02      | 7.21E+00   | 6.40E+03      | Z        | GEO-48        | 1.56E+03                          |                          |
| Phenol                 | 108-95-2   | 13            | 1              | 7.69            | 7.30E-02      | 9.90E+00      | 1.00E-01              | J        | 6.65E-01      | 1.34E-01   | 1.00E-01      | J        | GEO-20        | 1.46E+00                          |                          |
| Pyrene                 | 129-00-0   | 22            | 22             | 100             | 0.00E+00      | 0.00E+00      | 1.60E-02              | J        | 3.23E+02      | 9.24E+00   | 4.40E+03      | Z        | GEO-20        | 9.58E+02                          |                          |
| Volatiles              |            |               |                |                 |               |               |                       |          |               |            |               |          |               |                                   |                          |
| Acetone                | 67-64-1    | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 9.00E-03              | J        | 5.45E-02      | 3.00E-02   | 1.00E-01      | J        | SB-05         | 6.43E-02                          |                          |
| Benzene                | 71-43-2    | 2             | 1              | 50              | 1.00E-03      | 1.00E-03      | 7.00E-03              | J        | 3.75E-03      | 1.87E-03   | 7.00E-03      | J        | SB-05         | 4.60E-03                          |                          |
| Ethylbenzene           | 100-41-4   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 4.60E-02              | J        | 8.30E-02      | 7.43E-02   | 1.20E-01      | J        | SB-05         | 5.23E-02                          |                          |
| Styrene                | 100-42-5   | 2             | 1              | 50              | 1.00E-03      | 1.00E-03      | 1.00E-01              | J        | 5.01E-02      | 7.07E-03   | 1.00E-01      | J        | SB-05         | 7.04E-02                          |                          |
| Toluene                | 108-88-3   | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 1.30E-02              | J        | 7.65E-02      | 4.27E-02   | 1.40E-01      | J        | SB-05         | 8.98E-02                          |                          |
| Xylene (total)         | 1330-20-7  | 2             | 2              | 100             | 0.00E+00      | 0.00E+00      | 2.80E-01              | J        | 5.30E-01      | 4.67E-01   | 7.80E-01      | J        | SB-05         | 3.54E-01                          |                          |

**Table 13**  
**Statistical Summary and Selection of COPCs in EU4 Soil (0-20' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Tier I                          |                                   |  | Is the Maximum<br>Detected ><br>TRG? | Is the 95%<br>UCL ><br>TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|--------------------------------------|-----------------------------|
|                        |                  | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg |                                      |                             |
| Semivolatiles          |                  |                                 |                                   |  |                                      |                             |
| 2,4-dimethylphenol     | 2.53E+00         | 9.72E+00                        | Lognormal                         | 8.90E+00                                 | 4.08E+04                             | no                          |
| 2-methylnaphthalene    | 3.96E+02         | 2.39E+08                        | Lognormal                         | 1.50E+03                                 | 8.18E+04                             | no                          |
| 2-methylphenol         | 7.50E-01         | 3.10E+00                        | Lognormal                         | 7.30E-02                                 | 1.02E+05                             | no                          |
| 3- and 4-methylphenol  | 1.45E+00         | 5.96E+00                        | Lognormal                         | 2.10E-01                                 | 1.02E+04                             | no                          |
| Acenaphthene           | 2.30E+02         | 2.70E+05                        | Lognormal                         | 1.20E+03                                 | 1.23E+05                             | no                          |
| Acenaphthylene         | 9.90E+01         | 2.21E+03                        | Lognormal                         | 5.00E+01                                 | 1.23E+05                             | no                          |
| Anthracene             | 5.33E+02         | 1.25E+07                        | Lognormal                         | 3.00E+03                                 | 6.13E+05                             | no                          |
| Benz(a)anthracene      | 1.62E+02         | 1.54E+05                        | Lognormal                         | 9.30E+02                                 | 7.84E+00                             | YES - COPC                  |
| Benz(a)pyrene          | 9.16E+01         | 2.29E+04                        | Lognormal                         | 5.00E+02                                 | 7.84E-01                             | YES - COPC                  |
| Benz(b)fluoranthene    | 1.22E+02         | 5.93E+04                        | Lognormal                         | 5.30E+02                                 | 7.84E+00                             | YES - COPC                  |
| Benz(ghi)perylene      | 2.84E+01         | 4.41E+03                        | Lognormal                         | 1.30E+02                                 | 6.13E+04                             | no                          |
| Benz(k)fluoranthene    | 5.29E+01         | 2.30E+04                        | Lognormal                         | 2.90E+02                                 | 7.84E+01                             | YES - COPC                  |
| Carbazole              | 1.62E+02         | 2.13E+05                        | Lognormal                         | 6.20E+02                                 | 2.86E+02                             | YES - COPC                  |
| Chrysene               | 1.37E+02         | 2.60E+05                        | Lognormal                         | 6.90E+02                                 | 7.84E+02                             | *YES                        |
| Dibenz(a,h)anthracene  | 1.19E+01         | 1.83E+03                        | Lognormal                         | 6.40E+01                                 | 7.84E-01                             | YES - COPC                  |
| Dibenzofuran           | 2.83E+02         | 8.51E+06                        | Lognormal                         | 1.10E+03                                 | 8.18E+03                             | no                          |
| Fluoranthene           | 7.58E+02         | 1.38E+06                        | Lognormal                         | 4.60E+03                                 | 8.17E+04                             | no                          |
| Fluorene               | 3.68E+02         | 9.11E+06                        | Lognormal                         | 1.80E+03                                 | 8.17E+04                             | no                          |
| Indeno(1,2,3-cd)pyrene | 4.53E+01         | 1.15E+04                        | Lognormal                         | 2.50E+02                                 | 7.84E+00                             | YES - COPC                  |
| Naphthalene            | 6.76E+02         | 1.04E+07                        | Lognormal                         | 3.50E+03                                 | 8.24E+02                             | YES - COPC                  |
| N-nitrosodiphenylamine | 7.51E-01         | 4.36E+00                        | Unknown                           | 2.00E-01                                 | 1.17E+03                             | no                          |
| Phenanthrene           | 1.15E+03         | 6.70E+07                        | Lognormal                         | 6.40E+03                                 | 1.26E+02                             | YES - COPC                  |
| Phenol                 | 1.39E+00         | 3.63E+00                        | Unknown                           | 1.00E-01                                 | 1.23E+05                             | no                          |
| Pyrene                 | 6.75E+02         | 5.83E+05                        | Lognormal                         | 4.40E+03                                 | 2.60E+02                             | YES - COPC                  |
| Volatiles              |                  |                                 |                                   |  |                                      |                             |
| Acetone                | 3.42E-01         | 1.51E+34                        | Unknown                           | 1.00E-01                                 | 1.04E+05                             | no                          |
| Benzene                | 2.43E-02         | 1.48E+40                        | Unknown                           | 7.00E-03                                 | 1.36E+00                             | no                          |
| Ethylbenzene           | 3.17E-01         | 2.65E+04                        | Unknown                           | 1.20E-01                                 | 3.95E+02                             | no                          |
| Styrene                | 3.64E-01         | 1.64E+171                       | Unknown                           | 1.00E-01                                 | 3.84E+02                             | no                          |
| Toluene                | 4.77E-01         | 2.57E+33                        | Unknown                           | 1.40E-01                                 | 3.80E+01                             | no                          |
| Xylene (total)         | 2.11E+00         | 1.07E+06                        | Unknown                           | 7.80E-01                                 | 3.18E+02                             | no                          |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 14

*Statistical Summary and Selection of COPCs in EU Soil (0-1' bgs)*  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Maximum mg/kg | Detected | Maximum mg/kg | Detected | Maximum mg/kg | Logarithmic Mean mg/kg | Mean mg/kg | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|---------------|----------|---------------|----------|---------------|------------------------|------------|--------------------------|
| Semi-volatiles             |            |               |                |                 |               |               |                       |          |               |               |          |               |          |               |                        |            |                          |
| 2,2,4-dimethylphenol       | 105-67-9   | 6             | 1              | 16.67           | 7.60E-02      | 4.55E-01      | 1.10E-01              | J        | 1.07E-01      | 8.03E-02      | 1.10E-01 | J             | GEO-30   | 8.38E-02      |                        |            |                          |
| 2,2-methylnaphthalene      | 91-57-6    | 6             | 4              | 66.67           | 3.80E-02      | 5.10E-02      | 5.10E-02              | J        | 1.68E+00      | 1.71E-01      | 9.20E+00 | J             | GEO-30   | 3.69E+00      |                        |            |                          |
| 2,2-methylphenol           | 95-48-7    | 6             | 1              | 16.67           | 3.80E-02      | 2.09E-01      | 4.20E-02              | J        | 4.94E-02      | 3.77E-02      | 4.20E-02 | J             | GEO-30   | 3.92E-02      |                        |            |                          |
| 3,3'-3- and 4-methylphenol | 106-44-5   | 6             | 1              | 16.67           | 7.60E-02      | 4.13E-01      | 1.40E-01              | J        | 1.08E-01      | 8.22E-02      | 1.40E-01 | J             | GEO-30   | 7.94E-02      |                        |            |                          |
| Acenaphthene               | 83-32-9    | 8             | 2              | 25              | 3.80E-02      | 2.40E+00      | 1.60E-01              | J        | 6.91E+00      | 1.92E-01      | 5.35E+01 | J             | GEO-33   | 1.88E+01      |                        |            |                          |
| Acenaphthylene             | 208-96-8   | 8             | 5              | 62.5            | 3.80E-02      | 2.40E+00      | 4.80E-02              | J        | 2.61E+00      | 4.28E-01      | 1.60E+01 | J             | GEO-33   | 5.47E+00      |                        |            |                          |
| Anthracene                 | 120-12-7   | 8             | 4              | 50              | 1.00E-02      | 4.40E-02      | 1.30E-01              | J        | 1.07E+01      | 2.00E-01      | 7.97E+01 | J             | GEO-33   | 2.79E+01      |                        |            |                          |
| Benz(a)anthracene          | 56-55-3    | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 2.60E-01              | Z        | 1.30E+01      | 1.39E+00      | 8.35E+01 | J             | GEO-33   | 2.87E+01      |                        |            |                          |
| Benz(a)pyrene              | 50-32-8    | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 3.10E-01              | J        | 8.82E+00      | 1.38E+00      | 5.25E+01 | J             | GEO-33   | 1.79E+01      |                        |            |                          |
| Benz(b)fluoranthene        | 205-99-2   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 4.40E-01              | Z        | 1.38E+01      | 2.15E+00      | 7.95E+01 | J             | GEO-33   | 2.71E+01      |                        |            |                          |
| Benz(g,h)perylene          | 191-24-2   | 8             | 6              | 75              | 3.80E-02      | 3.80E-02      | 2.30E-01              | J        | 4.30E+00      | 5.42E-01      | 2.55E+01 | J             | GEO-33   | 8.71E+00      |                        |            |                          |
| Benz(k)fluoranthene        | 207-08-9   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 2.10E-01              | Z        | 5.07E+00      | 1.01E+00      | 2.85E+01 | J             | GEO-33   | 9.68E+00      |                        |            |                          |
| Carbazole                  | 86-74-8    | 6             | 3              | 50              | 3.80E-02      | 3.90E-02      | 5.30E-01              | J        | 2.63E+00      | 2.10E-01      | 1.35E+01 | J             | GEO-33   | 5.36E+00      |                        |            |                          |
| Chrysene                   | 218-01-9   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 2.70E-01              | Z        | 1.35E+01      | 1.62E+00      | 8.25E+01 | J             | GEO-33   | 2.83E+01      |                        |            |                          |
| Dibenz(a,h)anthracene      | 53-70-3    | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 6.30E-02              | J        | 1.33E+00      | 3.02E-01      | 7.45E+00 | J             | GEO-33   | 2.53E+00      |                        |            |                          |
| Dibenzofuran               | 132-64-9   | 6             | 4              | 66.67           | 3.80E-02      | 3.90E-02      | 3.90E-02              | J        | 4.94E+00      | 1.77E-01      | 2.90E+01 | J             | GEO-33   | 1.18E+01      |                        |            |                          |
| Fluoranthene               | 206-44-0   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 1.30E-01              | J        | 5.01E+01      | 2.31E+00      | 3.55E+02 | J             | GEO-33   | 1.23E+02      |                        |            |                          |
| Fluorene                   | 86-73-7    | 8             | 4              | 50              | 3.80E-02      | 5.20E-02      | 3.30E-01              | J        | 8.10E+00      | 1.90E-01      | 6.30E+01 | J             | GEO-33   | 2.22E+01      |                        |            |                          |
| Indeno(1,2,3-cd)pyrene     | 193-39-5   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 2.60E-01              | J        | 5.49E+00      | 1.02E+00      | 3.10E+01 | J             | GEO-33   | 1.05E+01      |                        |            |                          |
| Naphthalene                | 91-20-3    | 8             | 5              | 62.5            | 3.80E-02      | 5.60E-01      | 7.50E-02              | J        | 1.63E+00      | 2.97E-01      | 6.89E+00 | J             | GEO-33   | 2.65E+00      |                        |            |                          |
| Phenanthrene               | 85-01-8    | 8             | 6              | 75              | 3.80E-02      | 3.90E-02      | 1.50E-01              | J        | 3.25F+01      | 7.56E-01      | 2.45E+02 | J             | GEO-33   | 8.59E+01      |                        |            |                          |
| Phenol                     | 108-95-2   | 6             | 4              | 66.67           | 7.70E-02      | 4.13E-01      | 1.40E-01              | J        | 1.99E-01      | 1.64E-01      | 3.80E-01 | J             | GEO-29   | 1.13E-01      |                        |            |                          |
| Pyrene                     | 129-00-0   | 8             | 7              | 87.5            | 3.80E-02      | 3.80E-02      | 2.50E-01              | J        | 3.78E+01      | 2.54E+00      | 2.60E+02 | J             | GEO-33   | 9.01E+01      |                        |            |                          |

**Table 14**  
**Statistical Summary and Selection of COPCs in EU5 Soil (#-I' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I                            |   | Is the Maximum<br>Detected > TRG?<br>UCL > TRG? | Is the 95%<br>UCL > TRG? |
|------------------------|------------------|---------------------------------|-----------------------------------|--|-----------------------------------|---|---|--------------------------|
|                        |                  |                                 |                                   |  | Unrestricted Soil<br>TRG<br>mg/kg | Is the Maximum<br>Detected > TRG?<br>UCL > TRG? |   |                          |
| <b>Semivolatiles</b>   |                  |                                 |                                   |  |                                   |   |   |                          |
| 2,4-dimethylphenol     | 1.76E-01         | 4.42E-01                        | Normal/Lognormal                  | 1.10E-01                                 | 1.56E+03                          | no  |   |                          |
| 2-methylnaphthalene    | 4.71E+00         | 4.63E+04                        | Lognormal                         | 9.20E+00                                 | 3.13E+03                          | no  |   |                          |
| 2-methylphenol         | 8.17E-02         | 1.82E-01                        | Normal/Lognormal                  | 4.20E-02                                 | 3.91E+03                          | no  |   |                          |
| 3- and 4-methylphenol  | 1.74E-01         | 4.55E-01                        | Normal/Lognormal                  | 1.40E-01                                 | 3.91E+02                          | no  |   |                          |
| Acenaphthene           | 1.95E+01         | 3.47E+04                        | Lognormal                         | 5.35E+01                                 | 4.69E+03                          | no  |   |                          |
| Acenaphthylene         | 6.27E+00         | 1.25E+03                        | Lognormal                         | 1.60E+01                                 | 4.69E+03                          | no  |   |                          |
| Anthracene             | 2.94E+01         | 1.81E+07                        | Lognormal                         | 7.97E+01                                 | 2.35E+04                          | no  |   |                          |
| Benz(a)anthracene      | 3.23E+01         | 9.31E+04                        | Lognormal                         | 8.35E+01                                 | 8.75E-01                          | YES - COPC                                      |   |                          |
| Benz(a)pyrene          | 2.08E+01         | 2.16E+04                        | Lognormal                         | 5.25E+01                                 | 8.75E-02                          | YES - COPC                                      |   |                          |
| Benz(b)fluoranthene    | 3.20E+01         | 8.21E+04                        | Lognormal                         | 7.95E+01                                 | 8.75E-01                          | YES - COPC                                      |   |                          |
| Benz(g,h)perylene      | 1.01E+01         | 2.82E+04                        | Lognormal                         | 2.55E+01                                 | 2.35E+03                          | no  |   |                          |
| Benz(k)fluoranthene    | 1.16E+01         | 3.72E+03                        | Lognormal                         | 2.85E+01                                 | 8.75E+00                          | YES   |   |                          |
| Carbazole              | 7.04E+00         | 8.43E+06                        | Lognormal                         | 1.35E+01                                 | 3.19E+01                          | no  |   |                          |
| Chrysene               | 3.25E+01         | 1.14E+05                        | Lognormal                         | 8.25E+01                                 | 8.75E+01                          | YES*  |   |                          |
| Dibenz(a,h)anthracene  | 3.02E+00         | 1.76E+02                        | Lognormal                         | 7.45E+00                                 | 8.75E-02                          | YES   |   |                          |
| Dibenzofuran           | 1.46E+01         | 4.28E+06                        | Lognormal                         | 2.90E+01                                 | 3.13E+02                          | no  |   |                          |
| Fluoranthene           | 1.33E+02         | 2.93E+07                        | Lognormal                         | 3.55E+02                                 | 3.13E+03                          | no  |   |                          |
| Fluorene               | 2.30E+01         | 1.21E+05                        | Lognormal                         | 6.30E+01                                 | 3.13E+03                          | no  |   |                          |
| Indeno[1,2,3-cd]pyrene | 1.25E+01         | 5.34E+03                        | Lognormal                         | 2.10E+01                                 | 8.75E-01                          | YES   |   |                          |
| Naphthalene            | 3.41E+00         | 1.19E+03                        | Lognormal                         | 6.85E+00                                 | 6.45E+02                          | no  |   |                          |
| Phenanthrene           | 9.00E+01         | 2.29E+07                        | Lognormal                         | 2.45E+02                                 | 2.35E+03                          | no  |   |                          |
| Phenol                 | 2.92E-01         | 7.33E-01                        | Normal/Lognormal                  | 3.80E-01                                 | 4.69E+04                          | no  |   |                          |
| Pyrene                 | 9.82E+01         | 4.13E+06                        | Lognormal                         | 2.60E+02                                 | 2.35E+03                          | no  |   |                          |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

**Table 15**  
**Statistical Summary and Selection of COPCs in EU5 Soil (0-6' bgs)**  
**Kerr McGee, Hatfieldburg, MS**

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Maximum mg/kg | Mean mg/kg | Detected | Maximum mg/kg | Detected | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|---------------|------------|----------|---------------|----------|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>       |            |               |                |                 |               |               |                       |          |               |               |            |          |               |          |                                   |                          |
| 2,4-dimethylphenol         | 105-67-9   | 18            | 1              | 5.56            | 7.50E-02      | 4.55E-01      | 1.10E-01              | J        | 6.17E-02      | 4.98E-02      | 1.10E-01   | J        | GEO-30        | 5.60E-02 |                                   |                          |
| 2-methylnaphthalene        | 91-57-6    | 18            | 5              | 27.78           | 3.80E-02      | 4.10E-02      | 5.10E-02              | J        | 6.21E-01      | 4.99E-02      | 9.20E+00   | J        | GEO-33        | 2.15E+00 |                                   |                          |
| 2-methylphenol             | 95-48-7    | 18            | 1              | 5.56            | 3.80E-02      | 2.09E-01      | 4.20E-02              | J        | 2.95E-02      | 2.44E-02      | 4.20E-02   | J        | GEO-30        | 2.57E-02 |                                   |                          |
| 3- and 4-methylphenol      | 106-44-5   | 18            | 1              | 5.56            | 7.50E-02      | 4.13E-01      | 1.40E-01              | J        | 6.22E-02      | 5.02E-02      | 1.40E-01   | J        | GEO-30        | 5.46E-02 |                                   |                          |
| Acenaphthene               | 83-32-9    | 24            | 4              | 16.67           | 2.90E-02      | 2.40E+00      | 1.10E-01              | J        | 2.53E+00      | 6.24E-02      | 5.35E+01   | J        | GEO-33        | 1.09E+01 |                                   |                          |
| Acenaphthylene             | 208-96-8   | 24            | 6              | 25              | 3.80E-02      | 2.40E+00      | 4.80E-02              | J        | 9.59E-01      | 7.93E-02      | 1.60E+01   | J        | GEO-33        | 3.26E+00 |                                   |                          |
| Anthracene                 | 120-12-7   | 24            | 6              | 25              | 5.40E-04      | 6.00E-02      | 1.30E-01              | J        | 3.94E+00      | 4.33E-02      | 7.97E+01   | J        | GEO-33        | 1.63E+01 |                                   |                          |
| Benz(a)anthracene          | 56-55-3    | 24            | 13             | 54.17           | 3.80E-02      | 4.10E-02      | 6.80E-03              | Z        | 4.81E+00      | 1.11E-01      | 8.35E+01   | Z        | GEO-33        | 1.71E+01 |                                   |                          |
| Benz(a)pyrene              | 50-32-8    | 24            | 13             | 54.17           | 3.80E-02      | 4.10E-02      | 8.30E-03              | Z        | 3.17E+00      | 1.15E-01      | 5.25E+01   | Z        | GEO-33        | 1.07E+01 |                                   |                          |
| Benz(b)fluoranthene        | 205-99-2   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 9.00E-03              | Z        | 4.99E+00      | 1.59E-01      | 7.95E+01   | Z        | GEO-33        | 1.63E+01 |                                   |                          |
| Benz(g,h,i)perylene        | 191-24-2   | 24            | 12             | 50              | 3.80E-02      | 4.10E-02      | 6.70E-03              | J        | 1.55E+00      | 7.48E-02      | 2.55E+01   | J        | GEO-33        | 5.22E+00 |                                   |                          |
| Benz(k)fluoranthene        | 207-08-9   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 4.70E-03              | Z        | 1.84E+00      | 9.91E-02      | 2.85E+01   | Z        | GEO-33        | 5.86E+00 |                                   |                          |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 18            | 2              | 11.11           | 7.50E-02      | 4.13E-01      | 1.40E-01              | J        | 6.84E-02      | 5.41E-02      | 1.50E-01   | J        | GEO-32        | 5.80E-02 |                                   |                          |
| Carbazole                  | 86-74-8    | 18            | 4              | 22.22           | 3.80E-02      | 4.10E-02      | 5.30E-01              | J        | 9.78E-01      | 5.52E-02      | 1.35E+01   | J        | GEO-33        | 3.17E+00 |                                   |                          |
| Chrysene                   | 218-01-9   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 2.40E-03              | J        | 4.91E+00      | 1.16E-01      | 8.25E+01   | J        | GEO-33        | 1.69E+01 |                                   |                          |
| Dibenz(a,h)anthracene      | 53-70-3    | 24            | 12             | 50              | 3.80E-02      | 4.10E-02      | 1.70E-03              | J        | 4.89E-01      | 4.58E-02      | 7.45E+00   | J        | GEO-33        | 1.53E+00 |                                   |                          |
| Dibenzofuran               | 132-64-9   | 18            | 6              | 33.33           | 3.80E-02      | 4.10E-02      | 3.90E-02              | J        | 1.82E+00      | 5.87E-02      | 2.90E+01   | J        | GEO-33        | 6.81E+00 |                                   |                          |
| Fluoranthene               | 206-44-0   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 1.30E-02              | Z        | 1.85E+01      | 1.82E-01      | 3.55E+02   | Z        | GEO-33        | 7.23E+01 |                                   |                          |
| Fluorene                   | 86-73-7    | 24            | 8              | 33.33           | 2.90E-03      | 5.20E-02      | 3.60E-03              | J        | 2.98E+00      | 4.72E-02      | 6.30E+01   | J        | GEO-33        | 1.28E+01 |                                   |                          |
| Indeno(1,2,3-cd)pyrene     | 193-39-5   | 24            | 13             | 54.17           | 3.80E-02      | 4.10E-02      | 7.80E-03              | J        | 1.97E+00      | 9.68E-02      | 3.10E+01   | J        | GEO-33        | 6.37E+00 |                                   |                          |
| Naphthalene                | 91-20-3    | 24            | 6              | 25              | 2.90E-02      | 5.60E-01      | 7.50E-02              | J        | 5.93E-01      | 6.22E-02      | 6.83E+00   | J        | GEO-33        | 1.65E+00 |                                   |                          |
| Phenanthrene               | 85-01-8    | 24            | 13             | 54.17           | 3.80E-02      | 4.10E-02      | 6.80E-03              | J        | 1.21E+01      | 1.17E-01      | 2.45E+02   | J        | GEO-33        | 5.00E+01 |                                   |                          |
| Phenol                     | 108-95-2   | 18            | 13             | 72.22           | 7.50E-02      | 4.13E-01      | 1.00E-01              | J        | 1.48E-01      | 1.22E-01      | 3.80E-01   | J        | GEO-29        | 8.68E-02 |                                   |                          |
| Pyrene                     | 129-00-0   | 24            | 14             | 58.33           | 3.80E-02      | 4.10E-02      | 1.60E-02              | J        | 1.39E+01      | 1.87E-01      | 2.60E+02   | J        | GEO-33        | 5.30E+01 |                                   |                          |

**Table 15**  
**Statistical Summary and Selection of COPCs in EU5 Soil (0-6' bgs)**  
**Kerr McGee, Hanniesburg, MS**

| Constituent                | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|----------------------------|------------------|---------------------------------|-----------------------------------|--|--|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>       |                  |                                 |                                   |  |  |                                   |                          |
| 2,4-dimethylphenol         | 8.47E-02         | 7.85E-02                        | Unknown                           | 7.85E-02                                 | 4.08E+04                               | no                                |                          |
| 2-methylnaphthalene        | 1.50E+00         | 1.43E+00                        | Unknown                           | 1.43E+00                                 | 8.18E+04                               | no                                |                          |
| 2-methylphenol             | 4.01E-02         | 3.69E-02                        | Unknown                           | 3.69E-02                                 | 1.02E+05                               | no                                |                          |
| 3- and 4-methylphenol      | 8.46E-02         | 7.99E-02                        | Unknown                           | 7.99E-02                                 | 1.02E+04                               | no                                |                          |
| Acenaphthene               | 6.35E+00         | 3.21E+00                        | Unknown                           | 3.21E+00                                 | 1.23E+05                               | no                                |                          |
| Acenaphthylene             | 2.10E+00         | 2.70E+00                        | Unknown                           | 2.70E+00                                 | 1.23E+05                               | no                                |                          |
| Anthracene                 | 9.62E+00         | 6.15E+01                        | Unknown                           | 6.15E+01                                 | 6.13E+05                               | no                                |                          |
| Benz(a)anthracene          | 1.08E+01         | 7.77E+01                        | Unknown                           | 7.77E+01                                 | 7.84E+00                               | YES                               | YES - COPC               |
| Benz(a)pyrene              | 6.93E+00         | 4.10E+01                        | Unknown                           | 4.10E+01                                 | 7.84E-01                               | YES                               | YES - COPC               |
| Benz(b)fluoranthene        | 1.07E+01         | 1.30E+02                        | Unknown                           | 7.95E+01                                 | 7.84E+00                               | YES                               | YES - COPC               |
| Benz(ghi)perylene          | 3.38E+00         | 8.53E+00                        | Unknown                           | 8.53E+00                                 | 6.13E+04                               | no                                |                          |
| Benz(k)fluoranthene        | 3.89E+00         | 1.97E+01                        | Unknown                           | 1.97E+01                                 | 7.84E+01                               | no                                | YES*                     |
| Bis(2-ethylhexyl)phthalate | 9.21E-02         | 9.19E-02                        | Unknown                           | 9.19E-02                                 | 4.09E+02                               | no                                |                          |
| Carbazole                  | 2.28E+00         | 4.56E+00                        | Unknown                           | 4.56E+00                                 | 2.86E+02                               | no                                |                          |
| Chrysene                   | 1.08E+01         | 1.27E+02                        | Unknown                           | 8.25E+01                                 | 7.84E+02                               | no                                | YES*                     |
| Dibenz(a,h)anthracene      | 1.02E+00         | 2.04E+00                        | Unknown                           | 2.04E+00                                 | 7.84E-01                               | YES                               | YES - COPC               |
| Dibenzofuran               | 4.61E+00         | 4.75E+00                        | Unknown                           | 4.75E+00                                 | 8.18E+03                               | no                                |                          |
| Fluoranthene               | 4.38E+01         | 7.28E+02                        | Unknown                           | 3.55E+02                                 | 8.17E+04                               | no                                |                          |
| Fluorene                   | 7.47E+00         | 7.47E+00                        | Unknown                           | 7.47E+00                                 | 8.17E+04                               | no                                |                          |
| Indeno(1,2,3-cd)pyrene     | 4.20E+00         | 1.71E+01                        | Unknown                           | 1.71E+01                                 | 7.84E+00                               | YES                               | YES - COPC               |
| Naphthalene                | 1.17E+00         | 1.53E+00                        | Unknown                           | 1.53E+00                                 | 8.24E+02                               | no                                |                          |
| Phenanthrene               | 2.96E+01         | 1.26E+02                        | Unknown                           | 1.26E+02                                 | 6.13E+04                               | no                                |                          |
| Phenol                     | 1.84E-01         | 2.27E-01                        | Normal/Lognormal                  | 2.27E-01                                 | 1.23E+05                               | no                                |                          |
| Pyrene                     | 3.24E+01         | 4.64E+02                        | Unknown                           | 2.60E+02                                 | 6.13E+04                               | no                                |                          |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

**Table 16**  
**Statistical Summary and Selection of COPCs in EU5 Soil (0-20' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg   | Maximum mg/kg   | Logarithmic Mean mg/kg | Maximum mg/kg | Maximum mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|-----------------|-----------------|------------------------|---------------|---------------|-----------------------------------|--------------------------|
|                            |            |               |                |                 | Detection Limit | Detection Limit | Mean mg/kg             | Detected      | Detected      |                                   |                          |
|                            |            |               |                |                 | mg/kg           | mg/kg           | mg/kg                  | mg/kg         | mg/kg         |                                   |                          |
| <b>Semivolatiles</b>       |            |               |                |                 |                 |                 |                        |               |               |                                   |                          |
| 2,4-dimethylphenol         | 105-67-9   | 21            | 1              | 4.76            | 6.70E-02        | 4.55E-01        | 1.10E-01               | J             | 1.10E-01      | J                                 | 5.26E-02                 |
| 2-methylnaphthalene        | 91-57-6    | 21            | 6              | 28.57           | 3.30E-02        | 4.10E-02        | 5.10E-02               | J             | 1.25E+00      | SB-07                             | 3.73E+00                 |
| 2-methylphenol             | 95-48-7    | 21            | 1              | 4.76            | 3.80E-02        | 2.09E-01        | 4.20E-02               | J             | 3.01E-02      | 2.55E-02                          | 2.38E-02                 |
| 3- and 4-methylphenol      | 106-44-5   | 21            | 1              | 4.76            | 7.50E-02        | 4.13E-01        | 1.40E-01               | J             | 6.05E-02      | 5.01E-02                          | 5.05E-02                 |
| Acenaphthene               | 83-32-9    | 28            | 5              | 17.86           | 2.90E-02        | 2.40E+00        | 1.10E-01               | J             | 2.52E+00      | 7.17E-02                          | 1.02E+01                 |
| Acenaphthylene             | 208-96-8   | 28            | 7              | 25              | 3.30E-02        | 2.40E+00        | 4.80E-02               | J             | 8.45E-01      | 1.60E+01                          | GEO-33                   |
| Anthracene                 | 120-12-7   | 28            | 7              | 25              | 5.40E-04        | 6.00E-02        | 1.30E-01               | J             | 3.55E+00      | 4.43E-02                          | 7.97E+01                 |
| Benz(a)anthracene          | 56-55-3    | 28            | 15             | 53.57           | 3.30E-02        | 4.10E-02        | 6.80E-03               | Z             | 4.18E+00      | 1.08E-01                          | 8.35E+01                 |
| Benz(a)pyrene              | 50-32-8    | 28            | 15             | 53.57           | 3.80E-02        | 6.70E-02        | 8.30E-03               | Z             | 2.76E+00      | 1.15E-01                          | 5.25E+01                 |
| Benz(b)fluoranthene        | 205-99-2   | 28            | 16             | 57.14           | 3.80E-02        | 6.70E-02        | 9.00E-03               | Z             | 4.32E+00      | 1.55E-01                          | 9.96E+00                 |
| Benz(ghi)perylene          | 191-24-2   | 28            | 14             | 50              | 3.80E-02        | 6.70E-02        | 6.70E-03               | J             | 1.35E+00      | 2.55E-01                          | GEO-33                   |
| Benz(k)fluoranthene        | 207-08-9   | 28            | 16             | 57.14           | 3.80E-02        | 1.30E-01        | 4.70E-03               | Z             | 1.60E+00      | 1.02E-01                          | GEO-33                   |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 21            | 2              | 9.52            | 6.70E-02        | 4.13E-01        | 1.40E-01               | J             | 6.41E-02      | 5.14E-02                          | 1                        |
| Carbazole                  | 86-74-8    | 21            | 5              | 23.81           | 3.30E-02        | 4.10E-02        | 5.30E-01               | J             | 8.86E-01      | 1.35E+01                          | GEO-33                   |
| Chrysene                   | 218-01-9   | 28            | 16             | 57.14           | 3.30E-02        | 4.10E-02        | 2.40E-03               | J             | 4.27E+00      | 1.11E-01                          | 8.25E+01                 |
| Dibenz(a,h)anthracene      | 53-70-3    | 28            | 14             | 50              | 3.80E-02        | 6.70E-02        | 1.70E-03               | J             | 4.27E-01      | 4.60E-02                          | 7.45E+00                 |
| Dibenzofuran               | 132-64-9   | 21            | 7              | 33.33           | 3.30E-02        | 4.10E-02        | 3.90E-02               | J             | 2.02E+00      | 6.63E-02                          | GEO-32                   |
| Fluoranthene               | 206-44-0   | 28            | 16             | 57.14           | 3.30E-02        | 4.10E-02        | 1.30E-02               | Z             | 1.62E+01      | 1.79E-01                          | 5.46E-02                 |
| Fluorene                   | 86-73-7    | 28            | 9              | 32.14           | 2.90E-03        | 5.20E-02        | 3.60E-03               | J             | 2.86E+00      | 5.16E-02                          | GEO-33                   |
| Indeno[1,2,3-cd]pyrene     | 193-39-5   | 28            | 15             | 53.57           | 3.80E-02        | 6.70E-02        | 7.80E-03               | J             | 1.71E+00      | 9.65E-02                          | 3.10E+01                 |
| Naphthalene                | 91-20-3    | 28            | 7              | 25              | 2.90E-02        | 5.60E-01        | 7.50E-02               | J             | 1.81E+00      | 7.49E-02                          | 3.60E+01                 |
| Phenanthrene               | 85-01-8    | 28            | 15             | 53.57           | 3.30E-02        | 4.10E-02        | 6.80E-03               | J             | 1.11E+01      | 1.23E-01                          | 2.45E+02                 |
| Phenol                     | 108-95-2   | 21            | 13             | 61.9            | 3.30E-02        | 4.13E-01        | 1.00E-01               | J             | 1.29E-01      | 9.16E-02                          | GEO-29                   |
| Pyrene                     | 129-00-0   | 28            | 16             | 57.14           | 3.80E-02        | 6.70E-02        | 1.60E-02               | J             | 1.21E+01      | 1.91E-01                          | 2.60E+02                 |
| <b>Volatiles</b>           |            |               |                |                 |                 |                 |                        |               |               |                                   |                          |
| Acetone                    | 67-64-1    | 3             | 3              | 100             | 0.00E+00        | 0.00E+00        | 1.00E-01               | J             | 3.70E-02      | 2.92E-02                          | 5.90E-02                 |
| Benzene                    | 71-43-2    | 3             | 1              | 33.33           | 1.00E-03        | 1.00E-03        | 5.00E-03               | J             | 2.00E-03      | 1.08E-03                          | SB-07                    |
| Ethybenzene                | 100-41-4   | 3             | 1              | 33.33           | 1.00E-03        | 1.00E-03        | 2.40E-02               |               | 8.33E-03      | 1.82E-03                          | 2.40E-02                 |
| Toluene                    | 108-88-3   | 3             | 1              | 33.33           | 1.00E-03        | 1.00E-03        | 1.50E-02               |               | 5.33E-03      | 1.55E-03                          | 1.50E-02                 |
| Xylene (total)             | 1330-20-7  | 3             | 1              | 33.33           | 1.00E-03        | 1.00E-03        | 7.50E-02               |               | 2.53E-02      | 2.66E-03                          | 7.50E-02                 |



**Table 16**  
**Statistical Summary and Selection of COPCs in EU5 Soil (0-20' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? |
|----------------------------|------------------|---------------------------------|-----------------------------------|--|--|-----------------------------------|
| <b>Semi-volatiles</b>      |                  |                                 |                                   |  |  |                                   |
| 2,4-dimethylphenol         | 7.75E-02         | 7.01E-02                        | Unknown                           | 7.01E-02                                 | 4.08E+04                               | no                                |
| 2-methylnaphthalene        | 2.65E+00         | 4.34E+00                        | Unknown                           | 4.34E+00                                 | 8.18E+04                               | no                                |
| 2-methylphenol             | 3.91E-02         | 3.65E-02                        | Unknown                           | 3.65E-02                                 | 1.02E+05                               | no                                |
| 3- and 4-methylphenol      | 7.95E-02         | 7.38E-02                        | Unknown                           | 7.38E-02                                 | 1.02E+04                               | no                                |
| Acenaphthene               | 5.80E+00         | 4.62E+00                        | Unknown                           | 4.62E+00                                 | 1.23E+05                               | no                                |
| Acenaphthylene             | 1.82E+00         | 1.80E+00                        | Unknown                           | 1.80E+00                                 | 1.23E+05                               | no                                |
| Anthracene                 | 8.39E+00         | 3.95E+01                        | Unknown                           | 3.95E+01                                 | 6.13E+05                               | no                                |
| Benz(a)anthracene          | 9.27E+00         | 3.39E+01                        | Unknown                           | 3.39E+01                                 | 7.84E+00                               | YES                               |
| Benz(a)pyrene              | 5.96E+00         | 1.71E+01                        | Unknown                           | 1.71E+01                                 | 7.84E-01                               | YES                               |
| Benz(b)fluoranthene        | 9.21E+00         | 4.67E+01                        | Unknown                           | 4.67E+01                                 | 7.84E+00                               | YES                               |
| Benz(ghi)perylene          | 2.91E+00         | 4.10E+00                        | Unknown                           | 4.10E+00                                 | 6.13E+04                               | no                                |
| Benz(k)fluoranthene        | 3.35E+00         | 8.05E+00                        | Unknown                           | 8.05E+00                                 | 7.84E+01                               | no                                |
| Bis(2-ethylhexyl)phthalate | 8.47E-02         | 8.14E-02                        | Unknown                           | 8.14E-02                                 | 4.09E+02                               | no                                |
| Carbazole                  | 1.99E+00         | 3.30E+00                        | Unknown                           | 3.30E+00                                 | 2.86E+02                               | no                                |
| Chrysene                   | 9.32E+00         | 5.02E+01                        | Unknown                           | 5.02E+01                                 | 7.84E+02                               | *YES                              |
| Dibenz(a,h)anthracene      | 8.84E-01         | 1.04E+00                        | Unknown                           | 1.04E+00                                 | 7.84E-01                               | YES                               |
| Dibenzofuran               | 4.49E+00         | 8.71E+00                        | Unknown                           | 8.71E+00                                 | 8.18E+03                               | no                                |
| Fluoranthene               | 3.77E+01         | 3.46E+02                        | Unknown                           | 3.46E+02                                 | 8.17E+04                               | no                                |
| Fluorene                   | 6.70E+00         | 7.93E+00                        | Unknown                           | 7.93E+00                                 | 8.17E+04                               | no                                |
| Indeno(1,2,3-cd)pyrene     | 3.61E+00         | 7.55E+00                        | Unknown                           | 7.55E+00                                 | 7.84E+00                               | YES                               |
| Naphthalene                | 4.02E+00         | 4.18E+00                        | Unknown                           | 4.18E+00                                 | 8.24E+02                               | no                                |
| Phenanthrene               | 2.60E+01         | 1.03E+02                        | Unknown                           | 1.03E+02                                 | 6.13E+04                               | no                                |
| Phenol                     | 1.64E-01         | 2.50E-01                        | Normal/Lognormal                  | 2.50E-01                                 | 1.23E+05                               | no                                |
| Pyrene                     | 2.79E+01         | 1.80E+02                        | Unknown                           | 1.80E+02                                 | 6.13E+04                               | no                                |
| <b>Volatiles</b>           |                  |                                 |                                   |  |  |                                   |
| Acetone                    | 7.89E-02         | 1.64E+02                        | Normal/Lognormal                  | 5.90E-02                                 | 1.04E+05                               | no                                |
| Benzene                    | 6.38E-03         | 3.21E+04                        | Unknown                           | 5.00E-03                                 | 1.36E+00                               | no                                |
| Ethylbenzene               | 3.12E-02         | 2.50E+18                        | Unknown                           | 2.40E-02                                 | 3.95E+02                               | no                                |
| Toluene                    | 1.94E-02         | 3.20E+13                        | Unknown                           | 1.50E-02                                 | 3.80E+01                               | no                                |
| Xylene (total)             | 9.78E-02         | 7.06E+32                        | Unknown                           | 7.50E-02                                 | 3.18E+02                               | no                                |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

**Table 17**  
**Statistical Summary and Selection of COPCs in EU6 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                  | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Minimum Detected mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected Qualifier | Detected Concentration mg/kg | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|------------------------|------------|------------------------|---------------|--------------------|------------------------------|-----------------------------|-----------------------------------|--------------------------|
| Semivolatiles                |            |               |                |                 |               |               |                       |                        |            |                        |               |                    |                              |                             |                                   |                          |
| 1,2,4-trichlorobenzene       | 120-82-1   | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 1,2-dichlorobenzene          | 95-50-1    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 1,3-dichlorobenzene          | 541-73-1   | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 1,4-dichlorobenzene          | 106-46-7   | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 2,2'-oxypy (1-chloropropane) | 108-60-1   | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 2,4,5-trichlorophenol        | 95-95-4    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 2,4,6-trichlorophenol        | 88-06-2    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 2,4-dichlorophenol           | 120-83-2   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 2,4-dimethylphenol           | 105-67-9   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 2,4-dinitrophenol            | 51-28-5    | 3             | 1              | 33.33           | 2.40E-01      | 2.50E-01      | 2.30E+00              | 8.48E-01               | 3.26E-01   | 2.30E+00               | SD-04         | 1.26E+00           |                              |                             |                                   |                          |
| 2,4-dinitrotoluene           | 121-14-2   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 4.00E-01              | 1.62E-01               | 8.94E-02   | 4.00E-01               | SD-04         | 2.07E-01           |                              |                             |                                   |                          |
| 2,6-dinitrotoluene           | 606-20-2   | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 2-chloronaphthalene          | 91-58-7    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 2-chlorophenol               | 95-57-8    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 2-methylnaphthalene          | 91-57-6    | 3             | 3              | 100             | 0.00E+00      | 9.00E+00      | 9.10E-02              | J                      | 1.28E+01   | 1.15E+00               | 3.80E+01      | SD-04              | 2.18E+01                     |                             |                                   |                          |
| 2-methylphenol               | 95-48-7    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 2-nitroaniline               | 88-74-4    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 2-nitrophenol                | 88-75-5    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 3- and 4-methylphenol        | 106-44-5   | 3             | 3              | 100             | 0.00E+00      | 9.00E+00      | 9.30E-02              | J                      | 3.34E-01   | 2.02E-01               | 8.00E-01      | SD-04              | 4.03E-01                     |                             |                                   |                          |
| 3,3'-dichlorobenzidine       | 91-94-1    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 3-nitroaniline               | 99-09-2    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 4,6-dinitro-2-methylphenol   | 534-52-1   | 3             | 1              | 33.33           | 2.10E-01      | 2.10E-01      | 2.30E+00              | 7.37E-01               | 2.80E-01   | 2.00E+00               | SD-04         | 1.09E+00           |                              |                             |                                   |                          |
| 4-bromophenylphenylether     | 101-55-3   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 4-chloro-3-methylphenol      | 59-50-7    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 4-chloroaniline              | 106-47-8   | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 4-chlorophenylphenylether    | 7005-72-3  | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E-01           |                              |                             |                                   |                          |
| 4-nitroaniline               | 100-01-6   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| 4-nitrophenol                | 100-02-7   | 3             | 1              | 33.33           | 2.10E-01      | 2.10E-01      | 2.00E+00              | 7.37E-01               | 2.80E-01   | 2.00E+00               | J             | SD-04              | 1.09E+00                     |                             |                                   |                          |
| Acenaphthene                 | 83-32-9    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 1.00E-01              | J                      | 4.70E+01   | 2.32E+00               | 1.40E+02      | J                  | SD-04                        | 8.03E+01                    |                                   |                          |
| Acenaphthylene               | 208-96-8   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 1.70E-01              | J                      | 5.29E+00   | 2.17E+00               | 8.90E+00      | J                  | SD-03                        | 4.56E+00                    |                                   |                          |
| Anthracene                   | 120-12-7   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 8.80E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E-01           |                              |                             |                                   |                          |
| Benz(a)anthracene            | 56-55-3    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 9.30E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 5.14E+01           |                              |                             |                                   |                          |
| Benz(a)pyrene                | 50-32-8    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 9.70E-01              | J                      | 2.77E+01   | 1.16E+01               | 4.90E+01      | J                  | SD-03                        | 2.45E+01                    |                                   |                          |
| Benz(b)fluoranthene          | 205-99-2   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 1.40E+00              | 4.18E+01               | 1.71E+01   | 7.80E+00               | SD-03         | 3.85E+01           |                              |                             |                                   |                          |
| Benz(ghi)perylene            | 191-24-2   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 4.20E-01              | 1.40E+01               | 5.04E+00   | 3.20E+00               | SD-03         | 2.31E+00           |                              |                             |                                   |                          |
| Benz(k)fluoranthene          | 207-08-9   | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 5.00E-01              | 4.26E+01               | 1.36E+01   | 5.92E+00               | SD-03         | 1.65E+01           |                              |                             |                                   |                          |
| Bis(2-chloroethoxy)methane   | 111-91-1   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-03         | 1.18E+01           |                              |                             |                                   |                          |
| Bis(2-chloroethyl)ether      | 111-44-4   | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              | 1.48E-01               | 5.65E-02   | 4.00E-01               | SD-04         | 2.19E+01           |                              |                             |                                   |                          |
| Bis(2-ethylhexyl)phthalate   | 117-81-7   | 3             | 1              | 33.33           | 1.50E-01      | 2.50E-01      | 8.80E-01              | 3.60E-01               | 2.02E-01   | 8.80E-01               | SD-04         | 4.51E+01           |                              |                             |                                   |                          |
| Butylbenzylphthalate         | 85-68-7    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              | 2.95E-01               | 1.13E-01   | 8.00E-01               | SD-04         | 4.37E+01           |                              |                             |                                   |                          |
| Carbazole                    | 86-74-8    | 3             | 3              | 100             | 0.00E+00      | 0.00E+00      | 2.20E-01              | J                      | 3.37E+01   | 2.77E+00               | 1.00E+02      | J                  | SD-04                        | 5.74E+01                    |                                   |                          |

**Table 17**  
**Statistical Summary and Selection of COPCs in EU6 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                   | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Unrestricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|-------------------------------|------------------|---------------------------------|--------------------------------|--|--|-----------------------------------|--------------------------|
| <b>Semi挥发性</b>                |                  |                                 |                                |  |  |                                   |                          |
| 1,2,4-trichlorobenzene        | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 5.27E+02                                 | no                                |                          |
| 1,2-dichlorobenzene           | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 2.79E+02                                 | no                                |                          |
| 1,3-dichlorobenzene           | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 2.35E+03                                 | no                                |                          |
| 1,4-dichlorobenzene           | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 2.66E+01                                 | no                                |                          |
| 2,2'-oxybis (1-chloropropane) | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 5.93E+00                                 | no                                |                          |
| 2,4,5-trichlorophenol         | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 7.82E+03                                 | no                                |                          |
| 2,4,6-trichlorophenol         | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 5.81E+01                                 | no                                |                          |
| 2,4-dichlorophenol            | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 2.35E+02                                 | no                                |                          |
| 2,4-dimethylphenol            | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 1.56E+03                                 | no                                |                          |
| 2,4-dinitrophenol             | 2.97E+00         | 4.39E+11                        | Lognormal                      | 2.30E+00                                 | 1.56E+02                                 | no                                |                          |
| 2,4-dinitrotoluene            | 5.10E-01         | 1.19E+06                        | Lognormal                      | 4.00E-01                                 | 1.56E+02                                 | no                                |                          |
| 2,6-dinitrotoluene            | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 7.82E+01                                 | no                                |                          |
| 2-chloronaphthalene           | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 6.26E+03                                 | no                                |                          |
| 2-chlorophenol                | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 3.91E+02                                 | no                                |                          |
| 2-methylnaphthalene           | 4.96E-01         | 3.37E+41                        | Normal/Lognormal               | 3.80E+01                                 | 3.13E+03                                 | no                                |                          |
| 2-methylphenol                | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 3.91E+03                                 | no                                |                          |
| 2-nitroaniline*               | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | NA                                       | no                                |                          |
| 2-nitrophenol*                | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | NA                                       | no                                |                          |
| 3- and 4-methylphenol         | 1.01E+00         | 2.30E+05                        | Normal/Lognormal               | 8.00E-01                                 | 3.91E+02                                 | no                                |                          |
| 3,3'-dichlorobenzidine        | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 1.42E+00                                 | no                                |                          |
| 3-nitroaniline*               | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | NA                                       | no                                |                          |
| 4,6-dinitro-2-methylphenol    | 2.58E+00         | 4.94E+11                        | Unknown                        | 2.00E+00                                 | 7.82E+00                                 | no                                |                          |
| 4,6-dinitrophenylphenylether* | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | NA                                       | no                                |                          |
| 4-chloro-3-methylphenol*      | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | NA                                       | no                                |                          |
| 4-chloroaniline               | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 3.13E+02                                 | no                                |                          |
| 4-chlorophenylphenylether*    | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | NA                                       | no                                |                          |
| 4-nitroaniline*               | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | NA                                       | no                                |                          |
| 4-nitrophenol                 | 2.58E+00         | 4.94E+11                        | Unknown                        | 2.00E+00                                 | 6.26E+02                                 | no                                |                          |
| Acenaphthene                  | 1.83E-02         | 6.49E+38                        | Normal/Lognormal               | 1.40E+02                                 | 4.69E+03                                 | no                                |                          |
| Acenaphthylene                | 1.30E+01         | 1.08E+21                        | Normal/Lognormal               | 8.90E+00                                 | 4.69E+03                                 | no                                |                          |
| Anthracene                    | 7.12E-00         | 1.51E+04                        | Normal/Lognormal               | 5.50E+00                                 | 2.35E+04                                 | no                                |                          |
| Benz(a)anthracene             | 1.29E-02         | 6.04E+25                        | Normal/Lognormal               | 1.00E+02                                 | 8.75E-01                                 | YES                               | YES - COPC               |
| Benz(a)pyrene                 | 6.89E-01         | 6.28E+20                        | Normal/Lognormal               | 4.90E+01                                 | 8.75E-02                                 | YES                               | YES - COPC               |
| Benz(b)fluoranthene           | 1.07E-02         | 2.69E+21                        | Normal/Lognormal               | 7.80E+01                                 | 8.75E-01                                 | YES                               | YES - COPC               |
| Benz(ghi)perylene             | 4.14E-01         | 7.01E+21                        | Normal/Lognormal               | 3.20E+01                                 | 2.35E+03                                 | no                                |                          |
| Benz(k)fluoranthene           | 3.38E-01         | 1.62E+20                        | Normal/Lognormal               | 2.30E+01                                 | 8.75E+00                                 | YES                               | YES - COPC               |
| Bis(2-chloroethoxy)methane*   | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | NA                                       | no                                |                          |
| Bis(2-chloroethyl)ether       | 5.16E-01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 2.73E-01                                 | YES                               | YES - COPC               |
| Bis(2-ethylhexyl)phthalate    | 1.12E-00         | 2.81E+06                        | Normal/Lognormal               | 8.80E-01                                 | 4.56E+01                                 | no                                |                          |
| Butylbenzylphthalate          | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 9.28E+02                                 | no                                |                          |
| Carbazole                     | 1.30E-02         | 3.75E+43                        | Normal/Lognormal               | 1.00E+02                                 | 3.19E+01                                 | YES                               | YES - COPC               |

**Table 17**  
**Statistical Summary and Selection of COPCs in EU6 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent               | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Maximum mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Qualifier mg/kg | Detected | Maximum Qualifier mg/kg | Location | Standard Deviation mg/kg |
|---------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|---------------|------------|------------------------|-------------------------|----------|-------------------------|----------|--------------------------|
|                           |            |               |                |                 |               |               |                       |          |               |               |            |                        |                         |          |                         |          |                          |
| Chrysene                  | 218-01-9   | 2             | 3              | 100             | 0.00E+00      | 1.30E+00      | 1.30E+00              | J        | 3.98E+01      | 1.61E+01      | 7.60E+01   | J                      | SD-03                   | 3.74E+01 |                         |          |                          |
| Dibenz(a)anthracene       | 53-70-3    | 3             | 3              | 100             | 0.00E+00      | 1.50E+01      | 1.50E+01              | J        | 4.35E+00      | 1.68E+00      | 9.60E+00   | J                      | SD-04                   | 4.81E+00 |                         |          |                          |
| Dibenzofuran              | 132-64-9   | 3             | 3              | 100             | 0.00E+00      | 1.00E+01      | 1.00E+01              | J        | 5.02E+01      | 1.93E+00      | 1.50E+02   | J                      | SD-04                   | 8.64E+01 |                         |          |                          |
| Diethylphthalate          | 84-56-2    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              |          | 2.95E-01      | 1.13E-01      | 8.00E-01   |                        | SD-04                   | 4.37E+01 |                         |          |                          |
| Dimethylphthalate         | 131-11-3   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              |          | 2.95E-01      | 1.13E-01      | 8.00E-01   |                        | SD-04                   | 4.37E+01 |                         |          |                          |
| Di-n-butylphthalate       | 84-74-2    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              |          | 2.95E-01      | 1.13E-01      | 8.00E-01   |                        | SD-04                   | 4.37E+01 |                         |          |                          |
| Di-n-octylphthalate       | 117-84-0   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              |          | 2.95E-01      | 1.13E-01      | 8.00E-01   |                        | SD-04                   | 4.37E+01 |                         |          |                          |
| Fluoranthene              | 206-44-0   | 3             | 100            | 0.00E+00        | 2.00E+00      | 2.00E+00      | 1.80E-01              | J        | 1.64E+02      | 2.70E+01      | 4.70E+02   |                        | SD-04                   | 2.65E+02 |                         |          |                          |
| Fluorene                  | 86-73-7    | 3             | 3              | 100             | 0.00E+00      | 4.20E-02      | 4.30E-02              | 4.00E-01 |               | 8.71E+01      | 3.60E+00   | 2.60E+02               |                         | SD-04    | 1.50E+02                |          |                          |
| Hexachlorobenzene         | 118-74-1   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              |          | 1.48E-01      | 5.65E-02      | 4.00E-01   |                        | SD-04                   | 2.19E+01 |                         |          |                          |
| Hexachlorobutadiene       | 87-66-3    | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              |          | 2.95E-01      | 1.13E-01      | 8.00E-01   |                        | SD-04                   | 4.37E+01 |                         |          |                          |
| Hexachlorocyclopentadiene | 77-47-4    | 3             | 1              | 33.33           | 2.10E-01      | 2.10E-01      | 2.00E+00              |          | 7.37E-01      | 2.80E-01      | 2.00E+00   |                        | SD-04                   | 1.09E+00 |                         |          |                          |
| Hexachloroethane          | 67-72-1    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              |          | 1.48E-01      | 5.65E-02      | 4.00E-01   |                        | SD-04                   | 2.19E+01 |                         |          |                          |
| Indeno(1,2,3-cd)pyrene    | 193-39-5   | 3             | 3              | 100             | 0.00E+00      | 5.40E-01      | 5.40E-01              |          | 1.72E+01      | 6.32E+00      | 3.90E+01   |                        | SD-03                   | 1.97E+01 |                         |          |                          |
| Isophorone                | 78-59-1    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              |          | 1.48E-01      | 5.65E-02      | 4.00E-01   |                        | SD-04                   | 2.19E+01 |                         |          |                          |
| Naphthalene               | 91-20-3    | 3             | 100            | 0.00E+00        | 1.60E-01      | 1.60E-01      | 1.00E+00              | J        | 5.25E+00      | 1.53E+00      | 1.40E+01   | J                      | SD-04                   | 7.61E+00 |                         |          |                          |
| Nitrobenzene              | 98-95-3    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              |          | 1.48E-01      | 5.65E-02      | 4.00E-01   |                        | SD-04                   | 2.19E+01 |                         |          |                          |
| N-nitrosodi-n-propylamine | 621-64-7   | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              |          | 1.48E-01      | 5.65E-02      | 4.00E-01   |                        | SD-04                   | 2.19E+01 |                         |          |                          |
| N-nitrosodiphenylamine    | 86-30-6    | 3             | 1              | 33.33           | 4.20E-02      | 4.30E-02      | 4.00E-01              |          | 1.48E-01      | 5.65E-02      | 4.00E-01   |                        | SD-04                   | 2.19E+01 |                         |          |                          |
| Pentachlorophenol         | 87-86-5    | 3             | 1              | 33.33           | 2.10E-01      | 2.10E-01      | 2.00E+00              |          | 7.37E-01      | 2.80E-01      | 2.00E+00   |                        | SD-04                   | 1.09E+00 |                         |          |                          |
| Phenanthrene              | 85-01-8    | 3             | 3              | 100             | 0.00E+00      | 6.60E-01      | 6.60E-01              |          | 2.91E+02      | 1.27E+01      | 8.70E+02   |                        | SD-04                   | 5.01E+02 |                         |          |                          |
| Phenol                    | 108-95-2   | 3             | 1              | 33.33           | 8.40E-02      | 8.50E-02      | 8.00E-01              |          | 2.95E-01      | 1.13E-01      | 8.00E-01   |                        | SD-04                   | 4.37E+01 |                         |          |                          |
| Pyrene                    | 129-00-0   | 3             | 100            | 0.00E+00        | 1.60E+00      | 1.60E+00      | 1.11E+02              |          | 2.49E+01      | 3.00E+02      | SD-04      | 1.64E+02               |                         |          |                         |          |                          |

**Table 17**  
**Statistical Summary and Selection of COPCs in EU6 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent               | 95% UCL<br>mg/kg | Logarithmic<br>95% UCL<br>mg/kg | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Unrestricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|---------------------------|------------------|---------------------------------|--------------------------------|--|--|-----------------------------------|--------------------------|
| Chrysene                  | 1.03E+02         | 4.42E+21                        | Normal/Lognormal               | 7.60E+01                                 | 8.75E+01                                 | no                                | YES**                    |
| Dibenz(a,h)anthracene     | 1.25E+01         | 9.25E+19                        | Normal/Lognormal               | 9.60E+00                                 | 8.75E+02                                 | YES                               | YES - COPC               |
| Dibenzofuran              | 1.96E+02         | 1.09E+63                        | Lognormal                      | 1.50E+02                                 | 3.13E+02                                 | no                                |                          |
| Diethylphthalate          | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 1.97E+03                                 | no                                |                          |
| Dimethylphthalate         | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 7.82E+05                                 | no                                |                          |
| Di-n-butylphthalate       | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 2.28E+03                                 | no                                |                          |
| Di-n-octylphthalate       | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 1.56E+03                                 | no                                |                          |
| Fluoranthene              | 6.11E+02         | 1.50E+33                        | Normal/Lognormal               | 4.70E+02                                 | 3.13E+03                                 | no                                |                          |
| Fluorene                  | 3.40E+02         | 6.19E+61                        | Lognormal                      | 2.60E+02                                 | 3.13E+03                                 | no                                |                          |
| Hexachlorobenzene         | 5.16E+01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 3.99E-01                                 | YES                               | YES - COPC               |
| Hexachlorobutadiene       | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 8.82E-02                                 | YES                               | no                       |
| Hexachlorocyclopentadiene | 2.58E+00         | 4.94E+11                        | Unknown                        | 2.00E+00                                 | 9.51E-01                                 | YES                               | YES - COPC               |
| Hexachloroethane          | 5.16E+01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 4.56E+01                                 | no                                |                          |
| Indeno(1,2,3-cd)pyrene    | 5.05E+01         | 3.03E+21                        | Normal/Lognormal               | 3.90E+01                                 | 8.75E-01                                 | YES                               | YES***                   |
| Isophorone                | 5.16E+01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 6.72E+02                                 | no                                |                          |
| Naphthalene               | 1.81E+01         | 2.21E+21                        | Normal/Lognormal               | 1.40E+01                                 | 6.45E+02                                 | no                                |                          |
| Nitrobenzene              | 5.16E+01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 8.41E+00                                 | no                                |                          |
| N-nitrosodi-n-propylamine | 5.16E+01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 9.12E-02                                 | YES                               | YES - COPC               |
| N-nitrosodiphenylamine    | 5.16E+01         | 7.97E+10                        | Lognormal                      | 4.00E-01                                 | 1.30E+02                                 | no                                |                          |
| Pentachlorophenol         | 2.58E+00         | 4.94E+11                        | Unknown                        | 2.00E+00                                 | 2.66E+00                                 | no                                |                          |
| Phenanthrene              | 1.14E+03         | 6.20E+60                        | Lognormal                      | 8.70E+02                                 | 2.35E+03                                 | no                                |                          |
| Phenol                    | 1.03E+00         | 1.77E+11                        | Unknown                        | 8.00E-01                                 | 4.69E+04                                 | no                                |                          |
| Pyrene                    | 3.88E+02         | 3.84E+30                        | Normal/Lognormal               | 3.00E+02                                 | 2.35E+03                                 | no                                |                          |

NA - Not Available

\* Constituent will be retained as a COPC due to lack of Tier I TRG.

\*\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

\*\*\*Logarithmic 95% UCL is less than benchmark but retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.

Table 18

Statistical Summary and Selection of COPCs in EU6 Surface Water  
Kerr McGee, Hattiesburg, MS

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/L | Maximum mg/L | Mean mg/L | Logarithmic Mean mg/L | Maximum Detected mg/L | Detected Qualifier | Maximum Concentration mg/L | Location of Maximum Concentration | Standard Deviation mg/L |
|------------------------|------------|---------------|----------------|-----------------|--------------|--------------|-----------|-----------------------|-----------------------|--------------------|----------------------------|-----------------------------------|-------------------------|
| Semivolatiles          |            |               |                |                 |              |              |           |                       |                       |                    |                            |                                   |                         |
| Acenaphthene           | 83-32-9    | 2             | 1              | 50              | 1.00E-03     | 9.00E-03     | J         | 4.75E-03              | 2.12E-03              | 9.00E-03           | J                          | SW-03                             | 6.01E-03                |
| Benzo(a)anthracene     | 56-55-3    | .2            | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00           | NA                         | SW-03                             | 0.00E+00                |
| Benzo(b)pyrene         | 50-32-8    | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00           | NA                         | SW-03                             | 0.00E+00                |
| Benzo(b)fluoranthene   | 205-99-2   | 2             | 1              | 50              | 1.00E-03     | 1.00E-03     | J         | 4.75E-03              | 2.12E-03              | 9.00E-03           | J                          | SW-03                             | 6.01E-03                |
| Benzo(k)fluoranthene   | 207-08-9   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00           | NA                         | SW-03                             | 0.00E+00                |
| Chrysene               | 218-01-9   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00           | NA                         | SW-03                             | 0.00E+00                |
| Dibenz(a,h)anthracene  | 53-70-3    | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00           | NA                         | SW-03                             | 0.00E+00                |
| Fluoranthene           | 206-44-0   | 2             | 2              | 100             | 0.00E+00     | 0.00E+00     | 1.20E-02  | 1.25E-02              | 1.30E-02              |                    |                            | SW-03                             | 7.07E-04                |
| Fluorene               | 86-73-7    | 2             | 1              | 50              | 1.00E-03     | 1.00E-03     | 1.10E-02  | 5.75E-03              | 2.35E-03              | 1.10E-02           |                            | SW-03                             | 7.42E-03                |
| Indeno(1,2,3-cd)pyrene | 193-39-5   | 2             | 0              | 0               | 1.00E-03     | 1.00E-03     | NA        | 5.00E-04              | 5.00E-04              | 0.00E+00           | NA                         | SW-03                             | 0.00E+00                |

NA - Not applicable; constituent not detected in media.



Table 18

Statistical Summary and Selection of COPCs in EU6 Surface Water  
Kerr McGee, Hattiesburg, MS

| Constituent            | Human Health    |                                |                                   |   |  |                                   |
|------------------------|-----------------|--------------------------------|-----------------------------------|---|--|-----------------------------------|
|                        | 95% UCL<br>mg/L | Logarithmic<br>95% UCL<br>mg/L | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/L | Consumption of Water<br>& Organisms AWQC<br>mg/L | Is Maximum<br>Detected ><br>AWQC? |
| <b>Semivolatile</b>    |                 |                                |                                   |   |  |                                   |
| Acenaphthene           | 3.16E-02        | 6.39E+48                       | Unknown                           | 9.00E-03                                | 1.20E+00   | no                                |
| Benz(a)anthracene      | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |
| Benz(a)pyrene          | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |
| Benz(b)fluoranthene    | 3.16E-02        | 6.39E+48                       | Unknown                           | 9.00E-03                                | 4.40E-06   | YES - COPC                        |
| Benz(k)fluoranthene    | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |
| Chrysene               | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |
| Dibenz(a,h)anthracene  | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |
| Fluoranthene           | 1.57E-02        | 1.53E-02                       | Unknown                           | 1.30E-02                                | 3.00E-01   | no                                |
| Fluorene               | 3.89E-02        | 1.87E+56                       | Unknown                           | 1.10E-02                                | 1.30E+00   | no                                |
| Indeno[1,2,3-cd]pyrene | 5.00E-04        | 5.00E-04                       | Unknown                           | 5.00E-04                                | 4.40E-06   | YES*                              |

\*Retained as a COPC, as per MDEQ Comments (8/2/2000); constituent is a member of carcinogenic PAH family, one of which has been retained as a COPC.



**Table 19**  
**Summary of Human Health Exposure Parameters**  
**Kerr McGee, Hattiesburg, MS**

| Receptors:   | Adolescent<br>Visitor | Maintenance<br>Worker | Construction<br>Worker | Off-Site<br>Resident<br>Child | Off-Site<br>Resident<br>Adult |
|--|-----------------------|-----------------------|------------------------|-------------------------------|-------------------------------|
| <b>Parameter</b>                                     | <b>Units</b>          |                       |                        |                               |                               |
| Surface area available for exposure - soil           | cm <sup>2</sup> /day  | 3052                  | 1                      | 3000                          | 1                             |
| Surface area available for exposure - sed. & sw      | cm <sup>2</sup> /day  | 2311                  | 1                      | 3620                          | 1                             |
| Total skin surface area                              | cm <sup>2</sup>       | 12768.3               | 2                      | 20000                         | 2                             |
| Skin surface area available for exposure - soil      | %                     | 23.9%                 | 2                      | 15%                           | 2                             |
| Skin surface area available for exposure - sed. & sw | %                     | 18.1%                 | 2                      | 18.1%                         | 2                             |
| Adherence factor - soil                              | mg/cm <sup>2</sup>    | 0.026                 | 2                      | 0.038                         | 2                             |
| Adherence factor - sed.                              | mg/cm <sup>3</sup>    | 0.044                 | 2                      | 0.034                         | 2                             |
| Dermal absorption factor - benzo(a)pyrene            |                       | 0.03                  | 3                      | 0.03                          | 3                             |
| Dermal absorption factor - other PAHs                |                       | 0.1                   | 3                      | 0.1                           | 3                             |
| Exposure time  | hours/day             | 1                     | 5                      | 1                             | 5                             |
| Exposure frequency - soils                           | days/year             | 12                    | 5                      | 150                           | 5                             |
| Exposure frequency - sed. & sw                       | days/year             | 12                    | 5                      | 30                            | 5                             |
| Exposure frequency - EU 4 soil                       | days/year             | 12                    | 5                      | 30                            | 5                             |
| Exposure duration                                    | years                 | 10                    | 6                      | 25                            | 6                             |
| Body weight  | kg                    | 45                    | 6                      | 70                            | 6                             |
| Averaging time - noncarcinogenic                     | days                  | 3650                  | 7                      | 9125                          | 7                             |
| Averaging time - carcinogenic                        | days                  | 25550                 | 7                      | 25550                         | 7                             |
| Ingestion rate - soil                                | mg/day                | 100                   | 2                      | 100                           | 2                             |
| Ingestion rate - surface water                       | L/hour                | 0.01                  | 6                      | 0.01                          | 6                             |
| Matrix effect - PAHs                                 |                       | 0.29                  | 9                      | 0.29                          | 9                             |
| Inhalation rate                                      | m <sup>3</sup> /day   | NA                    | NA                     | 20                            | 6                             |
| Retention factor - semivolatiles                     |                       | NA                    | NA                     | 0.75                          | 8                             |

NA - Not Applicable

1 Calculated

2 USEPA 1997, Exposure Factors Handbook

3 USEPA 1995, Region III Technical Guidance Manual: Assessing Dermal Exposure to Soil

4 USEPA 1992, Dermal Exposure Assessment

5 Reasonable Maximum

6 USEPA 1995, Region IV

7 USEPA 1991, HHEM Supplemental Guidances

8 International Commission on Radiological Protection, 1968

9 Magee et al., 1996



**Table 20**

**Particulate Emission Rate for Vehicular Movement and Excavation**

**Kerr McGee, Hattiesburg, MS**

**Vehicular Movement**

$$E = k * (5.9) * (s/12)(S/30) * (W/3)^0.7((w/4)^0.5) * ((365-p)/365) = \quad 16.49 \quad \text{lbs/vehicle mile}$$

|     |       |   |
|-----|-------|---|
| E = | 16.49 | particulate emission rate (lbs/vehicle mile - 30 miles travelled total over 80 - 8 hr days) |
| k = | 0.5   | particle size multiplier  |
| s = | 50    | percent silt content  |
| S = | 15    | mean vehicle speed (mi/hr)  |
| W = | 12.5  | mean vehicle weight (ton)   |
| w = | 8     | mean number of wheels per vehicle   |
| p = | 110   | mean number of days with ≥ 0.01 inches of precipitation per year                            |

Emission Rate      lbs/sec =  $(E \text{ lbs/mi}) * (30 \text{ mi/job}) * (\text{job}/80 \text{ days}) * (1 \text{ day}/8 \text{ hrs}) * (1 \text{ hr}/3600 \text{ sec})$

2.15E-04      lbs/sec

9.74E-02      g/sec

0.00010      kg/sec

**Excavation**

$$E = (1.0 * s^{1.5})/M^{1.4} = \quad 7.90E+00 \quad \text{lbs/hour}$$

|     |          |                                      |
|-----|----------|--------------------------------------|
| E = | 7.90E+00 | particulate emission factor (lbs/hr) |
| s = | 50       | percent silt content                 |
| M = | 15.1     | percent soil moisture content        |

Emission Rate =      2.20E-03      lbs/sec

0.996      g/sec

0.000996      kg/sec



**Table 21**  
**Summary of Windrose Data**  
**Kerr McGee, Hattiesburg, MS**

**GRAPHICAL EXPOSURE MODELING SYSTEM**  
**STAR STATION JACKSON/THOMPSON, MS 1974-1978**

| DIRECTION | FREQUENCY | WINDSPEED | DIRECTION                          | FREQUENCY | WINDSPEED |
|-----------|-----------|-----------|------------------------------------|-----------|-----------|
| N         | 3.33325   | 0.03      | S                                  | 0.05336   | 3.08      |
| NNE       | 1.89301   | 0.03      | SSW                                | 0.09995   | 3.29      |
| NE        | 3.56791   | 0.07      | SW                                 | 0.10061   | 3.65      |
| ENE       | 0.12132   | 4.04      | WSW                                | 0.14723   | 3.93      |
| E         | 0.04843   | 3.39      | W                                  | 0.05047   | 3.7       |
| ESE       | 0.04328   | 3.12      | WNW                                | 0.04341   | 3.51      |
| SE        | 0.03686   | 3         | NW                                 | 0.02908   | 3.25      |
| SSE       | 0.05274   | 2.99      | NNW                                | 0.0406    | 3.26      |
| STABILITY | FREQUENCY | WINDSPEED | AUXILIARY VARIABLES                |           |           |
| 1         | 259.2     | 0.13      | Afternoon mixing height (meters)   | 1409      |           |
| 2         | 0.053     | 0.24      | Nocturnal mixing height (meters)   | 472       |           |
| 3         | 11.3      | 1         | Ambient air temperature (Kelvin)   | 303.6     |           |
| 4         | 0.01264   | 2.17      | Precipitation frequency (fraction) | 289.8     |           |
| 5         | 0.08137   | 2.98      | Precipitation intensity (mm/hour)  | 73.66     |           |
| 6         | 0.1315    | 3.91      | Grand average windspeed (m/s)      | 4.69      |           |



*Table 22  
Summary of Toxicity Values  
Kerr McGee, Hattiesburg, MS*

EPA-NCEA Regional Summary provisional value from Region III RBC Tables April 2000

H - Values are published in HEAST 10003

1 - Values are published in HEA 3.1, 1993.

KIS - Values are available in IKS, 2000

J. Values are withdrawn from other EPA data.

Region IV - Region IV default value, 1995

W - Withdrawn from IRIS or

4toxvals.xls

**Table 23**  
**Summary of Hazard and Risk Calculations**  
**Kerr McGee, Hattiesburg, MS**

| Source/Pathway                          | Potentially Exposed Population | Total Hazard Index | Total Cancer Risk | Driving Constituent | Table Referenced |
|---|--------------------------------|--------------------|-------------------|---------------------|------------------|
| Dermal Exposure to Sediment in EU1      | Visitor<br>Sub-Total           | NA<br>NA           | 1E-09<br>1E-09    |                     | 24               |
| Dermal Exposure to Surface Water in EU1 | Visitor                        | NA                 | 4E-07             |                     | 25               |
| Oral Exposure to Surface Water in EU1   | Visitor<br>Sub-Total           | 2E-07<br>2E-07     | NA<br>4E-07       |                     | 26               |
| Dermal Exposure to Surface Soil in EU2  | Visitor<br>Sub-Total           | NA<br>NA           | 1E-08<br>1E-08    |                     | 27               |
| Dermal Exposure to Surface Soil in EU3  | Visitor<br>Sub-Total           | NA<br>NA           | 2E-09<br>2E-09    |                     | 28               |
| Dermal Exposure to Sediment in EU4      | Visitor                        | 2E-03              | 5E-07             |                     | 29               |
| Oral Exposure to Sediment in EU4        | Visitor<br>Sub-Total           | 8E-03<br>1E-02     | 4E-08<br>5E-07    |                     | 30               |
| Dermal Exposure to Surface Water in EU4 | Visitor                        | 4E-05              | 3E-07             |                     | 31               |
| Oral Exposure to Surface Water in EU4   | Visitor<br>Sub-Total           | 2E-05<br>4E-05     | 4E-11<br>3E-07    |                     | 32               |
| Dermal Exposure to Surface Soil in EU4  | Visitor<br>Sub-Total           | 1E-03<br>9E-03     | 1E-06<br>1E-06    |                     | 33               |
| Oral Exposure to Surface Soil in EU4    | Visitor                        | 8E-03              | 1E-08             |                     | 34               |
| Dermal Exposure to Surface Soil in EU5  | Visitor<br>Sub-Total           | NA<br>NA           | 1E-07<br>1E-07    |                     | 35               |
| <b>Visitor Total:</b>                   |                                | <b>2E-02</b>       | <b>3E-06</b>      |                     |                  |

|   |                                 |                |                |                |    |
|---|---------------------------------|----------------|----------------|----------------|----|
| Dermal Exposure to Sediment in EU1      | Maintenance Worker<br>Sub-Total | NA<br>NA       | 7E-09<br>7E-09 |                | 36 |
| Dermal Exposure to Surface Water in EU1 | Maintenance Worker<br>Sub-Total | NA<br>NA       | 8E-07<br>8E-07 |                | 37 |
| Dermal Exposure to Surface Soil in EU2  | Maintenance Worker<br>Sub-Total | NA<br>NA       | 2E-07<br>2E-07 |                | 38 |
| Dermal Exposure to Sediment in EU4      | Maintenance Worker              | 3E-03          | 2E-06          | Benzo(a)pyrene | 39 |
| Oral Exposure to Sediment in EU4        | Maintenance Worker<br>Sub-Total | 1E-02<br>2E-02 | 1E-07<br>1E-07 |                | 40 |
| Dermal Exposure to Surface Water in EU4 | Maintenance Worker              | 1E-04          | 2E-06          | *              | 41 |
| Oral Exposure to Surface Water in EU4   | Maintenance Worker<br>Sub-Total | 3E-05<br>1E-04 | 2E-10<br>2E-06 |                | 42 |
| Dermal Exposure to Surface Soil in EU4  | Maintenance Worker<br>Sub-Total | 3E-03<br>1E-02 | 8E-06<br>2E-07 | Benzo(a)pyrene | 43 |
| Oral Exposure to Surface Soil in EU4    | Maintenance Worker<br>Sub-Total | 1E-02          | 8E-06          |                | 44 |
| Dermal Exposure to Surface Soil in EU5  | Maintenance Worker<br>Sub-Total | NA<br>NA       | 3E-06<br>3E-06 | Benzo(a)pyrene | 45 |
| <b>Maintenance Worker Total:</b>        |                                 | <b>3E-02</b>   | <b>1E-05</b>   |                |    |



**Table 23**  
**Summary of Hazard and Risk Calculations**  
**Kerr McGee, Hattiesburg, MS**

| Source/Pathway                          | Potentially Exposed Population | Total Hazard Index | Total Cancer Risk | Driving Constituent | Table Referenced |
|---|--------------------------------|--------------------|-------------------|---------------------|------------------|
| Dermal Exposure to Sediment in EU1      | Construction Worker            | NA                 | 2E-10             |                     | 46               |
|   | Sub-Total                      | NA                 | 2E-10             |                     |                  |
| Dermal Exposure to Surface Water in EU1 | Construction Worker            | NA                 | 9E-09             |                     | 47               |
|   | Sub-Total                      | NA                 | 9E-09             |                     |                  |
| Dermal Exposure to Soil in EU2          | Construction Worker            | NA                 | 2E-08             |                     | 48               |
| Inhalation of Fugitive Dust in EU2      | Construction Worker            | NA                 | 8E-09             |                     | 49               |
|   | Sub-Total                      | NA                 | 3E-08             |                     |                  |
| Dermal Exposure to Sediment in EU4      | Construction Worker            | NA                 | 7E-08             |                     | 50               |
| Oral Exposure to Sediment in EU4        | Construction Worker            | NA                 | 2E-09             |                     | 51               |
|   | Sub-Total                      | NA                 | 8E-08             |                     |                  |
| Dermal Exposure to Surface Water in EU4 | Construction Worker            | 3E-07              | 2E-08             |                     | 52               |
| Oral Exposure to Surface Water in EU4   | Construction Worker            | 9E-07              | 1E-11             |                     | 53               |
|   | Sub-Total                      | 1E-06              | 1E-11             |                     |                  |
| Dermal Exposure to Soil in EU4          | Construction Worker            | 2E-04              | 4E-07             |                     | 54               |
| Oral Exposure to Soil in EU4            | Construction Worker            | 1E-04              | 2E-09             |                     | 55               |
| Inhalation of Fugitive Dust in EU4      | Construction Worker            | NA                 | 1E-07             |                     | 56               |
|   | Sub-Total                      | 3E-04              | 6E-07             |                     |                  |
| Dermal Exposure to Soil in EU5          | Construction Worker            | NA                 | 1E-07             |                     | 57               |
| Inhalation of Fugitive Dust in EU5      | Construction Worker            | NA                 | 5E-08             |                     | 58               |
|   | Sub-Total                      | NA                 | 2E-07             |                     |                  |
| <b>Construction Worker Total:</b>       |                                | <b>3E-04</b>       | <b>9E-07</b>      |                     |                  |
|   |                                |                    |                   |                     |                  |
| Dermal Exposure to Sediment in EU6      | Child Off-Site Resident        | NA                 | 4E-07             |                     | 59               |
| Oral Exposure to Sediment in EU6        | Child Off-Site Resident        | 4E-05              | 2E-08             |                     | 60               |
|   | Sub-Total                      | 4E-05              | 4E-07             |                     |                  |
| Dermal Exposure to Sediment in EU6      | Adult Off-Site Resident        | NA                 | 2E-06             | *                   | 61               |
| Oral Exposure to Sediment in EU6        | Adult Off-Site Resident        | 4E-05              | 6E-08             |                     | 62               |
|   | Sub-Total                      | 4E-05              | 2E-06             |                     |                  |
| Dermal Exposure to Surface Water in EU6 | Child Off-Site Resident        | NA                 | 2E-06             | *                   | 63               |
|   | Sub-Total                      | NA                 | 2E-06             |                     |                  |
| Dermal Exposure to Surface Water in EU6 | Adult Off-Site Resident        | NA                 | 1E-06             |                     | 64               |
|   | Sub-Total                      | NA                 | 1E-06             |                     |                  |
| <b>Off-Site Resident Total:</b>         |                                | <b>7E-05</b>       | <b>5E-06</b>      |                     |                  |

\*Estimated carcinogenic risk level is below *de minimis* level as no single constituent exceeded  $1 \times 10^{-6}$  and the cumulative carcinogenic risk is below  $1 \times 10^{-4}$  (Section 501, MCEQ, 1999).



**Table 24****Dermal Exposure to EU1 Sediment by an Adolescent Visitor (Aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                        |  |
|---|----------------------|---|------------------------|--|
| Cs - Concentration in sediment =                            | mg/kg                | chem. spec.   |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 2311  | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 12768.3   | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 18.1%   | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.044   | USEPA 1997, EFH        |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                      | 0.03  | USEPA 1995, Region III |  |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |                      | 0.1   | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year            | 12  | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                | 10  | USEPA 1995, Region IV  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06  |                        |  |
| BW - Body weight =  | kg                   | 45  | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 3650  | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550   | USEPA 1991, HHEM       |  |

| Constituent            | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average                         |   |             |
|------------------------|---------------------------------|--------------------------------|------------------------------|--------------|---------------------------------|---|-------------|
|                        |                                 |                                |                              |              | Lifetime Daily Intake mg/kg-day | Daily Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>   |                                 |                                |                              |              |                                 |   |             |
| Benzo(a)anthracene     | 5.90E-01                        | 1.31E-09                       | NA                           | NA           | 1.88E-10                        | 7.30E-01                                | 1.37E-10    |
| Benzo(a)pyrene         | 3.90E-01                        | 8.69E-10                       | NA                           | NA           | 1.24E-10                        | 7.30E+00                                | 9.06E-10    |
| Benzo(b)fluoranthene   | 5.80E-01                        | 1.29E-09                       | NA                           | NA           | 1.85E-10                        | 7.30E-01                                | 1.35E-10    |
| Benzo(k)fluoranthene   | 1.90E-01                        | 4.23E-10                       | NA                           | NA           | 6.05E-11                        | 7.30E-02                                | 4.42E-12    |
| Chrysene               | 5.30E-01                        | 1.18E-09                       | NA                           | NA           | 1.69E-10                        | 7.30E-03                                | 1.23E-12    |
| Dibenz(a,h)anthracene  | 6.20E-02                        | 1.38E-10                       | NA                           | NA           | 1.97E-11                        | 7.30E+00                                | 1.44E-10    |
| Indeno(1,2,3-cd)pyrene | 2.20E-01                        | 4.90E-10                       | NA                           | NA           | 7.00E-11                        | 7.30E-01                                | 5.11E-11    |

NA - Not Available

Total Cancer Risk = 1.38E-09



Table 25

Dermal Exposure to EIU Surface Water by an Adolescent Visitor (aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                                     | $\frac{C_w * SA * K_p * ABS * ET * EF * ED * CF}{BW * AT}$ |  |                              |              |   |                                   |             |
|---|-------------------------------------|--|--|------------------------------|--------------|---|-----------------------------------|-------------|
| Cw - Concentration in surface water =                       | mg/L                                |  | see below                              |                              |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup>                     | 2311   | calculated                             |                              |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                     | 12768.3  | USEPA 1997, EFH                        |                              |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |                                     | 18.1%  | USEPA 1997, EFH                        |                              |              |   |                                   |             |
| Kp - Dermal permeability constant =                         | cm/hr                               | see below  |  |                              |              |   |                                   |             |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                                     | 0.03   | USEPA 1995, Region III                 |                              |              |   |                                   |             |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |                                     | 0.1  | USEPA 1995, Region III                 |                              |              |   |                                   |             |
| ET - Exposure time =  | hrs/day                             | 1  | USEPA 1992, Dermal Exposure Assessment |                              |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year                           | 12   | reasonable assumption                  |                              |              |   |                                   |             |
| ED - Exposure duration =                                    | years                               | 10   | USEPA 1995, Region IV                  |                              |              |   |                                   |             |
| CF - Conversion factor =                                    | L/cm <sup>3</sup>                   | 1.00E-03   |  |                              |              |   |                                   |             |
| BW - Body weight =  | kg                                  | 45   | USEPA 1995, Region IV                  |                              |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                                | 3650   | USEPA 1991, HHEM                       |                              |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                                | 25550  | USEPA 1991, HHEM                       |                              |              |   |                                   |             |
| Constituent   | Concentration in Surface Water mg/L | Kp cm/hr   | Average Daily Intake mg/kg-day         | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|   |                                     |  |  |                              |              |   |                                   |             |
| <b>Semivolatiles</b>  |                                     |  |  |                              |              |   |                                   |             |
| Benzo(a)anthracene  | 1.00E-03                            | 8.10E-01   | 1.37E-07                               | NA                           | NA           | 1.95E-08                                | 7.30E-01                          | 1.43E-08    |
| Benzo(a)pyrene  | 5.00E-04                            | 1.20E+00   | 1.01E-07                               | NA                           | NA           | 1.45E-08                                | 7.30E+00                          | 1.06E-07    |
| Benzo(b)fluoranthene  | 5.00E-04                            | 1.20E+00   | 1.01E-07                               | NA                           | NA           | 1.45E-08                                | 7.30E-01                          | 1.06E-08    |
| Benzo(k)fluoranthene  | 5.00E-04                            | 4.48E+01   | 3.78E-06                               | NA                           | NA           | 5.40E-07                                | 7.30E-02                          | 3.94E-08    |
| Chrysene  | 5.00E-04                            | 8.10E-01   | 6.84E-08                               | NA                           | NA           | 9.77E-09                                | 7.30E-03                          | 7.13E-11    |
| Dibenz(a,h)anthracene                                       | 5.00E-04                            | 2.70E+00   | 2.28E-07                               | NA                           | NA           | 3.26E-08                                | 7.30E+00                          | 2.38E-07    |
| Indeno(1,2,3-cd)pyrene                                      | 5.00E-04                            | 1.90E+00   | 1.60E-07                               | NA                           | NA           | 2.29E-08                                | 7.30E-01                          | 1.67E-08    |

NA - Not Available

Total Cancer Risk = 4.24E-07



**Table 26**

**Oral Exposure to EU1 Surface Water by an Adolescent Visitor (aged 7-16 years)**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u>Csw*IngR*EF*ED*ET</u> |  |  |
|--|-----------|--------------------------|--|--|
|  |           | BW*AT                    |  |  |
| Csw - Concentration in surface water =               | mg/L      | see below                |  |  |
| IngR - Ingestion rate for surface water =            | L/hour    | 0.01                     | USEPA 1995, Region IV                  |  |
| EF - Exposure frequency =                            | days/year | 12                       | reasonable assumption                  |  |
| ED - Exposure duration =                             | years     | 10                       | USEPA 1995, Region IV                  |  |
| ET - Exposure time =                                 | hrs/day   | 1                        | USEPA 1992, Dermal Exposure Assessment |  |
| BW - Body weight =                                   | kg        | 45                       | USEPA 1995, Region IV                  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650                     | USEPA 1991, HHEM                       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550                    | USEPA 1991, HHEM                       |  |

| Constituent          | Concentration         | Average Daily Intake | Oral Chronic RfD | Hazard Index | Average                         |  |             |
|----------------------|-----------------------|----------------------|------------------|--------------|---------------------------------|--|-------------|
|                      | in Surface Water mg/L |                      |                  |              | Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b> |                       |                      |                  |              |                                 |  |             |
| Pyrene               | 1.00E-03              | 7.31E-09             | 3.00E-02         | 2.44E-07     | 1.04E-09                        | NA                                     | NA          |

NA - Not Applicable

Total Hazard Index = 2.44E-07



Table 27

Dermal Exposure to EU2 Surface Soil (0-1') by an Adolescent Visitor (aged 10-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                      | <u><math>C_s * SA * AH * ABS * EF * ED * CF</math></u> |                        |  |  |
|---|----------------------|--|------------------------|--|--|
|   |                      | BW * AT  |                        |  |  |
| $C_s$ - Concentration in soil =                             | mg/kg                | chem. spec.  |                        |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 3052   | calculated             |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 12768.3  | USEPA 1997, EFH        |  |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 23.9%  | USEPA 1997, EFH        |  |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.026  | USEPA 1997, EFH        |  |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                      | 0.03   | USEPA 1995, Region III |  |  |
| EF - Exposure frequency =                                   | days/year            | 12   | reasonable assumption  |  |  |
| ED - Exposure duration =                                    | years                | 10   | USEPA 1995, Region IV  |  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06   |                        |  |  |
| BW - Body weight =  | kg                   | 45   | USEPA 1995, Region IV  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 3650   | USEPA 1991, HHEM       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550  | USEPA 1991, HHEM       |  |  |

| Constituent            | Concentration in<br>Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Dermal Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average                               |   |             |
|------------------------|-----------------------------------|--------------------------------------|------------------------------------|-----------------|---------------------------------------|---|-------------|
|                        |                                   |                                      |                                    |                 | Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>   |                                   |                                      |                                    |                 |                                       |   |             |
| Benzo(a)anthracene     | 6.70E+00                          | 1.17E-08                             | NA                                 | NA              | 1.66E-09                              | 7.30E-01                                | 1.22E-09    |
| Benzo(a)pyrene         | 5.08E+00                          | 8.8342E-09                           | NA                                 | NA              | 1.26E-09                              | 7.30E+00                                | 9.21E-09    |
| Benzo(b)fluoranthene   | 9.20E+00                          | 1.60E-08                             | NA                                 | NA              | 2.29E-09                              | 7.30E-01                                | 1.67E-09    |
| Benzo(k)fluoranthene   | 2.93E+00                          | 5.10E-09                             | NA                                 | NA              | 7.28E-10                              | 7.30E-02                                | 5.31E-11    |
| Chrysene               | 8.00E+00                          | 1.3912E-08                           | NA                                 | NA              | 1.99E-09                              | 7.30E-03                                | 1.45E-11    |
| Dibenz(a,h)anthracene  | 4.93E-01                          | 8.57E-10                             | NA                                 | NA              | 1.22E-10                              | 7.30E+00                                | 8.94E-10    |
| Indeno(1,2,3-cd)pyrene | 3.70E+00                          | 6.43E-09                             | NA                                 | NA              | 9.19E-10                              | 7.30E-01                                | 6.71E-10    |

NA - Not Available

Total Cancer Risk = 1.37E-08



Table 28

Dermal Exposure to EU3 Surface Soil (0-1') by an Adolescent Visitor (aged 10-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                                   | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                    |   |   |             |          |
|---|-----------------------------------|---|------------------------------------|---|---|-------------|----------|
| Cs - Concentration in soil =                                | mg/kg                             | chem. spec.   |                                    |   |   |             |          |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day              | 3052  | calculated                         |   |   |             |          |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                   | 12768.3   | USEPA 1997, EFH                    |   |   |             |          |
| Fs - Fraction of skin surface area available for exposure = |                                   | 23.9%   | USEPA 1997, EFH                    |   |   |             |          |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                | 0.026   | USEPA 1997, EFH                    |   |   |             |          |
| ABS <sub>top</sub> - Absorption - B(a)P =                   |                                   | 0.03  | USEPA 1995, Region III             |   |   |             |          |
| EF - Exposure frequency =                                   | days/year                         | 12  | reasonable assumption              |   |   |             |          |
| ED - Exposure duration =                                    | years                             | 10  | USEPA 1995, Region IV              |   |   |             |          |
| CF - Conversion factor =                                    | kg/mg                             | 1.00E-06  |                                    |   |   |             |          |
| BW - Body weight =  | kg                                | 45  | USEPA 1995, Region IV              |   |   |             |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                              | 3650  | USEPA 1991, HHEM                   |   |   |             |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                              | 25550   | USEPA 1991, HHEM                   |   |   |             |          |
|   |                                   |   |                                    |   |   |             |          |
| Constituent   | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day                | Dermal<br>Chronic RfD<br>mg/kg-day | Average Lifetime<br>Daily Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |          |
| Semivolatiles   |                                   |   |                                    |   |   |             |          |
| Benzo(a)anthracene  | 5.40E-01                          | 9.39E-10  | NA                                 | NA  | 1.34E-10                                | 7.30E-01    | 9.79E-11 |
| Benzo(a)pyrene  | 7.10E-01                          | 1.23E-09  | NA                                 | NA  | 1.76E-10                                | 7.30E+00    | 1.29E-09 |
| Benzo(b)fluoranthene  | 1.40E+00                          | 2.43E-09  | NA                                 | NA  | 3.48E-10                                | 7.30E-01    | 2.54E-10 |
| Benzo(k)fluoranthene  | 4.90E-01                          | 8.52E-10  | NA                                 | NA  | 1.22E-10                                | 7.30E-02    | 8.89E-12 |
| Chrysene  | 8.70E-01                          | 1.51E-09  | NA                                 | NA  | 2.16E-10                                | 7.30E-03    | 1.58E-12 |
| Dibenz(a,h)anthracene                                       | 1.60E-01                          | 2.78E-10  | NA                                 | NA  | 3.97E-11                                | 7.30E+00    | 2.90E-10 |
| Indeno(1,2,3-cd)pyrene                                      | 6.00E-01                          | 1.04E-09  | NA                                 | NA  | 1.49E-10                                | 7.30E-01    | 1.09E-10 |

NA - Not Available

Total Cancer Risk = 2.05E-09



**Table 29****Dermal Exposure to EU4 Sediment by an Adolescent Visitor (Aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                                 | $\frac{\text{Cs} * \text{SA} * \text{AH} * \text{ABS} * \text{EF} * \text{ED} * \text{CF}}{\text{BW} * \text{AT}}$ |                              |              |   |             |          |
|---|---------------------------------|--|------------------------------|--------------|---|-------------|----------|
| Cs - Concentration in sediment =                            | mg/kg                           | chem. spec.  |                              |              |   |             |          |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day            | 2311   | calculated                   |              |   |             |          |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                 | 12768.3  | USEPA 1997, EFH              |              |   |             |          |
| Fs - Fraction of skin surface area available for exposure = |                                 | 18.1%  | USEPA 1997, EFH              |              |   |             |          |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>              | 0.044  | USEPA 1997, EFH              |              |   |             |          |
| ABS <sub>pap</sub> - Absorption - B(a)P =                   |                                 | 0.03   | USEPA 1995, Region III       |              |   |             |          |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |                                 | 0.1  | USEPA 1995, Region III       |              |   |             |          |
| EF - Exposure frequency =                                   | days/year                       | 12   | reasonable assumption        |              |   |             |          |
| ED - Exposure duration =                                    | years                           | 10   | USEPA 1995, Region IV        |              |   |             |          |
| CF - Conversion factor =                                    | kg/mg                           | 1.00E-06   |                              |              |   |             |          |
| BW - Body weight =  | kg                              | 45   | USEPA 1995, Region IV        |              |   |             |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                            | 3650   | USEPA 1991, HHEM             |              |   |             |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                            | 25550  | USEPA 1991, HHEM             |              |   |             |          |
| Constituent   | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day   | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |             |          |
|   |                                 |  |                              |              | Cancer Slope Factor 1/(mg/kg-day)       | Cancer Risk |          |
| <b>Semivolatiles</b>  |                                 |  |                              |              |   |             |          |
| Benzo(a)anthracene  | 3.30E+02                        | 7.35E-07   | NA                           | NA           | 1.05E-07                                | 7.30E-01    | 7.67E-08 |
| Benzo(a)pyrene  | 1.30E+02                        | 2.90E-07   | NA                           | NA           | 4.14E-08                                | 7.30E+00    | 3.02E-07 |
| Benzo(b)fluoranthene  | 1.80E+02                        | 4.01E-07   | NA                           | NA           | 5.73E-08                                | 7.30E-01    | 4.18E-08 |
| Benzo(k)fluoranthene  | 6.40E+01                        | 1.43E-07   | NA                           | NA           | 2.04E-08                                | 7.30E-02    | 1.49E-09 |
| Carbazole   | 5.90E+02                        | 1.31E-06   | NA                           | NA           | 1.88E-07                                | 2.00E-02    | 3.76E-09 |
| Chrysene  | 2.90E+02                        | 6.46E-07   | NA                           | NA           | 9.23E-08                                | 7.30E-03    | 6.74E-10 |
| Dibenz(a,h)anthracene                                       | 1.20E+01                        | 2.67E-08   | NA                           | NA           | 3.82E-09                                | 7.30E+00    | 2.79E-08 |
| Dibenzofuran  | 9.40E+02                        | 2.10E-06   | 2.00E-03                     | 1.05E-03     | 2.99E-07                                | NA          | NA       |
| Indeno(1,2,3-cd)pyrene                                      | 4.70E+01                        | 1.05E-07   | NA                           | NA           | 1.50E-08                                | 7.30E-01    | 1.09E-08 |
| Naphthalene   | 3.00E+03                        | 6.69E-06   | 1.00E-02                     | 6.69E-04     | 9.55E-07                                | NA          | NA       |
| Phenanthrene  | 3.20E+03                        | 7.13E-06   | NA                           | NA           | 1.02E-06                                | NA          | NA       |

NA - Not Available

Total Hazard Index = 1.72E-03

Total Cancer Risk = 4.65E-07

**Table 30**  
**Oral Exposure to EU4 Sediment by an Adolescent Visitor (Aged 7-16 years)**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u>Cd*IngR*EF*ED*CF*ME</u> |                       |  |
|--|-----------|----------------------------|-----------------------|--|
|  |           | <u>BW*AT</u>               |                       |  |
| Cd - Concentration in sediment =                     | mg/kg     | see below                  |                       |  |
| IngR - Ingestion rate for sediment =                 | mg/day    | 100                        | USEPA 1997, EFH       |  |
| EF - Exposure frequency =                            | days/year | 12                         | reasonable assumption |  |
| ED - Exposure duration =                             | years     | 10                         | USEPA 1995, Region IV |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06                   |                       |  |
| ME <sub>s</sub> - Matrix effect - PAHs =             |           | 0.29                       | Magee, et al., 1996   |  |
| BW - Body weight =                                   | kg        | 45                         | USEPA 1995, Region IV |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650                       | USEPA 1991, HHEM      |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550                      | USEPA 1991, HHEM      |  |

| Constituent          | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Oral Chronic RfD mg/kg-day | Hazard Index | Average                         |                     |                          |
|----------------------|---------------------------------|--------------------------------|----------------------------|--------------|---------------------------------|---------------------|--------------------------|
|                      |                                 |                                |                            |              | Lifetime Daily Intake mg/kg-day | Daily 1/(mg/kg-day) | Oral Cancer Slope Factor |
| <b>Semivolatiles</b> |                                 |                                |                            |              |                                 |                     |                          |
| Carbazole            | 5.90E+02                        | 1.25E-05                       | NA                         | NA           | 1.79E-06                        | 2.00E-02            | 3.57E-08                 |
| Dibenzofuran         | 9.40E+02                        | 1.99E-05                       | 4.00E-03                   | 4.98E-03     | 2.85E-06                        | NA                  | NA                       |
| Naphthalene          | 3.00E+03                        | 6.36E-05                       | 2.00E-02                   | 3.18E-03     | 9.08E-06                        | NA                  | NA                       |
| Phenanthrene         | 3.20E+03                        | 6.78E-05                       | NA                         | NA           | 9.69E-06                        | NA                  | NA                       |

NA - Not Applicable

Total Hazard Index = 8.16E-03

Total Cancer Risk = 3.57E-08



**Table 31****Dermal Exposure to EU4 Surface Water by an Adolescent Visitor (aged 7-16 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                   | <u>Cw*SA*Kp*ABS*ET*EF*ED*CF</u><br>BW*AT |  |                                     |                 |                        |             |          |
|---|-------------------|--|--|-------------------------------------|-----------------|------------------------|-------------|----------|
| Cw - Concentration in surface water =                                   | mg/L              | see below                                |  |                                     |                 |                        |             |          |
| SA - Surface area available for exposure =                              | cm <sup>2</sup>   | 2311                                     | calculated                             |                                     |                 |                        |             |          |
| SA <sub>t</sub> - Total skin surface area =                             | cm <sup>2</sup>   | 12768.3                                  | USEPA 1997, EFH                        |                                     |                 |                        |             |          |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = |                   | 18.1%                                    | USEPA 1997, EFH                        |                                     |                 |                        |             |          |
| K <sub>p</sub> - Dermal permeability constant =                         | cm/hr             | see below                                |  |                                     |                 |                        |             |          |
| ABS <sub>bap</sub> - Absorption - B(a)P =                               |                   | 0.03                                     | USEPA 1995, Region III                 |                                     |                 |                        |             |          |
| ABS <sub>pah</sub> - Absorption - PAHs =                                |                   | 0.1                                      | USEPA 1995, Region III                 |                                     |                 |                        |             |          |
| ET - Exposure time =  | hrs/day           | 1  | USEPA 1992, Dermal Exposure Assessment |                                     |                 |                        |             |          |
| EF - Exposure frequency =   | days/year         | 12                                       | reasonable assumption                  |                                     |                 |                        |             |          |
| ED - Exposure duration =  | years             | 10                                       | USEPS 1995, Region IV                  |                                     |                 |                        |             |          |
| CF - Conversion factor =  | L/cm <sup>3</sup> | 1.00E-03                                 |  |                                     |                 |                        |             |          |
| BW - Body weight =  | kg                | 45                                       | USEPA 1995, Region IV                  |                                     |                 |                        |             |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                    | days              | 3650                                     | USEPA 1991, HHEM                       |                                     |                 |                        |             |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =                       | days              | 25550                                    | USEPA 1991, HHEM                       |                                     |                 |                        |             |          |
| Concentration<br>in Surface<br>Water                                    |                   | Average<br>Daily<br>Intake               | Dermal<br>Chronic<br>RfD               | Average<br>Lifetime<br>Daily Intake |                 | Cancer Slope<br>Factor |             |          |
| Constituent   | mg/L              | Kp<br>cm/hr                              | mg/kg-day                              | mg/kg-day                           | Hazard<br>Index | 1/(mg/kg-day)          | Cancer Risk |          |
| <b>Semivolatiles</b>  |                   |  |  |                                     |                 |                        |             |          |
| Benzo(a)anthracene  | 5.00E-03          | 8.10E-01                                 | 2.05E-07                               | NA                                  | NA              | 2.93E-08               | 7.30E-01    | 2.14E-08 |
| Benzo(a)pyrene  | 5.00E-04          | 1.20E+00                                 | 3.04E-08                               | NA                                  | NA              | 4.34E-09               | 7.30E+00    | 3.17E-08 |
| Benzo(b)fluoranthene  | 1.20E-02          | 1.20E+00                                 | 7.29E-07                               | NA                                  | NA              | 1.04E-07               | 7.30E-01    | 7.61E-08 |
| Benzo(k)fluoranthene  | 2.00E-03          | 4.48E+01                                 | 4.54E-06                               | NA                                  | NA              | 6.48E-07               | 7.30E-02    | 4.73E-08 |
| Bis(2-ethylhexyl)phthalate  | 3.00E-03          | 3.30E-02                                 | 5.01E-09                               | 1.00E-02                            | 5.01E-07        | 7.16E-10               | 1.40E-02    | 1.00E-11 |
| Carbazole   | 1.00E-02          | 3.57E-02                                 | 1.81E-08                               | NA                                  | NA              | 2.59E-09               | 2.00E-02    | 5.17E-11 |
| Chrysene  | 6.00E-03          | 8.10E-01                                 | 2.46E-07                               | NA                                  | NA              | 3.52E-08               | 7.30E-03    | 2.57E-10 |
| Dibenz(a,h)anthracene   | 5.00E-04          | 2.70E+00                                 | 6.84E-08                               | NA                                  | NA              | 9.77E-09               | 7.30E+00    | 7.13E-08 |
| Dibenzofuran  | 1.10E-02          | 1.51E-01                                 | 8.42E-08                               | 2.00E-03                            | 4.21E-05        | 1.20E-08               | NA          | NA       |
| Indeno(1,2,3-cd)pyrene  | 5.00E-04          | 1.90E+00                                 | 4.81E-08                               | NA                                  | NA              | 6.87E-09               | 7.30E-01    | 5.02E-09 |
| Phenanthrene  | 1.70E-02          | 2.30E-01                                 | 1.98E-07                               | NA                                  | NA              | 2.83E-08               | NA          | NA       |

NA - Not Available

Total Hazard Index = 4.26E-05

Total Cancer Risk = 2.53E-07



**Table 32**

**Oral Exposure to EU4 Surface Water by an Adolescent Visitor (aged 7-16 years)**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                     |           | $\frac{C_{sw} \cdot Ingr \cdot EF \cdot ED \cdot ET}{BW \cdot AT}$ |  |  |  |
|--|-----------|--|--|--|--|
| <i>C<sub>sw</sub></i> - Concentration in surface water = | mg/L      | see below  |  |  |  |
| Ingr - Ingestion rate for surface water =                | L/hour    | 0.01   | USEPA 1995, Region IV                  |  |  |
| EF - Exposure frequency =                                | days/year | 12   | reasonable assumption                  |  |  |
| ED - Exposure duration =                                 | years     | 10   | USEPA 1995, Region IV                  |  |  |
| ET - Exposure time =                                     | hrs/day   | 1  | USEPA 1992, Dermal Exposure Assessment |  |  |
| BW - Body weight =                                       | kg        | 45   | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =     | days      | 3650   | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =        | days      | 25550  | USEPA 1991, HHEM                       |  |  |

| Constituent                | Concentration<br>in Surface<br>Water<br>mg/L | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average                               |  |             |
|----------------------------|--|--------------------------------------|----------------------------------|-----------------|---------------------------------------|--|-------------|
|                            |  |                                      |                                  |                 | Lifetime Daily<br>Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>       |  |                                      |                                  |                 |                                       |  |             |
| Bis(2-ethylhexyl)phthalate | 3.00E-03                                     | 2.19E-08                             | 2.00E-02                         | 1.10E-06        | 3.13E-09                              | 1.40E-02                                     | 4.38E-11    |
| Carbazole                  | 1.00E-02                                     | 7.31E-08                             | NA                               | NA              | 1.04E-08                              | NA   | NA          |
| Dibenzofuran               | 1.10E-02                                     | 8.04E-08                             | 4.00E-03                         | 2.01E-05        | 1.15E-08                              | NA   | NA          |
| Phenanthrene               | 1.70E-02                                     | 1.24E-07                             | NA                               | NA              | 1.77E-08                              | NA   | NA          |

NA - Not Applicable

Total Hazard Index = 2.12E-05

Total Cancer Risk = 4.38E-11



Table 33

Dermal Exposure to EU4 Surface Soil (0-1') by an Adolescent Visitor (Aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                                   | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                    |                 |  |   |             |
|---|-----------------------------------|---|------------------------------------|-----------------|--|---|-------------|
| Cs - Concentration in soil =                                | mg/kg                             | chem. spec.   |                                    |                 |  |   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day              | 3052  | calculated                         |                 |  |   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                   | 12768.3   | USEPA 1997, EFH                    |                 |  |   |             |
| Fs - Fraction of skin surface area available for exposure = |                                   | 23.9%   | USEPA 1997, EFH                    |                 |  |   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                | 0.026   | USEPA 1997, EFH                    |                 |  |   |             |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                                   | 0.03  | USEPA 1995, Region III             |                 |  |   |             |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |                                   | 0.1   | USEPA 1995, Region III             |                 |  |   |             |
| EF - Exposure frequency =                                   | days/year                         | 12  | reasonable assumption              |                 |  |   |             |
| ED - Exposure duration =                                    | years                             | 10  | USEPA 1995, Region IV              |                 |  |   |             |
| CF - Conversion factor =                                    | kg/mg                             | 1.00E-06  |                                    |                 |  |   |             |
| BW - Body weight =  | kg                                | 45  | USEPA 1995, Region IV              |                 |  |   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                              | 3650  | USEPA 1991, HHEM                   |                 |  |   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                              | 25550   | USEPA 1991, HHEM                   |                 |  |   |             |
| Constituent   | Concentration<br>in Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day                | Dermal Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime<br>Daily Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |                                   |   |                                    |                 |  |   |             |
| Benzo(a)anthracene  | 9.30E+02                          | 1.62E-06  | NA                                 | NA              | 2.31E-07   | 7.30E-01                                | 1.69E-07    |
| Benzo(a)pyrene  | 5.00E+02                          | 8.70E-07  | NA                                 | NA              | 1.24E-07   | 7.30E+00                                | 9.07E-07    |
| Benzo(b)fluoranthene  | 5.30E+02                          | 9.22E-07  | NA                                 | NA              | 1.32E-07   | 7.30E-01                                | 9.61E-08    |
| Benzo(k)fluoranthene  | 2.90E+02                          | 5.04E-07  | NA                                 | NA              | 7.20E-08   | 7.30E-02                                | 5.26E-09    |
| Carbazole   | 2.30E+02                          | 4.00E-07  | NA                                 | NA              | 5.71E-08   | 2.00E-02                                | 1.14E-09    |
| Chrysene  | 6.90E+02                          | 1.20E-06  | NA                                 | NA              | 1.71E-07   | 7.30E-03                                | 1.25E-09    |
| Dibenz(a,h)anthracene                                       | 6.40E+01                          | 1.11E-07  | NA                                 | NA              | 1.59E-08   | 7.30E+00                                | 1.16E-07    |
| Fluoranthene  | 4.60E+03                          | 8.00E-06  | 2.00E-02                           | 4.00E-04        | 1.14E-06   | NA                                      | NA          |
| Indeno(1,2,3-cd)pyrene                                      | 2.50E+02                          | 4.35E-07  | NA                                 | NA              | 6.21E-08   | 7.30E-01                                | 4.53E-08    |
| Naphthalene   | 2.20E+03                          | 3.83E-06  | 1.00E-02                           | 3.83E-04        | 5.47E-07   | NA                                      | NA          |
| Phenanthrene  | 6.40E+03                          | 1.11E-05  | NA                                 | NA              | 1.59E-06   | NA                                      | NA          |
| Pyrene  | 4.40E+03                          | 7.65E-06  | 1.50E-02                           | 5.10E-04        | 1.09E-06   | NA                                      | NA          |

NA - Not Available

Total Hazard Index = 1.29E-03

Total Cancer Risk = 1.34E-06



**Table 34**

**Oral Exposure to EU4 Surface Soil (0-1') by an Adolescent Visitor (Aged 7-16 years)  
Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u>Cd*IngR*EF*ED*CF*ME</u> |                       |  |  |
|--|-----------|----------------------------|-----------------------|--|--|
|  |           | BW*AT                      |                       |  |  |
| Cd - Concentration in soil =                         | mg/kg     | see below                  |                       |  |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 100                        | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 12                         | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 10                         | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06                   |                       |  |  |
| ME <sub>s</sub> - Matrix effect - PAHs =             |           | 0.29                       | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 45                         | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 3650                       | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550                      | USEPA 1991, HHEM      |  |  |

| Constituent          | Concentration in<br>Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average                               |                              |  |
|----------------------|-----------------------------------|--------------------------------------|----------------------------------|-----------------|---------------------------------------|------------------------------|--|
|                      |                                   |                                      |                                  |                 | Lifetime Daily<br>Intake<br>mg/kg-day | Daily<br>Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) |
| <b>Semivolatiles</b> |                                   |                                      |                                  |                 |                                       |                              |  |
| Carbazole            | 2.30E+02                          | 4.87E-06                             | NA                               | NA              | 6.96E-07                              | 2.00E-02                     | 1.39E-08                                     |
| Fluoranthene         | 4.60E+03                          | 9.75E-05                             | 4.00E-02                         | 2.44E-03        | 1.39E-05                              | NA                           | NA   |
| Naphthalene          | 2.20E+03                          | 4.66E-05                             | 2.00E-02                         | 2.33E-03        | 6.66E-06                              | NA                           | NA   |
| Phenanthrene         | 3.20E+03                          | 6.78E-05                             | NA                               | NA              | 9.69E-06                              | NA                           | NA   |
| Pyrene               | 4.40E+03                          | 9.32E-05                             | 3.00E-02                         | 3.11E-03        | 1.33E-05                              | NA                           | NA   |

NA - Not Applicable

Total Hazard Index = 7.87E-03

Total Cancer Risk = 1.39E-08



Table 35

Dermal Exposure to EUS Surface Soil (0-1') by an Adolescent Visitor (Aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                |                               |              |   |   |             |
|---|---|--------------------------------|-------------------------------|--------------|---|---|-------------|
| Cs - Concentration in soil =                                | mg/kg   | chem. spec.                    |                               |              |   |   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                                | 3052                           | calculated                    |              |   |   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                                     | 12768.3                        | USEPA 1997, EFH               |              |   |   |             |
| Fs - Fraction of skin surface area available for exposure = |   | 23.9%                          | USEPA 1997, EFH               |              |   |   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                                  | 0.026                          | USEPA 1997, EFH               |              |   |   |             |
| ABS <sub>baP</sub> - Absorption - B(a)P =                   |   | 0.03                           | USEPA 1995, Region III        |              |   |   |             |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |   | 0.1                            | USEPA 1995, Region III        |              |   |   |             |
| EF - Exposure frequency =                                   | days/year   | 12                             | reasonable assumption         |              |   |   |             |
| ED - Exposure duration =                                    | years   | 10                             | USEPA 1995, Region IV         |              |   |   |             |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                       |                               |              |   |   |             |
| BW - Body weight =  | kg  | 45                             | USEPA 1995, Region IV         |              |   |   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 3650                           | USEPA 1991, HHEM              |              |   |   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                          | USEPA 1991, HHEM              |              |   |   |             |
| Constituent   | Concentration in Soil mg/kg                         | Average Daily Intake mg/kg-day | Dermal Chronic Rf/D mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Average Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |   |                                |                               |              |   |   |             |
| Benzo(a)anthracene  | 8.35E+01  | 1.45E-07                       | NA                            | NA           | 2.07E-08                                | 7.30E-01                                  | 1.51E-08    |
| Benzo(a)pyrene  | 5.25E+01  | 9.13E-08                       | NA                            | NA           | 1.30E-08                                | 7.30E+00                                  | 9.52E-08    |
| Benzo(b)fluoranthene  | 7.95E+01  | 1.38E-07                       | NA                            | NA           | 1.98E-08                                | 7.30E-01                                  | 1.44E-08    |
| Benzo(k)fluoranthene  | 2.85E+01  | 4.96E-08                       | NA                            | NA           | 7.08E-09                                | 7.30E-02                                  | 5.17E-10    |
| Chrysene  | 8.25E+01  | 1.43E-07                       | NA                            | NA           | 2.05E-08                                | 7.30E-03                                  | 1.50E-10    |
| Dibenz(a,h)anthracene                                       | 7.45E+00  | 1.30E-08                       | NA                            | NA           | 1.85E-09                                | 7.30E+00                                  | 1.35E-08    |
| Indeno(1,2,3-cd)pyrene                                      | 3.10E+01  | 5.39E-08                       | NA                            | NA           | 7.70E-09                                | 7.30E-01                                  | 5.62E-09    |

NA - Not Available

Total Cancer Risk = 1.45E-07



**Table 36**  
**Dermal Exposure to EU1 Sediment by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u><math>\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}</math></u> |                                |                              |              |   |                                   |             |
|---|---|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| Cs - Concentration in soil =                                | mg/kg   | chem. spec.                    |                              |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day  | 3620                           | calculated                   |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>   | 20000                          | USEPA 1997, EFH              |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |   | 18.1%                          | USEPA 1997, EFH              |              |   |                                   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>  | 0.034                          | USEPA 1997, EFH              |              |   |                                   |             |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |   | 0.03                           | USEPA 1995, Region III       |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year   | 30                             | reasonable assumption        |              |   |                                   |             |
| ED - Exposure duration =                                    | years   | 25                             | USEPA 1995, Region IV        |              |   |                                   |             |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                       |                              |              |   |                                   |             |
| BW - Body weight =  | kg  | 70                             | USEPA 1995, Region IV        |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 9125                           | USEPA 1991, HHEM             |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                          | USEPA 1991, HHEM             |              |   |                                   |             |
|   |   |                                |                              |              |   |                                   |             |
| Constituent   | Concentration in Sediment mg/kg                                       | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |   |                                |                              |              |   |                                   |             |
| Benzo(a)anthracene  | 5.90E-01  | 2.56E-09                       | NA                           | NA           | 9.14E-10                                | 7.30E-01                          | 6.67E-10    |
| Benzo(a)pyrene  | 3.90E-01  | 1.69E-09                       | NA                           | NA           | 6.04E-10                                | 7.30E+00                          | 4.41E-09    |
| Benzo(b)fluoranthene  | 5.80E-01  | 2.51E-09                       | NA                           | NA           | 8.98E-10                                | 7.30E-01                          | 6.56E-10    |
| Benzo(k)fluoranthene  | 1.90E-01  | 8.24E-10                       | NA                           | NA           | 2.94E-10                                | 7.30E-02                          | 2.15E-11    |
| Chrysene  | 5.30E-01  | 2.30E-09                       | NA                           | NA           | 8.21E-10                                | 7.30E-03                          | 5.99E-12    |
| Dibenz(a,h)anthracene                                       | 6.20E-02  | 2.69E-10                       | NA                           | NA           | 9.60E-11                                | 7.30E+00                          | 7.01E-10    |
| Indeno(1,2,3-cd)pyrene                                      | 2.20E-01  | 9.54E-10                       | NA                           | NA           | 3.41E-10                                | 7.30E-01                          | 2.49E-10    |

NA - Not Available

Total Cancer Risk = 6.71E-09



Table 37

*Dermal Exposure to EUI Surface Water by a Maintenance Worker*

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | $\frac{C_w \cdot SA \cdot K_p \cdot ABS \cdot ET \cdot EF \cdot ED \cdot CF}{BW \cdot AT}$ |           |  |  |  |
|---|--|-----------|--|--|--|
| Cw - Concentration in surface water =                       | mg/L   | see below |  |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup>  | 3620      | calculated                             |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>  | 20000     | USEPA 1997, EFH                        |  |  |
| Fs - Fraction of skin surface area available for exposure = |  | 18.1%     | USEPA 1997, EFH                        |  |  |
| Kp - Dermal permeability constant =                         | cm/hr  | see below |  |  |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |  | 0.03      | USEPA 1995, Region III                 |  |  |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |  | 0.1       | USEPA 1995, Region III                 |  |  |
| ET - Exposure time =  | hrs/day  | 1         | USEPA 1992, Dermal Exposure Assessment |  |  |
| EF - Exposure frequency =                                   | days/year  | 30        | reasonable assumption                  |  |  |
| ED - Exposure duration =                                    | years  | 25        | USEPA 1995, Region IV                  |  |  |
| CF - Conversion factor =                                    | L/cm <sup>3</sup>  | 1.00E-03  |  |  |  |
| BW - Body weight =  | kg   | 70        | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days   | 9125      | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days   | 25550     | USEPA 1991, HHEM                       |  |  |

| Constituent            | Concentration in Surface Water mg/L | Kp cm/hr | Average Daily Intake | Dermal Chronic RfD | Hazard Index | Average Lifetime Daily Intake | Cancer Slope Factor | Cancer Risk |
|------------------------|-------------------------------------|----------|----------------------|--------------------|--------------|-------------------------------|---------------------|-------------|
|                        |                                     |          | mg/kg-day            | mg/kg-day          |              | mg/kg-day                     | 1/(mg/kg-day)       |             |
| <b>Semivolatiles</b>   |                                     |          |                      |                    |              |                               |                     |             |
| Benzo(a)anthracene     | 1.00E-03                            | 8.10E-01 | 1.03E-07             | NA                 | NA           | 3.69E-08                      | 7.30E-01            | 2.69E-08    |
| Benzo(a)pyrene         | 5.00E-04                            | 1.20E+00 | 7.65E-08             | NA                 | NA           | 2.73E-08                      | 7.30E+00            | 1.99E-07    |
| Benzo(b)fluoranthene   | 5.00E-04                            | 1.20E+00 | 7.65E-08             | NA                 | NA           | 2.73E-08                      | 7.30E-01            | 1.99E-08    |
| Benzo(k)fluoranthene   | 5.00E-04                            | 4.48E+01 | 2.86E-06             | NA                 | NA           | 1.02E-06                      | 7.30E-02            | 7.45E-08    |
| Chrysene               | 5.00E-04                            | 8.10E-01 | 5.16E-08             | NA                 | NA           | 1.84E-08                      | 7.30E-03            | 1.35E-10    |
| Dibenz(a,h)anthracene  | 5.00E-04                            | 2.70E+00 | 1.72E-07             | NA                 | NA           | 6.15E-08                      | 7.30E+00            | 4.49E-07    |
| Indeno(1,2,3-cd)pyrene | 5.00E-04                            | 1.90E+00 | 1.21E-07             | NA                 | NA           | 4.33E-08                      | 7.30E-01            | 3.16E-08    |

NA - Not Available

Total Cancer Risk = 8.01E-07



Table 38

Dermal Exposure to EU2 Surface Soil (0-6') by a Maintenance Worker

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                      |                            |                 |   |   |                |
|---|---|--------------------------------------|----------------------------|-----------------|---|---|----------------|
| Cs - Concentration in soil =                                | mg/kg   | chem. spec.                          |                            |                 |   |   |                |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                                | 3000                                 | calculated                 |                 |   |   |                |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                                     | 20000                                | USEPA 1997, EFH            |                 |   |   |                |
| Fs - Fraction of skin surface area available for exposure = |   | 15%                                  | USEPA 1997, EFH            |                 |   |   |                |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                                  | 0.038                                | USEPA 1997, EFH            |                 |   |   |                |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |   | 0.03                                 | USEPA 1995, Region III     |                 |   |   |                |
| EF - Exposure frequency =                                   | days/year   | 150                                  | reasonable assumption      |                 |   |   |                |
| ED - Exposure duration =                                    | years   | 25                                   | USEPA 1995, Region IV      |                 |   |   |                |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                             |                            |                 |   |   |                |
| BW - Body weight =  | kg  | 70                                   | USEPA 1995, Region IV      |                 |   |   |                |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 9125                                 | USEPA 1991, HHEM           |                 |   |   |                |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                                | USEPA 1991, HHEM           |                 |   |   |                |
|   |   |                                      |                            |                 |   |   |                |
| Constituent   | Concentration<br>in Soil<br>mg/kg                   | Average Daily<br>Intake<br>mg/kg-day | Dermal<br>RfD<br>mg/kg-day | Hazard<br>Index | Average Lifetime<br>Daily Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
| Semivolatiles   |   |                                      |                            |                 |   |   |                |
| Benzo(a)anthracene  | 2.80E+00  | 5.62E-08                             | NA                         | NA              | 2.01E-08                                      | 7.30E-01                                | 1.47E-08       |
| Benzo(a)pyrene  | 2.64E+00  | 5.30E-08                             | NA                         | NA              | 1.89E-08                                      | 7.30E+00                                | 1.38E-07       |
| Benzo(b)fluoranthene  | 9.20E+00  | 1.85E-07                             | NA                         | NA              | 6.60E-08                                      | 7.30E-01                                | 4.82E-08       |
| Benzo(k)fluoranthene  | 1.84E+00  | 3.69E-08                             | NA                         | NA              | 1.32E-08                                      | 7.30E-02                                | 9.63E-10       |
| Chrysene  | 5.33E+00  | 1.07E-07                             | NA                         | NA              | 3.82E-08                                      | 7.30E-03                                | 2.79E-10       |
| Dibenz(a,h)anthracene                                       | 2.39E-01  | 4.80E-09                             | NA                         | NA              | 1.71E-09                                      | 7.30E+00                                | 1.25E-08       |
| Indeno(1,2,3-cd)pyrene                                      | 1.97E+00  | 3.96E-08                             | NA                         | NA              | 1.41E-08                                      | 7.30E-01                                | 1.03E-08       |

NA - Not Available

Total Cancer Risk = 2.25E-07



**Table 39****Dermal Exposure to EU4 Sediment by a Maintenance Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u><math>\frac{\text{Cs} * \text{SA} * \text{AH} * \text{ABS} * \text{EF} * \text{ED} * \text{CF}}{\text{BW} * \text{AT}}</math></u> |                                |                              |              |   |                                   |             |
|---|--|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| Cs - Concentration in sediment =                            | mg/kg  | chem. spec.                    |                              |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day   | 3620                           | calculated                   |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>  | 20000                          | USEPA 1997, EFH              |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |  | 18.1%                          | USEPA 1997, EFH              |              |   |                                   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.034                          | USEPA 1997, EFH              |              |   |                                   |             |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |  | 0.03                           | USEPA 1995, Region III       |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year  | 30                             | reasonable assumption        |              |   |                                   |             |
| ED - Exposure duration =                                    | years  | 25                             | USEPA 1995, Region IV        |              |   |                                   |             |
| CF - Conversion factor =                                    | kg/mg  | 1.00E-06                       |                              |              |   |                                   |             |
| BW - Body weight =  | kg   | 70                             | USEPA 1995, Region IV        |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days   | 9125                           | USEPA 1991, HHEM             |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days   | 25550                          | USEPA 1991, HHEM             |              |   |                                   |             |
|   |  |                                |                              |              |   |                                   |             |
| Constituent   | Concentration in Sediment mg/kg  | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |  |                                |                              |              |   |                                   |             |
| Benzo(a)anthracene  | 3.30E+02   | 1.43E-06                       | NA                           | NA           | 5.11E-07                                | 7.30E-01                          | 3.73E-07    |
| Benzo(a)pyrene  | 1.30E+02   | 5.64E-07                       | NA                           | NA           | 2.01E-07                                | 7.30E+00                          | 1.47E-06    |
| Benzo(b)fluoranthene  | 1.80E+02   | 7.80E-07                       | NA                           | NA           | 2.79E-07                                | 7.30E-01                          | 2.03E-07    |
| Benzo(k)fluoranthene  | 6.40E+01   | 2.77E-07                       | NA                           | NA           | 9.91E-08                                | 7.30E-02                          | 7.23E-09    |
| Carbazole   | 5.90E+02   | 2.56E-06                       | NA                           | NA           | 9.14E-07                                | 2.00E-02                          | 1.83E-08    |
| Chrysene  | 2.90E+02   | 1.26E-06                       | NA                           | NA           | 4.49E-07                                | 7.30E-03                          | 3.28E-09    |
| Dibenz(a,h)anthracene                                       | 1.20E+01   | 5.20E-08                       | NA                           | NA           | 1.86E-08                                | 7.30E+00                          | 1.36E-07    |
| Dibenzofuran  | 9.40E+02   | 4.08E-06                       | 2.00E-03                     | 2.04E-03     | 1.46E-06                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene                                      | 4.70E+01   | 2.04E-07                       | NA                           | NA           | 7.28E-08                                | 7.30E-01                          | 5.31E-08    |
| Naphthalene   | 3.00E+03   | 1.30E-05                       | 1.00E-02                     | 1.30E-03     | 4.65E-06                                | NA                                | NA          |
| Phenanthrene  | 3.20E+03   | 1.39E-05                       | NA                           | NA           | 4.95E-06                                | NA                                | NA          |

NA - Not Available

Total Hazard Index = 3.34E-03

Total Cancer Risk = 2.26E-06



**Table 40**  
**Oral Exposure to EU4 Sediment by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |                                       | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                                  |  |
|--|---------------------------------------|---|----------------------------------|--|
| Cd - Concentration in sediment =                     | mg/kg                                 | see below   |                                  |  |
| IngR - Ingestion rate for soil =                     | mg/day                                | 100   | USEPA 1997, EFH                  |  |
| EF - Exposure frequency =                            | days/year                             | 30  | reasonable assumption            |  |
| ED - Exposure duration =                             | years                                 | 25  | USEPA 1995, Region IV            |  |
| CF - Conversion factor =                             | kg/mg                                 | 1.00E-06  |                                  |  |
| ME <sub>s</sub> - Matrix effect - PAHs =             |                                       | 0.29  | Magee, et al., 1996              |  |
| BW - Body weight =                                   | kg                                    | 70  | USEPA 1995, Region IV            |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                                  | 9125  | USEPA 1991, HHEM                 |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                                  | 25550   | USEPA 1991, HHEM                 |  |
| Constituent  | Concentration<br>in Sediment<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day  | Oral Chronic<br>RfD<br>mg/kg-day | Average                                      |
|  |                                       |   |                                  | Lifetime Daily<br>Intake<br>mg/kg-day        |
| Semivolatiles  |                                       |   | Hazard<br>Index                  | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) |
|  |                                       |   |                                  | Cancer Risk                                  |
|  |                                       |   |                                  |  |
| Carbazole  | 5.90E+02                              | 2.01E-05  | NA                               | 7.18E-06                                     |
| Dibenzofuran   | 9.40E+02                              | 3.20E-05  | 4.00E-03                         | 1.14E-05                                     |
| Naphthalene  | 3.00E+03                              | 1.02E-04  | 2.00E-02                         | 5.11E-03                                     |
| Phenanthrene   | 3.20E+03                              | 1.09E-04  | NA                               | 3.65E-05                                     |
|  |                                       |   |                                  | NA   |
|  |                                       |   |                                  | NA   |
|  |                                       |   |                                  | NA   |
| NA - Not Applicable                                  |                                       | Total Hazard Index =  | 1.31E-02                         | Total Cancer Risk =                          |
|  |                                       |   |                                  | 1.44E-07                                     |



**Table 41**  
**Dermal Exposure to EU4 Surface Water by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                   | $\frac{C_w * SA * K_p * ABS * ET * EF * ED * CF}{BW * AT}$ |  |  |  |
|---|-------------------|--|--|--|--|
| C <sub>w</sub> - Concentration in surface water =                       | mg/L              | see below  |  |  |  |
| SA - Surface area available for exposure =                              | cm <sup>2</sup>   | 3620   | calculated                             |  |  |
| SA <sub>t</sub> - Total skin surface area =                             | cm <sup>2</sup>   | 20000  | USEPA 1997, EFH                        |  |  |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = |                   | 18.1%  | USEPA 1997, EFH                        |  |  |
| K <sub>p</sub> - Dermal permeability constant =                         | cm/hr             | see below  |  |  |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                               |                   | 0.03   | USEPA 1995, Region III                 |  |  |
| ABS <sub>pah</sub> - Absorption - PAHs =                                |                   | 0.1  | USEPA 1995, Region III                 |  |  |
| ET - Exposure time =  | hrs/day           | 1  | USEPA 1992, Dermal Exposure Assessment |  |  |
| EF - Exposure frequency =   | days/year         | 30   | reasonable assumption                  |  |  |
| ED - Exposure duration =  | years             | 25   | USEPA 1995, Region IV                  |  |  |
| CF - Conversion factor =  | L/cm <sup>3</sup> | 1.00E-03   |  |  |  |
| BW - Body weight =  | kg                | 70   | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                    | days              | 9125   | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =                       | days              | 25550  | USEPA 1991, HHEM                       |  |  |

| Constituent                | Concentration in Surface Water |                      | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |          |          | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|----------------------------|--------------------------------|----------------------|--------------------------------|------------------------------|--------------|---|----------|----------|-----------------------------------|-------------|
|                            | mg/L                           | K <sub>p</sub> cm/hr |                                |                              |              | Lifetime                                | Daily    |          |                                   |             |
| <b>Semivolatiles</b>       |                                |                      |                                |                              |              |   |          |          |                                   |             |
| Benzo(a)anthracene         | 5.00E-03                       | 8.10E-01             | 5.16E-07                       | NA                           | NA           | 1.84E-07                                | 7.30E-01 | 1.35E-07 |                                   |             |
| Benzo(a)pyrene             | 5.00E-04                       | 1.20E+00             | 7.65E-08                       | NA                           | NA           | 2.73E-08                                | 7.30E+00 | 1.99E-07 |                                   |             |
| Benzo(b)fluoranthene       | 1.20E-02                       | 1.20E+00             | 1.84E-06                       | NA                           | NA           | 6.56E-07                                | 7.30E-01 | 4.79E-07 |                                   |             |
| Benzo(k)fluoranthene       | 2.00E-03                       | 4.48E+01             | 1.14E-05                       | NA                           | NA           | 4.08E-06                                | 7.30E-02 | 2.98E-07 |                                   |             |
| Bis(2-ethylhexyl)phthalate | 3.00E-03                       | 3.30E-02             | 1.26E-08                       | 1.00E-02                     | 1.26E-06     | 4.51E-09                                | 1.40E-02 | 6.31E-11 |                                   |             |
| Carbazole                  | 1.00E-02                       | 3.57E-02             | 4.56E-08                       | NA                           | NA           | 1.63E-08                                | 2.00E-02 | 3.26E-10 |                                   |             |
| Chrysene                   | 6.00E-03                       | 8.10E-01             | 6.20E-07                       | NA                           | NA           | 2.21E-07                                | 7.30E-03 | 1.62E-09 |                                   |             |
| Dibenz(a,h)anthracene      | 5.00E-04                       | 2.70E+00             | 1.72E-07                       | NA                           | NA           | 6.15E-08                                | 7.30E+00 | 4.49E-07 |                                   |             |
| Dibenzofuran               | 1.10E-02                       | 1.51E-01             | 2.12E-07                       | 2.00E-03                     | 1.06E-04     | 7.57E-08                                | NA       | NA       |                                   |             |
| Indeno(1,2,3-cd)pyrene     | 5.00E-04                       | 1.90E+00             | 1.21E-07                       | NA                           | NA           | 4.33E-08                                | 7.30E-01 | 3.16E-08 |                                   |             |
| Phenanthrene               | 1.70E-02                       | 2.30E-01             | 4.99E-07                       | NA                           | NA           | 1.78E-07                                | NA       | NA       |                                   |             |

NA - Not Available

Total Hazard Index = 1.07E-04

Total Cancer Risk = 1.59E-06



**Table 42**  
**Oral Exposure to EU4 Surface Water by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |  | $\frac{C_{sw} * IngR * EF * ED * ET}{BW * AT}$ |                                  |  |  |
|--|--|--|----------------------------------|--|--|
| Csw - Concentration in surface water =               |  | mg/L   | see below                        |  |  |
| IngR - Ingestion rate for surface water =            |  | L/hour   | 0.01                             | USEPA 1995, Region IV                  |  |
| EF - Exposure frequency =                            |  | days/year                                      | 30                               | reasonable assumption                  |  |
| ED - Exposure duration =                             |  | years  | 25                               | USEPA 1995, Region IV                  |  |
| ET - Exposure time =                                 |  | hrs/day  | 1                                | USEPA 1992, Dermal Exposure Assessment |  |
| BW - Body weight =                                   |  | kg   | 70                               | USEPA 1995, Region IV                  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = |  | days   | 9125                             | USEPA 1991, HHEM                       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    |  | days   | 25550                            | USEPA 1991, HHEM                       |  |
| Constituent  | Concentration<br>in Surface<br>Water<br>mg/L | Average<br>Daily Intake<br>mg/kg-day           | Oral Chronic<br>RfD<br>mg/kg-day | Hazard<br>Index                        | Average                                      |
|  |  |  |                                  |  | Lifetime Daily<br>Intake<br>mg/kg-day        |
| Semivolatiles  |  |  |                                  |  | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) |
|  |  |  |                                  |  | Cancer Risk                                  |
| Bis(2-ethylhexyl)phthalate                           | 3.00E-03                                     | 3.52E-08                                       | 2.00E-02                         | 1.76E-06                               | 1.26E-08                                     |
| Carbazole  | 1.00E-02                                     | 1.17E-07                                       | NA                               | NA                                     | 4.19E-08                                     |
| Dibenzofuran   | 1.10E-02                                     | 1.29E-07                                       | 4.00E-03                         | 3.23E-05                               | 4.61E-08                                     |
| Phenanthrene   | 1.70E-02                                     | 2.00E-07                                       | NA                               | NA                                     | 7.13E-08                                     |

NA - Not Applicable

Total Hazard Index = 3.41E-05

Total Cancer Risk = 1.76E-10



Table 43

*Dermal Exposure to EU4 Surface Soil (0-6') by a Maintenance Worker*

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                      | <u><math>Cs * SA * AH * ABS * EF * ED * CF</math></u> |                        |  |  |
|---|----------------------|---|------------------------|--|--|
|   |                      | BW * AT   |                        |  |  |
| $C_s$ - Concentration in soil =                             | mg/kg                | chem. spec.   |                        |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 3000  | calculated             |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000   | USEPA 1997, EPH        |  |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 15%   | USEPA 1997, EPH        |  |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.038   | USEPA 1997, EPH        |  |  |
| ABS <sub>top</sub> - Absorption - B(a)P =                   |                      | 0.03  | USEPA 1995, Region III |  |  |
| EF - Exposure frequency =                                   | days/year            | 30  | reasonable assumption  |  |  |
| ED - Exposure duration =                                    | years                | 25  | USEPA 1995, Region IV  |  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06  |                        |  |  |
| BW - Body weight =  | kg                   | 70  | USEPA 1995, Region IV  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 9125  | USEPA 1991, HHEM       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550   | USEPA 1991, HHEM       |  |  |

| Constituent            | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal          |                 | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------------|--------------------------------------|-----------------|-----------------|--|---|-------------|
|                        |                                   |                                      | Chronic<br>RFID | Hazard<br>Index |  |   |             |
| <b>Semivolatiles</b>   |                                   |                                      |                 |                 |  |   |             |
| Benzo(a)anthracene     | 9.30E+02                          | 3.73E-06                             | NA              | NA              | 1.33E-06   | 7.30E-01                                | 9.74E-07    |
| Benzo(a)pyrene         | 5.00E+02                          | 2.01E-06                             | NA              | NA              | 7.17E-07   | 7.30E+00                                | 5.23E-06    |
| Benzo(b)fluoranthene   | 5.30E+02                          | 2.13E-06                             | NA              | NA              | 7.60E-07   | 7.30E-01                                | 5.55E-07    |
| Benzo(k)fluoranthene   | 2.90E+02                          | 1.16E-06                             | NA              | NA              | 4.16E-07   | 7.30E-02                                | 3.04E-08    |
| Carbazole              | 6.20E+02                          | 2.49E-06                             | NA              | NA              | 8.89E-07   | 2.00E-02                                | 1.78E-08    |
| Chrysene               | 6.90E+02                          | 2.77E-06                             | NA              | NA              | 9.90E-07   | 7.30E-03                                | 7.22E-09    |
| Dibenz(a,h)anthracene  | 6.40E+01                          | 2.57E-07                             | NA              | NA              | 9.18E-08   | 7.30E+00                                | 6.70E-07    |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                          | 1.00E-06                             | NA              | NA              | 3.59E-07   | 7.30E-01                                | 2.62E-07    |
| Naphthalene            | 3.50E+03                          | 1.41E-05                             | 1.00E-02        | 1.41E-03        | 5.02E-06   | NA                                      | NA          |
| Phenanthrene           | 6.40E+03                          | 2.57E-05                             | NA              | NA              | 9.18E-06   | NA                                      | NA          |
| Pyrene                 | 4.40E+03                          | 1.77E-05                             | 1.50E-02        | 1.18E-03        | 6.31E-06   | NA                                      | NA          |

NA - Not Available

Total Hazard Index = 2.58E-03

Total Cancer Risk = 7.75E-06



**Table 44**  
**Oral Exposure to EU4 Surface Soil (0-6') by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| intake (mg/kg-day) =                                 |           | <u>Cd*IngR*EF*ED*CF*ME</u> |                       |  |
|--|-----------|----------------------------|-----------------------|--|
|  |           | BW*AT                      |                       |  |
| Cd - Concentration in soil =                         | mg/kg     | see below                  |                       |  |
| IngR - Ingestion rate for soil =                     | mg/day    | 100                        | USEPA 1997, EFH       |  |
| EF - Exposure frequency =                            | days/year | 30                         | reasonable assumption |  |
| ED - Exposure duration =                             | years     | 25                         | USEPA 1995, Region IV |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06                   |                       |  |
| ME <sub>s</sub> - Matrix effect - PAHs =             |           | 0.29                       | Magee, et al., 1996   |  |
| BW - Body weight =                                   | kg        | 70                         | USEPA 1995, Region IV |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 9125                       | USEPA 1991, HHEM      |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550                      | USEPA 1991, HHEM      |  |

| Constituent          | Concentration in<br>Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard Index | Average                               |  |             |
|----------------------|-----------------------------------|--------------------------------------|----------------------------------|--------------|---------------------------------------|--|-------------|
|                      |                                   |                                      |                                  |              | Lifetime Daily<br>Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b> |                                   |                                      |                                  |              |                                       |  |             |
| Carbazole            | 6.20E+02                          | 2.11E-05                             | NA                               | NA           | 7.54E-06                              | 2.00E-02                                     | 1.51E-07    |
| Naphthalene          | 3.50E+03                          | 1.19E-04                             | 2.00E-02                         | 5.96E-03     | 4.26E-05                              | NA   | NA          |
| Phenanthrene         | 6.40E+03                          | 2.18E-04                             | NA                               | NA           | 7.78E-05                              | NA   | NA          |
| Pyrene               | 4.40E+03                          | 1.50E-04                             | 3.00E-02                         | 4.99E-03     | 5.35E-05                              | NA   | NA          |

NA - Not Applicable

Total Hazard Index = 1.10E-02

Total Cancer Risk = 1.51E-07



**Table 45**

**Dermal Exposure to EUS Surface Soil (0-6') by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) = $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                   |                                      |                                    |              |  |   |                |
|--|-----------------------------------|--------------------------------------|------------------------------------|--------------|--|---|----------------|
| Cs - Concentration in soil =   | mg/kg                             | chem. spec.                          |                                    |              |  |   |                |
| SA - Surface area available for exposure =                               | cm <sup>2</sup> /day              | 3000                                 | calculated                         |              |  |   |                |
| SA <sub>t</sub> - Total skin surface area =                              | cm <sup>2</sup>                   | 20000                                | USEPA 1997, EFH                    |              |  |   |                |
| Fs - Fraction of skin surface area available for exposure =              |                                   | 15%                                  | USEPA 1997, EFH                    |              |  |   |                |
| AH - Adherence factor =  | mg/cm <sup>2</sup>                | 0.038                                | USEPA 1997, EFH                    |              |  |   |                |
| ABS <sub>bap</sub> - Absorption - B(a)P =                                |                                   | 0.03                                 | USEPA 1995, Region III             |              |  |   |                |
| EF - Exposure frequency =  | days/year                         | 150                                  | reasonable assumption              |              |  |   |                |
| ED - Exposure duration =   | years                             | 25                                   | USEPA 1995, Region IV              |              |  |   |                |
| CF - Conversion factor =   | kg/mg                             | 1.00E-06                             |                                    |              |  |   |                |
| BW - Body weight =   | kg                                | 70                                   | USEPA 1995, Region IV              |              |  |   |                |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                     | days                              | 9125                                 | USEPA 1991, HHEM                   |              |  |   |                |
| AT <sub>c</sub> - Averaging time - carcinogenic =                        | days                              | 25550                                | USEPA 1991, HHEM                   |              |  |   |                |
|  |                                   |                                      |                                    |              |  |   |                |
| Constituent  | Concentration<br>in Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Dermal<br>Chronic RfD<br>mg/kg-day | Hazard Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
|  |                                   |                                      |                                    |              |  |   |                |
| <b>Semivolatiles</b>   |                                   |                                      |                                    |              |  |   |                |
| Benzo(a)anthracene   | 7.77E+01                          | 1.56E-06                             | NA                                 | NA           | 5.57E-07   | 7.30E-01                                | 4.07E-07       |
| Benzo(a)pyrene   | 4.10E+01                          | 8.23E-07                             | NA                                 | NA           | 2.94E-07   | 7.30E+00                                | 2.15E-06       |
| Benzo(b)fluoranthene   | 7.95E+01                          | 1.60E-06                             | NA                                 | NA           | 5.70E-07   | 7.30E-01                                | 4.16E-07       |
| Benzo(k)fluoranthene   | 1.97E+01                          | 3.96E-07                             | NA                                 | NA           | 1.41E-07   | 7.30E-02                                | 1.03E-08       |
| Chrysene   | 8.25E+01                          | 1.66E-06                             | NA                                 | NA           | 5.92E-07   | 7.30E-03                                | 4.32E-09       |
| Dibenz(a,h)anthracene  | 2.04E+00                          | 4.10E-08                             | NA                                 | NA           | 1.46E-08   | 7.30E+00                                | 1.07E-07       |
| Indeno(1,2,3-cd)pyrene   | 1.71E+01                          | 3.43E-07                             | NA                                 | NA           | 1.23E-07   | 7.30E-01                                | 8.95E-08       |

NA - Not Available

Total Cancer Risk = 3.18E-06



**Table 46****Dermal Exposure to EU1 Sediment by a Construction Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                                       | <u><math>Cs * SA * AH * ABS * EF * ED * CF</math></u> |                                       |              |   |
|---|---------------------------------------|---|---------------------------------------|--------------|---|
|   |                                       | BW * AT   |                                       |              |   |
| Cs - Concentration in sediment =                            | mg/kg                                 | chem. spec.   |                                       |              |   |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                  | 3620  | calculated                            |              |   |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                       | 20000   | USEPA 1997, EFH                       |              |   |
| Fs - Fraction of skin surface area available for exposure = |                                       | 18.1%   | USEPA 1997, EFH                       |              |   |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                    | 0.105   | USEPA 1997, EFH                       |              |   |
| ABS <sub>dep</sub> - Absorption - B(a)P =                   |                                       | 0.03  | USEPA 1995, Region III                |              |   |
| EF - Exposure frequency =                                   | days/year                             | 8   | reasonable assumption                 |              |   |
| ED - Exposure duration =                                    | years                                 | 1   | reasonable assumption                 |              |   |
| CF - Conversion factor =                                    | kg/mg                                 | 1.00E-06  |                                       |              |   |
| BW - Body weight =  | kg                                    | 70  | USEPA 1995, Region IV                 |              |   |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                                  | 365   | USEPA 1991, HHEM                      |              |   |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                                  | 25550   | USEPA 1991, HHEM                      |              |   |
| Constituent   | Concentration in<br>Sediment<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day                  | Dermal<br>Subchronic RfD<br>mg/kg-day | Hazard Index | Average                                     |
|   |                                       |   |                                       |              | Lifetime Daily<br>Intake<br>mg/kg-day       |
| Semivolatiles   |                                       |   |                                       |              | Cancer<br>Slope Factor<br>1/(mg/kg-<br>day) |
|   |                                       |   |                                       |              | Cancer<br>Risk                              |
|   |                                       |   |                                       |              |   |
| Benzo(a)anthracene  | 5.90E-01                              | 2.11E-09  | NA                                    | NA           | 3.01E-11                                    |
| Benzo(a)pyrene  | 3.90E-01                              | 1.39E-09  | NA                                    | NA           | 1.99E-11                                    |
| Benzo(b)fluoranthene  | 5.80E-01                              | 2.07E-09  | NA                                    | NA           | 2.96E-11                                    |
| Benzo(k)fluoranthene  | 1.90E-01                              | 6.78E-10  | NA                                    | NA           | 9.69E-12                                    |
| Chrysene  | 5.30E-01                              | 1.89E-09  | NA                                    | NA           | 2.70E-11                                    |
| Dibenz(a,h)anthracene                                       | 6.20E-02                              | 2.21E-10  | NA                                    | NA           | 3.16E-12                                    |
| Indeno(1,2,3-cd)pyrene                                      | 2.20E-01                              | 7.85E-10  | NA                                    | NA           | 1.12E-11                                    |

NA - Not Available

Total Cancer Risk = 2.21E-10



**Table 47****Dermal Exposure to EUI Surface Water by a Construction Worker****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                                     | $\frac{C_w * SA * K_p * ABS * ET * EF * ED * CF}{BW * AT}$ |  |                                 |              |   |                                   |             |
|---|-------------------------------------|--|--|---------------------------------|--------------|---|-----------------------------------|-------------|
| Cw - Concentration in surface water =                                   |                                     | mg/L   | see below                              |                                 |              |   |                                   |             |
| SA - Surface area available for exposure =                              |                                     | cm <sup>2</sup>  | 3620                                   | calculated                      |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                             |                                     | cm <sup>2</sup>  | 20000                                  | USEPA 1997, EFH                 |              |   |                                   |             |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = |                                     |  | 18.1%                                  | USEPA 1997, EFH                 |              |   |                                   |             |
| K <sub>p</sub> - Dermal permeability constant =                         |                                     | cm/hr  | see below                              |                                 |              |   |                                   |             |
| ABS <sub>bap</sub> - Absorption - B(a)P =                               |                                     |  | 0.03                                   | USEPA 1995, Region III          |              |   |                                   |             |
| ABS <sub>pah</sub> - Absorption - PAHs =                                |                                     |  | 0.1                                    | USEPA 1995, Region III          |              |   |                                   |             |
| ET - Exposure time =  | hrs/day                             | 1  | USEPA 1992, Dermal Exposure Assessment |                                 |              |   |                                   |             |
| EF - Exposure frequency =   | days/year                           | 8  | reasonable assumption                  |                                 |              |   |                                   |             |
| ED - Exposure duration =  | years                               | 1  | reasonable assumption                  |                                 |              |   |                                   |             |
| CF - Conversion factor =  | L/cm <sup>3</sup>                   | 1.00E-03   |  |                                 |              |   |                                   |             |
| BW - Body weight =  | kg                                  | 70   | USEPA 1995, Region IV                  |                                 |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                    | days                                | 365  | USEPA 1991, HHEM                       |                                 |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =                       | days                                | 25550  | USEPA 1991, HHEM                       |                                 |              |   |                                   |             |
| Constituent   | Concentration in Surface Water mg/L | K <sub>p</sub> cm/hr                                       | Average Daily Intake mg/kg-day         | Dermal Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>  |                                     |  |  |                                 |              |   |                                   |             |
| Benzo(a)anthracene  | 1.00E-03                            | 8.10E-01   | 2.75E-08                               | NA                              | NA           | 3.93E-10                                | 7.30E-01                          | 2.87E-10    |
| Benzo(a)pyrene  | 5.00E-04                            | 1.20E+00   | 2.04E-08                               | NA                              | NA           | 2.91E-10                                | 7.30E+00                          | 2.13E-09    |
| Benzo(b)fluoranthene  | 5.00E-04                            | 1.20E+00   | 2.04E-08                               | NA                              | NA           | 2.91E-10                                | 7.30E-01                          | 2.13E-10    |
| Benzo(k)fluoranthene  | 5.00E-04                            | 4.48E+01   | 7.62E-07                               | NA                              | NA           | 1.09E-08                                | 7.30E-02                          | 7.94E-10    |
| Chrysene  | 5.00E-04                            | 8.10E-01   | 1.38E-08                               | NA                              | NA           | 1.97E-10                                | 7.30E-03                          | 1.44E-12    |
| Dibenz(a,h)anthracene   | 5.00E-04                            | 2.70E+00   | 4.59E-08                               | NA                              | NA           | 6.56E-10                                | 7.30E+00                          | 4.79E-09    |
| Indeno(1,2,3-cd)pyrene  | 5.00E-04                            | 1.90E+00   | 3.23E-08                               | NA                              | NA           | 4.61E-10                                | 7.30E-01                          | 3.37E-10    |

NA - Not Available

Total Cancer Risk = 8.55E-09



**Table 48**  
**Dermal Exposure to EU2 Soil (0-10') by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |                        |  |
|---|----------------------|------------------------------|------------------------|--|
|   |                      | BW*AT                        |                        |  |
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.                  |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 5560                         | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000                        | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 27.8%                        | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.1                          | USEPA 1997, EFH        |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                      | 0.03                         | USEPA 1995, Region III |  |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |                      | 0.1                          | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year            | 80                           | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                | 1                            | reasonable assumption  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06                     |                        |  |
| BW - Body weight =  | kg                   | 70                           | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 365                          | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550                        | USEPA 1991, HHEM       |  |

| Constituent            | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal                         |                 | Average                               |   |                |
|------------------------|-----------------------------------|--------------------------------------|--------------------------------|-----------------|---------------------------------------|---|----------------|
|                        |                                   |                                      | Subchronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
| <b>Semivolatiles</b>   |                                   |                                      |                                |                 |                                       |   |                |
| Benzo(a)anthracene     | 2.80E+00                          | 1.46E-07                             | NA                             | NA              | 2.09E-09                              | 7.30E-01                                | 1.53E-09       |
| Benzo(a)pyrene         | 2.64E+00                          | 1.38E-07                             | NA                             | NA              | 1.97E-09                              | 7.30E+00                                | 1.44E-08       |
| Benzo(b)fluoranthene   | 9.20E+00                          | 4.80E-07                             | NA                             | NA              | 6.86E-09                              | 7.30E-01                                | 5.01E-09       |
| Benzo(k)fluoranthene   | 1.84E+00                          | 9.61E-08                             | NA                             | NA              | 1.37E-09                              | 7.30E-02                                | 1.00E-10       |
| Chrysene               | 5.33E+00                          | 2.78E-07                             | NA                             | NA              | 3.98E-09                              | 7.30E-03                                | 2.90E-11       |
| Dibenz(a,h)anthracene  | 2.39E-01                          | 1.25E-08                             | NA                             | NA              | 1.78E-10                              | 7.30E+00                                | 1.30E-09       |
| Indeno(1,2,3-cd)pyrene | 1.97E+00                          | 1.03E-07                             | NA                             | NA              | 1.47E-09                              | 7.30E-01                                | 1.07E-09       |

NA - Not Available

Total Cancer Risk = 2.34E-08





**Table 49**  
**Exposure to Construction Workers from Inhalation of Fugitive Dust in EU2**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) = $C_a * InhR * EF * ED * RF$   |                             |                      |  |                                |               |              |                                |                              |             |
|--|-----------------------------|----------------------|--|--------------------------------|---------------|--------------|--------------------------------|------------------------------|-------------|
| $BW * AT$<br>$C_a$ - Concentration in air = mg/m <sup>3</sup> chem.spec.<br>$InhR$ - Inhalation Rate = m <sup>3</sup> /shift 20<br>$EF$ - Exposure Frequency = shifts/year 80<br>$ED$ - Exposure Duration = years 1<br>$RF_s$ - Retention Factor - semivolatiles = 0.75<br>$AT_n$ - Averaging Time noncarcinogenic = days 365<br>$AT_c$ - Averaging Time carcinogenic = days 25550<br>$BW$ - Body Weight = kg 70<br>$E_i$ - Emission Rate (mg/sec) = $C_s * (PER_v + PER_e)$<br>$C_s$ - Concentration in soil = mg/kg chem.spec. |                             |                      |  |                                |               |              |                                |                              |             |
| Chemicals  | Concentration in Soil mg/kg | Emission Rate mg/sec | Concentration in Air mg/m <sup>3</sup> | Average Daily Intake mg/kg-day | Inhalation    |              | Average Daily Intake mg/kg-day | Inhalation                   |             |
|  |                             |                      |  |                                | Subchronic RD | Hazard Index |                                | Lifetime Cancer Slope Factor | Cancer Risk |
| <b>Semivolatiles</b>   |                             |                      |  |                                |               |              |                                |                              |             |
| Benzo(a)anthracene   | 2.80E+00                    | 3.06E-03             | 2.72E-06                               | 1.28E-07                       | NA            | NA           | 1.82E-09                       | 3.10E-01                     | 5.65E-10    |
| Benzo(a)pyrene   | 2.64E+00                    | 2.89E-03             | 2.56E-06                               | 1.20E-07                       | NA            | NA           | 1.72E-09                       | 3.10E+00                     | 5.33E-09    |
| Benzo(b)fluoranthene   | 9.20E+00                    | 1.01E-02             | 8.93E-06                               | 4.19E-07                       | NA            | NA           | 5.99E-09                       | 3.10E-01                     | 1.86E-09    |
| Benzo(k)fluoranthene   | 1.84E+00                    | 2.01E-03             | 1.79E-06                               | 8.39E-08                       | NA            | NA           | 1.20E-09                       | 3.10E-02                     | 3.71E-11    |
| Chrysene   | 5.33E+00                    | 5.83E-03             | 5.17E-06                               | 2.43E-07                       | NA            | NA           | 3.47E-09                       | 3.10E-03                     | 1.08E-11    |
| Dibenz(a,h)anthracene  | 2.39E-01                    | 2.61E-04             | 2.32E-07                               | 1.09E-08                       | NA            | NA           | 1.56E-10                       | 3.10E+00                     | 4.82E-10    |
| Indeno(1,2,3-cd)pyrene   | 1.97E+00                    | 2.15E-03             | 1.91E-06                               | 8.98E-08                       | NA            | NA           | 1.28E-09                       | 3.10E-01                     | 3.98E-10    |
| NA - Not Available   |                             |                      |  |                                |               |              |                                |                              |             |
| Total Cancer Risk:   |                             |                      |  |                                |               |              |                                |                              |             |
| 8.11E-09   |                             |                      |  |                                |               |              |                                |                              |             |

**Table 50****Dermal Exposure to EU4 Sediment by a Construction Worker**

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u><math>\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}</math></u> |                          |                        |                             |                            |             |          |
|---|---|--------------------------|------------------------|-----------------------------|----------------------------|-------------|----------|
| Cs - Concentration in sediment =                            | mg/kg   | chem. spec.              |                        |                             |                            |             |          |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day  | 3620                     | calculated             |                             |                            |             |          |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>   | 20000                    | USEPA 1997, EFH        |                             |                            |             |          |
| Fs - Fraction of skin surface area available for exposure = |   | 18.1%                    | USEPA 1997, EFH        |                             |                            |             |          |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>  | 0.105                    | USEPA 1997, EFH        |                             |                            |             |          |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |   | 0.03                     | USEPA 1995, Region III |                             |                            |             |          |
| EF - Exposure frequency =                                   | days/year   | 8                        | reasonable assumption  |                             |                            |             |          |
| ED - Exposure duration =                                    | years   | 1                        | reasonable assumption  |                             |                            |             |          |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                 |                        |                             |                            |             |          |
| BW - Body weight =  | kg  | 70                       | USEPA 1995, Region IV  |                             |                            |             |          |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 365                      | USEPA 1991, HHEM       |                             |                            |             |          |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                    | USEPA 1991, HHEM       |                             |                            |             |          |
| Dermal  |   |                          |                        |                             |                            |             |          |
| Constituent   | Average Daily Intake  | Subchronic RID mg/kg-day | Lifetime Daily Intake  | Average Cancer Slope Factor | Cancer Slope 1/(mg/kg-day) | Cancer Risk |          |
| Semivolatiles   | Concentration in Sediment mg/kg                                       |                          | Hazard Index           |                             |                            |             |          |
| Benzo(a)anthracene  | 3.30E+02  | 1.18E-06                 | NA                     | NA                          | 1.68E-08                   | 7.30E-01    | 1.23E-08 |
| Benzo(a)pyrene  | 1.30E+02  | 4.64E-07                 | NA                     | NA                          | 6.63E-09                   | 7.30E+00    | 4.84E-08 |
| Benzo(b)fluoranthene  | 1.80E+02  | 6.43E-07                 | NA                     | NA                          | 9.18E-09                   | 7.30E-01    | 6.70E-09 |
| Benzo(k)fluoranthene  | 6.40E+01  | 2.29E-07                 | NA                     | NA                          | 3.26E-09                   | 7.30E-02    | 2.38E-10 |
| Carbazole   | 5.90E+02  | 2.11E-06                 | NA                     | NA                          | 3.01E-08                   | 2.00E-02    | 6.02E-10 |
| Chrysene  | 2.90E+02  | 1.04E-06                 | NA                     | NA                          | 1.48E-08                   | 7.30E-03    | 1.08E-10 |
| Dibenz(a,h)anthracene                                       | 1.20E+01  | 4.28E-08                 | NA                     | NA                          | 6.12E-10                   | 7.30E+00    | 4.47E-09 |
| Dibenzofuran  | 9.40E+02  | 3.36E-06                 | NA                     | NA                          | 4.79E-08                   | NA          | NA       |
| Indeno(1,2,3-cd)pyrene                                      | 4.70E+01  | 1.68E-07                 | NA                     | NA                          | 2.40E-09                   | 7.30E-01    | 1.75E-09 |
| Naphthalene   | 3.00E+03  | 1.07E-05                 | NA                     | NA                          | 1.53E-07                   | NA          | NA       |
| Phenanthrene  | 3.20E+03  | 1.14E-05                 | NA                     | NA                          | 1.63E-07                   | NA          | NA       |

NA - Not Available

Total Cancer Risk = 7.46E-08



**Table 51**  
**Oral Exposure to EU4 Sediment by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |                                       | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                           |              |  |
|--|---------------------------------------|---|---------------------------|--------------|--|
| Cd - Concentration in sediment =                     | mg/kg                                 | see below   |                           |              |  |
| IngR - Ingestion rate for sediment =                 | mg/day                                | 100   | USEPA 1997, EFH           |              |  |
| EF - Exposure frequency =                            | days/year                             | 8   | reasonable assumption     |              |  |
| ED - Exposure duration =                             | years                                 | 1   | USEPA 1995, Region IV     |              |  |
| CF - Conversion factor =                             | kg/mg                                 | 1.00E-06  |                           |              |  |
| ME <sub>s</sub> - Matrix effect - PAHs =             |                                       | 0.29  | Magee, et al., 1996       |              |  |
| BW - Body weight =                                   | kg                                    | 70  | USEPA 1995, Region IV     |              |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                                  | 365   | USEPA 1991, HHEM          |              |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                                  | 25550   | USEPA 1991, HHEM          |              |  |
| Constituent  | Concentration<br>in Sediment<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day  | Oral<br>Subchronic<br>RfD |              | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day |
|  |                                       |   | mg/kg-day                 | Hazard Index | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day)     |
| <b>Semivolatiles</b>                                 |                                       |   |                           |              |  |
| Carbazole  | 5.90E+02                              | 5.36E-06  | NA                        | NA           | 7.65E-08   |
| Dibenzofuran   | 9.40E+02                              | 8.54E-06  | NA                        | NA           | 1.22E-07   |
| Naphthalene  | 3.00E+03                              | 2.72E-05  | NA                        | NA           | 3.89E-07   |
| Phenanthrene   | 3.20E+03                              | 2.91E-05  | NA                        | NA           | 4.15E-07   |

NA - Not Applicable

Total Cancer Risk = 1.53E-09



**Table 52**

**Dermal Exposure to EU4 Surface Water by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                   | <u>Cw*SA*Kp*ABS*ET*EF*ED*CF</u> |  |  |  |
|---|-------------------|---------------------------------|--|--|--|
|   |                   | <u>BW*AT</u>                    |  |  |  |
| Cw - Concentration in surface water =                                   | mg/L              | see below                       |  |  |  |
| SA - Surface area available for exposure =                              | cm <sup>2</sup>   | 3620                            | calculated                             |  |  |
| SA <sub>t</sub> - Total skin surface area =                             | cm <sup>2</sup>   | 20000                           | USEPA 1997, EFH                        |  |  |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = |                   | 18.1%                           | USEPA 1997, EFH                        |  |  |
| K <sub>p</sub> - Dermal permeability constant =                         | cm/hr             | see below                       |  |  |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                               |                   | 0.03                            | USEPA 1995, Region III                 |  |  |
| ABS <sub>pbh</sub> - Absorption - PAHs =                                |                   | 0.1                             | USEPA 1995, Region III                 |  |  |
| ET - Exposure time =  | hrs/day           | 1                               | USEPA 1992, Dermal Exposure Assessment |  |  |
| EF - Exposure frequency =   | days/year         | 8                               | reasonable assumption                  |  |  |
| ED - Exposure duration =  | years             | 1                               | reasonable assumption                  |  |  |
| CF - Conversion factor =  | L/cm <sup>3</sup> | 1.00E-03                        |  |  |  |
| BW - Body weight =  | kg                | 70                              | USEPA 1995, Region IV                  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                    | days              | 365                             | USEPA 1991, HHEM                       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =                       | days              | 25550                           | USEPA 1991, HHEM                       |  |  |

| Constituent                | Concentration in Surface Water |                      | Average Daily Intake mg/kg-day | Dermal Subchronic RfD mg/kg-day |          | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|----------------------------|--------------------------------|----------------------|--------------------------------|---------------------------------|----------|--------------|---|-----------------------------------|-------------|
|                            | Water mg/L                     | K <sub>p</sub> cm/hr |                                |                                 |          |              |   |                                   |             |
| <b>Semivolatiles</b>       |                                |                      |                                |                                 |          |              |   |                                   |             |
| Benzo(a)anthracene         | 5.00E-03                       | 8.10E-01             | 1.38E-07                       | NA                              | NA       | NA           | 1.97E-09                                | 7.30E-01                          | 1.44E-09    |
| Benzo(a)pyrene             | 5.00E-04                       | 1.20E+00             | 2.04E-08                       | NA                              | NA       | NA           | 2.91E-10                                | 7.30E+00                          | 2.13E-09    |
| Benzo(b)fluoranthene       | 1.20E-02                       | 1.20E+00             | 4.90E-07                       | NA                              | NA       | NA           | 7.00E-09                                | 7.30E-01                          | 5.11E-09    |
| Benzo(k)fluoranthene       | 2.00E-03                       | 4.48E+01             | 3.04E-06                       | NA                              | NA       | NA           | 4.35E-08                                | 7.30E-02                          | 3.18E-09    |
| Bis(2-ethylhexyl)phthalate | 3.00E-03                       | 3.30E-02             | 3.37E-09                       | 1.00E-02                        | 3.37E-07 | 3.37E-07     | 4.81E-11                                | 1.40E-02                          | 6.73E-13    |
| Carbazole                  | 1.00E-02                       | 3.57E-02             | 1.22E-08                       | NA                              | NA       | NA           | 1.74E-10                                | 2.00E-02                          | 3.47E-12    |
| Chrysene                   | 6.00E-03                       | 8.10E-01             | 1.65E-07                       | NA                              | NA       | NA           | 2.36E-09                                | 7.30E-03                          | 1.72E-11    |
| Dibenz(a,h)anthracene      | 5.00E-04                       | 2.70E+00             | 4.59E-08                       | NA                              | NA       | NA           | 6.56E-10                                | 7.30E+00                          | 4.79E-09    |
| Dibenzofuran               | 1.10E-02                       | 1.51E-01             | 5.65E-08                       | NA                              | NA       | NA           | 8.07E-10                                | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene     | 5.00E-04                       | 1.90E+00             | 3.23E-08                       | NA                              | NA       | NA           | 4.61E-10                                | 7.30E-01                          | 3.37E-10    |
| Phenanthrene               | 1.70E-02                       | 2.30E-01             | 1.33E-07                       | NA                              | NA       | NA           | 1.90E-09                                | NA                                | NA          |

NA - Not Available

Total Hazard Index = 3.37E-07

Total Cancer Risk = 1.70E-08



**Table 53**

**Oral Exposure to EU4 Surface Water by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | $\frac{C_{sw} * IngR * EF * ED * ET}{BW * AT}$ |  |                                  |
|--|-----------|--|--|----------------------------------|
| C <sub>sw</sub> - Concentration in surface water =   | mg/L      | see below                                      |  |                                  |
| IngR - Ingestion rate for surface water =            | L/hour    | 0.01   | USEPA 1995, Region IV                  |                                  |
| EF - Exposure frequency =                            | days/year | 8  | reasonable assumption                  |                                  |
| ED - Exposure duration =                             | years     | 1  | USEPA 1995, Region IV                  |                                  |
| ET - Exposure time =                                 | hrs/day   | 1  | USEPA 1992, Dermal Exposure Assessment |                                  |
| BW - Body weight =                                   | kg        | 70   | USEPA 1995, Region IV                  |                                  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 365  | USEPA 1991, HHEM                       |                                  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550  | USEPA 1991, HHEM                       |                                  |
| Concentration<br>in Surface<br>Water                 |           |  |  |                                  |
| Constituent  | mg/L      | Average<br>Daily Intake<br>mg/kg-day           | Oral<br>Subchronic RfD<br>mg/kg-day    | Average Lifetime<br>Hazard Index |
| Semivolatiles  |           |  |  |                                  |
| Bis(2-ethylhexyl)phthalate                           | 3.00E-03  | 9.39E-09                                       | 1.00E-02                               | 9.39E-07                         |
| Carbazole  | 1.00E-02  | 3.13E-08                                       | NA                                     | NA                               |
| Dibenzofuran   | 1.10E-02  | 3.44E-08                                       | NA                                     | NA                               |
| Phenanthrene   | 1.70E-02  | 5.32E-08                                       | NA                                     | NA                               |

NA - Not Applicable

Total Hazard Index = 9.39E-07

Total Cancer Risk = 1.08E-11



**Table 54**  
**Dermal Exposure to EU4 Soil (0-20') by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |                        |  |
|---|----------------------|------------------------------|------------------------|--|
|   |                      | BW*AT                        |                        |  |
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.                  |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 5560                         | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000                        | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 27.8%                        | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.1                          | USEPA 1997, EFH        |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                      | 0.03                         | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year            | 8                            | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                | 1                            | reasonable assumption  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06                     |                        |  |
| BW - Body weight =  | kg                   | 70                           | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 365                          | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550                        | USEPA 1991, HHEM       |  |

| Constituent            | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day | Dermal                   |              | Average                         |                                   |             |
|------------------------|-----------------------------|--------------------------------|--------------------------|--------------|---------------------------------|-----------------------------------|-------------|
|                        |                             |                                | Subchronic RfD mg/kg-day | Hazard Index | Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>   |                             |                                |                          |              |                                 |                                   |             |
| Benzo(a)anthracene     | 9.30E+02                    | 4.86E-06                       | NA                       | NA           | 6.94E-08                        | 7.30E-01                          | 5.07E-08    |
| Benzo(a)pyrene         | 5.00E+02                    | 2.61E-06                       | NA                       | NA           | 3.73E-08                        | 7.30E+00                          | 2.72E-07    |
| Benzo(b)fluoranthene   | 5.30E+02                    | 2.77E-06                       | NA                       | NA           | 3.95E-08                        | 7.30E-01                          | 2.89E-08    |
| Benzo(k)fluoranthene   | 2.90E+02                    | 1.51E-06                       | NA                       | NA           | 2.16E-08                        | 7.30E-02                          | 1.58E-09    |
| Carbazole              | 6.20E+02                    | 3.24E-06                       | NA                       | NA           | 4.63E-08                        | 2.00E-02                          | 9.25E-10    |
| Chrysene               | 6.90E+02                    | 3.60E-06                       | NA                       | NA           | 5.15E-08                        | 7.30E-03                          | 3.76E-10    |
| Dibenz(a,h)anthracene  | 6.40E+01                    | 3.34E-07                       | NA                       | NA           | 4.78E-09                        | 7.30E+00                          | 3.49E-08    |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                    | 1.31E-06                       | NA                       | NA           | 1.87E-08                        | 7.30E-01                          | 1.36E-08    |
| Naphthalene            | 3.50E+03                    | 1.83E-05                       | NA                       | NA           | 2.61E-07                        | NA                                | NA          |
| Phenanthrene           | 6.40E+03                    | 3.34E-05                       | NA                       | NA           | 4.78E-07                        | NA                                | NA          |
| Pyrene                 | 4.40E+03                    | 2.30E-05                       | 1.50E-01                 | 1.53E-04     | 3.28E-07                        | NA                                | NA          |

NA - Not Available

Total Hazard Index = 1.53E-04

Total Cancer Risk = 4.03E-07

**Table 55**  
**Oral Exposure to EU4 Soil (0-20') by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                                |                               |              |   |  |             |
|--|---|--------------------------------|-------------------------------|--------------|---|--|-------------|
| Cd - Concentration in soil =                         | mg/kg   | see below                      |                               |              |   |  |             |
| IngR - Ingestion rate for soil =                     | mg/day  | 100                            | USEPA 1997, EFH               |              |   |  |             |
| EF - Exposure frequency =                            | days/year   | 8                              | reasonable assumption         |              |   |  |             |
| ED - Exposure duration =                             | years   | 1                              | USEPA 1995, Region IV         |              |   |  |             |
| CF - Conversion factor =                             | kg/mg   | 1.00E-06                       |                               |              |   |  |             |
| ME <sub>s</sub> - Matrix effect - PAHs =             |   | 0.29                           | Magee, et al., 1996           |              |   |  |             |
| BW - Body weight =                                   | kg  | 70                             | USEPA 1995, Region IV         |              |   |  |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days  | 365                            | USEPA 1991, HHEM              |              |   |  |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days  | 25550                          | USEPA 1991, HHEM              |              |   |  |             |
|  |   |                                |                               |              |   |  |             |
| Constituent  | Concentration in Soil mg/kg   | Average Daily Intake mg/kg-day | Oral Subchronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles  |   |                                |                               |              |   |  |             |
| Carbazole  | 6.20E+02  | 5.63E-06                       | NA                            | NA           | 8.04E-08                                | 2.00E-02                               | 1.61E-09    |
| Naphthalene  | 3.50E+03  | 3.18E-05                       | NA                            | NA           | 4.54E-07                                | NA                                     | NA          |
| Phenanthrene   | 6.40E+03  | 5.81E-05                       | NA                            | NA           | 8.30E-07                                | NA                                     | NA          |
| Pyrene   | 4.40E+03  | 4.00E-05                       | 3.00E-01                      | 1.33E-04     | 5.71E-07                                | NA                                     | NA          |
| NA - Not Applicable                                  |   | Total Hazard Index = 1.33E-04  |                               |              | Total Cancer Risk = 1.61E-09            |  |             |



**Table 56**  
**Exposure to Construction Workers from Inhalation of Fugitive Dust in EU4**  
**Kerr McGee, Hattiesburg, MS**

| Chemicals              | Concentration in<br>Soil<br>mg/kg | Emission<br>Rate<br>mg/sec | Concentration in<br>Air<br>mg/m <sup>3</sup> | Average Daily Intake<br>mg/kg-day | Inhalation<br>Subchronic RD<br>mg/kg-day | Hazard<br>Index | Average Lifetime<br>Daily Intake<br>mg/kg-day | Inhalation<br>Cancer Slope<br>Factor | Cancer<br>Risk |
|------------------------|-----------------------------------|----------------------------|--|-----------------------------------|--|-----------------|---|--------------------------------------|----------------|
| <b>Semivolatiles</b>   |                                   |                            |  |                                   |  |                 |   |                                      |                |
| Benz(a)anthracene      | 9.30E+02                          | 1.02E+00                   | 9.02E-04                                     | 4.24E-06                          | NA                                       | NA              | 6.05E-08                                      | 3.10E-01                             | 1.88E-08       |
| Benz(a)pyrene          | 5.00E+02                          | 5.47E-01                   | 4.85E-04                                     | 2.28E-06                          | NA                                       | NA              | 3.26E-08                                      | 3.10E+00                             | 1.01E-07       |
| Benz(b)fluoranthene    | 5.30E+02                          | 5.79E-01                   | 5.14E-04                                     | 2.42E-06                          | NA                                       | NA              | 3.45E-08                                      | 3.10E-01                             | 1.07E-08       |
| Benz(k)fluoranthene    | 2.90E+02                          | 3.17E-01                   | 2.81E-04                                     | 1.32E-06                          | NA                                       | NA              | 1.89E-08                                      | 3.10E-02                             | 5.84E-10       |
| Carbazole              | 6.20E+02                          | 6.78E-01                   | 6.02E-04                                     | 2.83E-06                          | NA                                       | NA              | 4.04E-08                                      | NA                                   | NA             |
| Chrysene               | 6.90E+02                          | 7.54E-01                   | 6.70E-04                                     | 3.14E-06                          | NA                                       | NA              | 4.49E-08                                      | 3.10E-03                             | 1.39E-10       |
| Dibenz(a,h)anthracene  | 6.40E+01                          | 7.00E-02                   | 6.21E-05                                     | 2.92E-07                          | NA                                       | NA              | 4.17E-09                                      | 3.10E+00                             | 1.29E-08       |
| Indeno(1,2,3-cd)pyrene | 2.50E+02                          | 2.73E-01                   | 2.43E-04                                     | 1.14E-06                          | NA                                       | NA              | 1.63E-08                                      | 3.10E-01                             | 5.05E-09       |
| Naphthalene            | 3.50E+03                          | 3.83E+00                   | 3.40E-03                                     | 1.60E-05                          | NA                                       | NA              | 2.28E-07                                      | NA                                   | NA             |
| Phenanthrene           | 6.40E+03                          | 7.00E+00                   | 6.21E-03                                     | 2.92E-05                          | NA                                       | NA              | 4.17E-07                                      | NA                                   | NA             |
| Pyrene                 | 4.40E+03                          | 4.81E+00                   | 4.27E-03                                     | 2.01E-05                          | NA                                       | NA              | 2.86E-07                                      | NA                                   | NA             |
| NA - Not Available     |                                   |                            |  |                                   |  |                 |   |                                      |                |
| Total Cancer Risk:     |                                   |                            |  |                                   |  |                 |   |                                      | 1.49E-07       |

$C_a = \frac{C_a * InhR * EF * ED * RF}{BW * AT}$   
 $E_i = \frac{E_i / (Hb * W * V)}{Hb - Downwind Hit (m) = 4.81}$   
 $W - Width (m) = 50$   
 $V - Wind speed (m/sec) = 4.69$   
 $Length (downwind distance) (m) = 50$   
 $r - Roughness Ht. (m) = 0.20$   
 $z - downwind distance (m) = 50$   
 $z = 6.25[r(Hb/r + Ln(Hb/r) - 1.58)]$

$E_i - Emission Rate (mg/sec) = C_s * (PERv + PERe)$   
 $C_s - Concentration in soil = mg/kg see below$



**Table 57**

**Dermal Exposure to EUS Soil (0-20') by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                      | <u>Cs*SA*AH*ABS*EF*ED*CF</u>      |                                      |  |  |
|---|----------------------|-----------------------------------|--------------------------------------|--|--|
|   |                      | <u>BW*AT</u>                      |                                      |  |  |
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.                       |                                      |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 5560                              | calculated                           |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000                             | USEPA 1997, EFH                      |  |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 27.8%                             | USEPA 1997, EFH                      |  |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.1                               | USEPA 1997, EFH                      |  |  |
| ABS <sub>top</sub> - Absorption - B(a)P =                   |                      | 0.03                              | USEPA 1995, Region III               |  |  |
| EF - Exposure frequency =                                   | days/year            | 80                                | reasonable assumption                |  |  |
| ED - Exposure duration =                                    | years                | 1                                 | reasonable assumption                |  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06                          |                                      |  |  |
| BW - Body weight =  | kg                   | 70                                | USEPA 1995, Region IV                |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 365                               | USEPA 1991, HHEM                     |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550                             | USEPA 1991, HHEM                     |  |  |
| Constituent   |                      | Concentration in<br>Soil<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Dermal<br>Subchronic<br>RfD<br>mg/kg-day | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day |
|   |                      |                                   |                                      | Hazard Index                             | Cancer Slope<br>Factor<br>1/(mg/kg-day)          |
| Constituent   |                      |                                   |                                      |  | Cancer<br>Risk                                   |
| <b>Semivolatiles</b>  |                      |                                   |                                      |  |  |
| Benzo(a)anthracene  | 3.39E+01             | 1.77E-06                          | NA                                   | NA                                       | 2.53E-08   |
| Benzo(a)pyrene  | 1.71E+01             | 8.93E-07                          | NA                                   | NA                                       | 1.28E-08   |
| Benzo(b)fluoranthene  | 4.67E+01             | 2.44E-06                          | NA                                   | NA                                       | 3.48E-08   |
| Benzo(k)fluoranthene  | 8.05E+00             | 4.20E-07                          | NA                                   | NA                                       | 6.01E-09   |
| Chrysene  | 5.02E+01             | 2.62E-06                          | NA                                   | NA                                       | 3.75E-08   |
| Dibenz(a,h)anthracene                                       | 1.04E+00             | 5.43E-08                          | NA                                   | NA                                       | 7.76E-10   |
| Indeno(1,2,3-cd)pyrene                                      | 7.55E+00             | 3.94E-07                          | NA                                   | NA                                       | 5.63E-09   |

NA - Not Available

Total Cancer Risk = 1.48E-07



**Table 58**  
**Exposure to Construction Workers from Inhalation of Fugitive Dust in EUS**  
**Kerr McGee, Hattiesburg, MS**

| Chemicals              | Concentration in Soil mg/kg | Emission Rate mg/sec | Concentration in Air mg/m <sup>3</sup> | Average Daily Intake mg/kg-day | Inhalation RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Inhalation Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|-----------------------------|----------------------|--|--------------------------------|--------------------------|--------------|---|--|-------------|
| <b>Semivolatiles</b>   |                             |                      |  |                                |                          |              |   |  |             |
| Benz(a)anthracene      | 3.39E+01                    | 3.71E-02             | 3.29E-05                               | 1.54E-06                       | NA                       | NA           | 2.21E-08                                | 3.10E-01                                     | 6.84E-09    |
| Benz(a)pyrene          | 1.71E+01                    | 1.87E-02             | 1.66E-05                               | 7.79E-07                       | NA                       | NA           | 1.11E-08                                | 3.10E+00                                     | 3.45E-08    |
| Benz(b)fluoranthene    | 4.67E+01                    | 5.11E-02             | 4.53E-05                               | 2.13E-06                       | NA                       | NA           | 3.04E-08                                | 3.10E-01                                     | 9.43E-09    |
| Benz(k)fluoranthene    | 8.05E+00                    | 8.80E-03             | 7.81E-06                               | 3.67E-07                       | NA                       | NA           | 5.24E-09                                | 3.10E-02                                     | 1.62E-10    |
| Chrysene               | 5.02E+01                    | 5.49E-02             | 4.87E-05                               | 2.29E-06                       | NA                       | NA           | 3.27E-08                                | 3.10E-03                                     | 1.01E-10    |
| Dibenz(a,h)anthracene  | 1.04E+00                    | 1.14E-03             | 1.01E-06                               | 4.74E-08                       | NA                       | NA           | 6.77E-10                                | 3.10E+00                                     | 2.10E-09    |
| Indeno(1,2,3-cd)pyrene | 7.55E+00                    | 8.25E-03             | 7.33E-06                               | 3.44E-07                       | NA                       | NA           | 4.92E-09                                | 3.10E-01                                     | 1.52E-09    |

NA - Not Available

Total Cancer Risk: 5.47E-08

Ca = Concentration in Air (mg/m<sup>3</sup>) =  $E_i / (Hb * W * V)$   
 Ei - Emission Rate of Component (mg/sec) = see below  
 Hb - Downwind Ht (m) = 4.81  
 W - Width (m) = 50  
 V - Wind speed (m/sec) = 4.69  
 Length (downwind distance) (m) = 50  
 r - Roughness Ht. (m) = 0.20  
 z - downwind distance (m) = 50  

$$x = 6.25[Hbr * Ln(Hbr/r + 1.58)]$$

$E_i = \text{Emission Rate (mg/sec)} = C_s * (\text{PERv} + \text{PERc})$   
 Cs - Concentration in soil = mg/kg see below

**Table 59****Dermal Exposure to EU6 Sediment by a Child Resident (Aged 1 to 6 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  |                                       | <u>Cs*SA*AH*ABS*EF*ED*CF</u><br>BW*AT |                                |                 |   |   |                |
|---|---------------------------------------|---------------------------------------|--------------------------------|-----------------|---|---|----------------|
| Cs - Concentration in sediment =                            | mg/kg                                 | chem. spec.                           |                                |                 |   |   |                |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                  | 3620                                  | calculated                     |                 |   |   |                |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                       | 20000                                 | USEPA 1997, EFH                |                 |   |   |                |
| Fs - Fraction of skin surface area available for exposure = |                                       | 18.1%                                 | USEPA 1997, EFH                |                 |   |   |                |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                    | 0.044                                 | USEPA 1997, EFH                |                 |   |   |                |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                                       | 0.03                                  | USEPA 1995, Region III         |                 |   |   |                |
| EF - Exposure frequency =                                   | days/year                             | 40                                    | reasonable assumption          |                 |   |   |                |
| ED - Exposure duration =                                    | years                                 | 6                                     | USEPA 1995, Region IV          |                 |   |   |                |
| CF - Conversion factor =                                    | kg/mg                                 | 1.00E-06                              |                                |                 |   |   |                |
| BW - Body weight =  | kg                                    | 70                                    | USEPA 1995, Region IV          |                 |   |   |                |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                                  | 2190                                  | USEPA 1991, HHEM               |                 |   |   |                |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                                  | 25550                                 | USEPA 1991, HHEM               |                 |   |   |                |
| Dermal  |                                       |                                       |                                |                 |   |   |                |
| Constituent   | Concentration<br>in Sediment<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day  | Subchronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average Lifetime<br>Daily Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer<br>Risk |
| <b>Semivolatiles</b>  |                                       |                                       |                                |                 |   |   |                |
| 2-Nitroaniline  | 4.00E-01                              | 2.99E-09                              | NA                             | NA              | 2.56E-10                                      | NA                                      | NA             |
| 2-Nitrophenol   | 8.00E-01                              | 5.98E-09                              | NA                             | NA              | 5.13E-10                                      | NA                                      | NA             |
| 3-Nitroaniline  | 8.00E-01                              | 5.98E-09                              | NA                             | NA              | 5.13E-10                                      | NA                                      | NA             |
| 4-Bromophenylphenoxyether                                   | 8.00E-01                              | 5.98E-09                              | NA                             | NA              | 5.13E-10                                      | NA                                      | NA             |
| 4-Chloro-3-methylphenol                                     | 8.00E-01                              | 5.98E-09                              | NA                             | NA              | 5.13E-10                                      | NA                                      | NA             |
| 4-Chlorophenylphenoxyether                                  | 4.00E-01                              | 2.99E-09                              | NA                             | NA              | 2.56E-10                                      | NA                                      | NA             |
| 4-Nitroaniline  | 8.00E-01                              | 5.98E-09                              | NA                             | NA              | 5.13E-10                                      | NA                                      | NA             |
| Benzo(a)anthracene  | 1.00E+02                              | 7.48E-07                              | NA                             | NA              | 6.41E-08                                      | 7.30E-01                                | 4.68E-08       |
| Benzo(a)pyrene  | 4.90E+01                              | 3.67E-07                              | NA                             | NA              | 3.14E-08                                      | 7.30E+00                                | 2.29E-07       |
| Benzo(b)fluoranthene  | 7.80E+01                              | 5.84E-07                              | NA                             | NA              | 5.00E-08                                      | 7.30E-01                                | 3.65E-08       |
| Benzo(k)fluoranthene  | 2.30E+01                              | 1.72E-07                              | NA                             | NA              | 1.47E-08                                      | 7.30E-02                                | 1.08E-09       |
| Bis(2-chloroethoxy)methane                                  | 8.00E-01                              | 5.98E-09                              | NA                             | NA              | 5.13E-10                                      | NA                                      | NA             |
| Bis(2-chloroethyl)ether                                     | 4.00E-01                              | 2.99E-09                              | NA                             | NA              | 2.56E-10                                      | 1.10E+00                                | 2.82E-10       |
| Carbazole   | 1.00E+02                              | 7.48E-07                              | NA                             | NA              | 6.41E-08                                      | 2.00E-02                                | 1.28E-09       |
| Chrysene  | 7.60E+01                              | 5.69E-07                              | NA                             | NA              | 4.87E-08                                      | 7.30E-03                                |                |
| Dibenz(a,h)anthracene                                       | 9.60E+00                              | 7.18E-08                              | NA                             | NA              | 6.16E-09                                      | 7.30E+00                                | 4.49E-08       |
| Hexachlorobenzene   | 4.00E-01                              | 2.99E-09                              | NA                             | NA              | 2.56E-10                                      | 1.60E+00                                | 4.10E-10       |
| Hexachlorocyclopentadiene                                   | 2.00E+00                              | 1.50E-08                              | NA                             | NA              | 1.28E-09                                      | NA                                      | NA             |
| Indeno(1,2,3-cd)pyrene                                      | 3.90E+01                              | 2.92E-07                              | NA                             | NA              | 2.50E-08                                      | 7.30E-01                                |                |
| N-nitrosodi-n-propylamine                                   | 4.00E-01                              | 2.99E-09                              | NA                             | NA              | 2.56E-10                                      | 7.00E+00                                | 1.80E-09       |

NA - Not Available

Total Cancer Risk = 3.62E-07



**Table 60****Oral Exposure to EU6 Sediment by a Child Resident (Aged 1 to 6 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =       |                                 | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                            |              |   |  |             |
|----------------------------|---------------------------------|---|----------------------------|--------------|---|--|-------------|
| Constituent                | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day  | Oral Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Oral Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles              |                                 |   |                            |              |   |  |             |
| 2-Nitroaniline             | 4.00E-01                        | 1.82E-08  | NA                         | NA           | 1.56E-09                                | NA                                     | NA          |
| 2-Nitrophenol              | 8.00E-01                        | 3.63E-08  | NA                         | NA           | 3.11E-09                                | NA                                     | NA          |
| 3-Nitroaniline             | 8.00E-01                        | 3.63E-08  | NA                         | NA           | 3.11E-09                                | NA                                     | NA          |
| 4-Bromophenylphenylether   | 8.00E-01                        | 3.63E-08  | NA                         | NA           | 3.11E-09                                | NA                                     | NA          |
| 4-Chloro-3-methylphenol    | 8.00E-01                        | 3.63E-08  | NA                         | NA           | 3.11E-09                                | NA                                     | NA          |
| 4-Chlorophenylphenylether  | 4.00E-01                        | 1.82E-08  | NA                         | NA           | 1.56E-09                                | NA                                     | NA          |
| 4-Nitroaniline             | 8.00E-01                        | 3.63E-08  | NA                         | NA           | 3.11E-09                                | NA                                     | NA          |
| Bis(2-chloroethoxy)methane | 8.00E-01                        | 3.63E-08  | NA                         | NA           | 3.11E-09                                | NA                                     | NA          |
| Bis(2-chloroethyl)ether    | 4.00E-01                        | 1.82E-08  | NA                         | NA           | 1.56E-09                                | 1.10E+00                               | 1.71E-09    |
| Carbazole                  | 1.00E+02                        | 4.54E-06  | NA                         | NA           | 3.89E-07                                | 2.00E-02                               | NA          |
| Hexachlorobenzene          | 4.00E-01                        | 1.82E-08  | 8.00E-04                   | 2.27E-05     | 1.56E-09                                | 1.60E+00                               | 2.49E-09    |
| Hexachlorocyclopentadiene  | 2.00E+00                        | 9.08E-08  | 7.00E-03                   | 1.30E-05     | 7.78E-09                                | NA                                     | NA          |
| N-nitrosodi-n-propylamine  | 4.00E-01                        | 1.82E-08  | NA                         | NA           | 1.56E-09                                | 7.00E+00                               | 1.09E-08    |

NA - Not Applicable

Total Hazard Index = 3.57E-05

Total Cancer Risk = 1.51E-08



**Table 61**

**Dermal Exposure to EU6 Sediment by an Adult Resident (Aged 7 to 30 years)**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) = $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |   |                                |                          |                        |          |          |   |           |                                   |             |
|--|---|--------------------------------|--------------------------|------------------------|----------|----------|---|-----------|-----------------------------------|-------------|
| Constituent  | Cs - Concentration in sediment =                            | mg/kg                          | chem. spec.              |                        |          |          |   |           |                                   |             |
|  | SA - Surface area available for exposure =                  | cm <sup>2</sup> /day           | 3620                     | calculated             |          |          |   |           |                                   |             |
|  | SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                | 20000                    | USEPA 1997, EFH        |          |          |   |           |                                   |             |
|  | Fs - Fraction of skin surface area available for exposure = |                                | 18.1%                    | USEPA 1997, EFH        |          |          |   |           |                                   |             |
|  | AH - Adherence factor =                                     | mg/cm <sup>2</sup>             | 0.044                    | USEPA 1997, EFH        |          |          |   |           |                                   |             |
|  | ABS <sub>bap</sub> - Absorption - B(a)P =                   |                                | 0.03                     | USEPA 1995, Region III |          |          |   |           |                                   |             |
|  | EF - Exposure frequency =                                   | days/year                      | 40                       | reasonable assumption  |          |          |   |           |                                   |             |
|  | ED - Exposure duration =                                    | years                          | 24                       | USEPA 1995, Region IV  |          |          |   |           |                                   |             |
|  | CF - Conversion factor =                                    | kg/mg                          | 1.00E-06                 |                        |          |          |   |           |                                   |             |
|  | BW - Body weight =  | kg                             | 70                       | USEPA 1995, Region IV  |          |          |   |           |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                     | days  | 8760                           |                          | USEPA 1991, HHEM       |          |          |   |           |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =                        | days  | 25550                          |                          | USEPA 1991, HHEM       |          |          |   |           |                                   |             |
|  |   |                                |                          |                        |          |          |   |           |                                   |             |
| Constituent  | Concentration in Sediment mg/kg                             | Average Daily Intake mg/kg-day | Subchronic RfD mg/kg-day | Hazard Index           | Dermal   |          | Average Lifetime Daily Intake mg/kg-day |           | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|  |   |                                |                          |                        | Lifetime | Daily    | Intake                                  | mg/kg-day |                                   |             |
| <b>Semivolatiles</b>   |   |                                |                          |                        |          |          |   |           |                                   |             |
| 2-Nitroaniline   | 4.00E-01  | 2.99E-09                       | NA                       | NA                     | 1.03E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| 2-Nitrophenol  | 8.00E-01  | 5.98E-09                       | NA                       | NA                     | 2.05E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| 3-Nitroaniline   | 8.00E-01  | 5.98E-09                       | NA                       | NA                     | 2.05E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| 4-Bromophenylphenylether   | 8.00E-01  | 5.98E-09                       | NA                       | NA                     | 2.05E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| 4-Chloro-3-methylphenol  | 8.00E-01  | 5.98E-09                       | NA                       | NA                     | 2.05E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| 4-Chlorophenylphenylether  | 4.00E-01  | 2.99E-09                       | NA                       | NA                     | 1.03E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| 4-Nitroaniline   | 8.00E-01  | 5.98E-09                       | NA                       | NA                     | 2.05E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| Benzo(a)anthracene   | 1.00E+02  | 7.48E-07                       | NA                       | NA                     | 2.56E-07 | 7.30E-01 | 1.87E-07                                | NA        | NA                                | NA          |
| Benzo(a)pyrene   | 4.90E+01  | 3.67E-07                       | NA                       | NA                     | 1.26E-07 | 7.30E+00 | 9.17E-07                                | NA        | NA                                | NA          |
| Benzo(b)fluoranthene   | 7.80E+01  | 5.84E-07                       | NA                       | NA                     | 2.00E-07 | 7.30E-01 | 1.46E-07                                | NA        | NA                                | NA          |
| Benzo(k)fluoranthene   | 2.30E+01  | 1.72E-07                       | NA                       | NA                     | 5.90E-08 | 7.30E-02 | 4.31E-09                                | NA        | NA                                | NA          |
| Bis(2-chloroethoxy)methane   | 8.00E-01  | 5.98E-09                       | NA                       | NA                     | 2.05E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| Bis(2-chloroethyl)ether  | 4.00E-01  | 2.99E-09                       | NA                       | NA                     | 1.03E-09 | 1.10E+00 | 1.13E-09                                | NA        | NA                                | NA          |
| Carbazole  | 1.00E+02  | 7.48E-07                       | NA                       | NA                     | 2.56E-07 | 2.00E-02 | 5.13E-09                                | NA        | NA                                | NA          |
| Chrysene   | 7.60E+01  | 5.69E-07                       | NA                       | NA                     | 1.95E-07 | 7.30E-03 | 1.42E-09                                | NA        | NA                                | NA          |
| Dibenz(a,h)anthracene  | 9.60E+00  | 7.18E-08                       | NA                       | NA                     | 2.46E-08 | 7.30E+00 | 1.80E-07                                | NA        | NA                                | NA          |
| Hexachlorobenzene  | 4.00E-01  | 2.99E-09                       | NA                       | NA                     | 1.03E-09 | 1.60E+00 | 1.64E-09                                | NA        | NA                                | NA          |
| Hexachlorocyclopentadiene  | 2.00E+00  | 1.50E-08                       | NA                       | NA                     | 5.13E-09 | NA       | NA                                      | NA        | NA                                | NA          |
| Indeno(1,2,3-cd)pyrene   | 3.90E+01  | 2.92E-07                       | NA                       | NA                     | 1.00E-07 | 7.30E-01 | 7.30E-08                                | NA        | NA                                | NA          |
| N-nitrosodi-n-propylamine  | 4.00E-01  | 2.99E-09                       | NA                       | NA                     | 1.03E-09 | 7.00E+00 | 7.18E-09                                | NA        | NA                                | NA          |

NA - Not Available

Total Cancer Risk = 1.52E-06



**Table 62****Oral Exposure to EU6 Sediment by an Adult Resident (Aged 7 to 30 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =                                 |           | <u><math>Cd * IngR * EF * ED * CF * ME</math></u> |                       |  |  |
|--|-----------|---|-----------------------|--|--|
|  |           | <u><math>BW * AT</math></u>                       |                       |  |  |
| Cd - Concentration in sediment =                     | mg/kg     | see below   |                       |  |  |
| IngR - Ingestion rate for sediment =                 | mg/day    | 100   | USEPA 1997, EFH       |  |  |
| EF - Exposure frequency =                            | days/year | 40  | reasonable assumption |  |  |
| ED - Exposure duration =                             | years     | 24  | USEPA 1995, Region IV |  |  |
| CF - Conversion factor =                             | kg/mg     | 1.00E-06  |                       |  |  |
| ME <sub>s</sub> - Matrix effect - PAHs =             |           | 0.29  | Magee, et al., 1996   |  |  |
| BW - Body weight =                                   | kg        | 70  | USEPA 1995, Region IV |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days      | 8760  | USEPA 1991, HHEM      |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days      | 25550   | USEPA 1991, HHEM      |  |  |

| Constituent                | Concentration<br>in Sediment<br>mg/kg | Average<br>Daily Intake<br>mg/kg-day | Oral Chronic<br>RfD<br>mg/kg-day | Hazard Index | Average                               |  |             |
|----------------------------|---------------------------------------|--------------------------------------|----------------------------------|--------------|---------------------------------------|--|-------------|
|                            |                                       |                                      |                                  |              | Lifetime<br>Daily Intake<br>mg/kg-day | Oral Cancer<br>Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>       |                                       |                                      |                                  |              |                                       |  |             |
| 2-Nitroaniline             | 4.00E-01                              | 1.82E-08                             | NA                               | NA           | 6.23E-09                              | NA   | NA          |
| 2-Nitrophenol              | 8.00E-01                              | 3.63E-08                             | NA                               | NA           | 1.25E-08                              | NA   | NA          |
| 3-Nitroaniline             | 8.00E-01                              | 3.63E-08                             | NA                               | NA           | 1.25E-08                              | NA   | NA          |
| 4-Bromophenylphenylether   | 8.00E-01                              | 3.63E-08                             | NA                               | NA           | 1.25E-08                              | NA   | NA          |
| 4-Chloro-3-methylphenol    | 8.00E-01                              | 3.63E-08                             | NA                               | NA           | 1.25E-08                              | NA   | NA          |
| 4-Chlorophenylphenylether  | 4.00E-01                              | 1.82E-08                             | NA                               | NA           | 6.23E-09                              | NA   | NA          |
| 4-Nitroaniline             | 8.00E-01                              | 3.63E-08                             | NA                               | NA           | 1.25E-08                              | NA   | NA          |
| Bis(2-chloroethoxy)methane | 8.00E-01                              | 3.63E-08                             | NA                               | NA           | 1.25E-08                              | NA   | NA          |
| Bis(2-chloroethyl)ether    | 4.00E-01                              | 1.82E-08                             | NA                               | NA           | 6.23E-09                              | 1.10E+00                                     | 6.85E-09    |
| Carbazole                  | 1.00E+02                              | 4.54E-06                             | NA                               | NA           | 1.56E-06                              | 2.00E-02                                     | NA          |
| Hexachlorobenzene          | 4.00E-01                              | 1.82E-08                             | 8.00E-04                         | 2.27E-05     | 6.23E-09                              | 1.60E+00                                     | 9.96E-09    |
| Hexachlorocyclopentadiene  | 2.00E+00                              | 9.08E-08                             | 7.00E-03                         | 1.30E-05     | 3.11E-08                              | NA   | NA          |
| N-nitrosodi-n-propylamine  | 4.00E-01                              | 1.82E-08                             | NA                               | NA           | 6.23E-09                              | 7.00E+00                                     | 4.36E-08    |

NA - Not Applicable

Total Hazard Index = 3.57E-05

Total Cancer Risk = 6.04E-08



**Table 63****Dermal Exposure to EU6 Surface Water by a Child Resident (Aged 1 to 6 years)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u>C<sub>w</sub>*S<sub>A</sub>*K<sub>p</sub>*ABS*ET*EF*ED*CF</u> |
|---|--|
|   | <u>BW*AT</u>   |
| C <sub>w</sub> - Concentration in surface water =                       | mg/L   |
| S <sub>A</sub> - Surface area available for exposure =                  | cm <sup>2</sup>  |
| S <sub>A<sub>t</sub></sub> - Total skin surface area =                  | cm <sup>2</sup>  |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = | 18.1%  |
| K <sub>p</sub> - Dermal permeability constant =                         | cm/hr  |
| A <sub>B<sub>bap</sub></sub> - Absorption - B(a)P =                     | 0.03   |
| A <sub>B<sub>pah</sub></sub> - Absorption - PAHs =                      | 0.1  |
| E <sub>T</sub> - Exposure time =  | hrs/day  |
| E <sub>F</sub> - Exposure frequency =                                   | days/year  |
| E <sub>D</sub> - Exposure duration =                                    | years  |
| C <sub>F</sub> - Conversion factor =                                    | L/cm <sup>3</sup>  |
| B <sub>W</sub> - Body weight =  | kg   |
| A <sub>T<sub>n</sub></sub> - Averaging time - noncarcinogenic =         | days   |
| A <sub>T<sub>c</sub></sub> - Averaging time - carcinogenic =            | days   |
| see below   |  |
| calculated  |  |
| USEPA 1997, EFH   |  |
| USEPA 1997, EFH   |  |
| USEPA 1995, Region III  |  |
| USEPA 1995, Region III  |  |
| USEPA 1992, Dermal Exposure Assessment reasonable assumption            |  |
| USEPA 1995, Region IV   |  |
| USEPA 1995, Region IV   |  |
| USEPA 1991, HHEM  |  |
| USEPA 1991, HHEM  |  |

| Constituent            | Concentration in Surface Water |          | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |                                   |          | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|------------------------|--------------------------------|----------|--------------------------------|------------------------------|--------------|---|-----------------------------------|----------|-----------------------------------|-------------|
|                        | mg/L                           | Kp cm/hr |                                |                              |              | Lifetime Daily Intake mg/kg-day         | Cancer Slope Factor 1/(mg/kg-day) |          |                                   |             |
| <b>Semivolatiles</b>   |                                |          |                                |                              |              |   |                                   |          |                                   |             |
| Benzo(a)anthracene     | 5.00E-04                       | 8.10E-01 | 3.21E-07                       | NA                           | NA           | 2.75E-08                                | 7.30E-01                          | 2.01E-08 |                                   |             |
| Benzo(a)pyrene         | 5.00E-04                       | 1.20E+00 | 4.76E-07                       | NA                           | NA           | 4.08E-08                                | 7.30E+00                          | 2.98E-07 |                                   |             |
| Benzo(b)fluoranthene   | 9.00E-03                       | 1.20E+00 | 8.57E-06                       | NA                           | NA           | 7.34E-07                                | 7.30E-01                          | 5.36E-07 |                                   |             |
| Benzo(k)fluoranthene   | 5.00E-04                       | 4.48E+01 | 1.78E-05                       | NA                           | NA           | 1.52E-06                                | 7.30E-02                          | 1.11E-07 |                                   |             |
| Chrysene               | 5.00E-04                       | 8.10E-01 | 3.21E-07                       | NA                           | NA           | 2.75E-08                                | 7.30E-03                          | 2.01E-10 |                                   |             |
| Dibenz(a,h)anthracene  | 5.00E-04                       | 2.70E+00 | 1.07E-06                       | NA                           | NA           | 9.18E-08                                | 7.30E+00                          | 6.70E-07 |                                   |             |
| Indeno(1,2,3-ed)pyrene | 5.00E-04                       | 1.90E+00 | 7.54E-07                       | NA                           | NA           | 6.46E-08                                | 7.30E-01                          | 4.72E-08 |                                   |             |

NA - Not Available

Total Cancer Risk = 1.68E-06



**Table 64**

**Dermal Exposure to EU6 Surface Water by an Adult Resident (Aged 7 to 30 years)**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u>C<sub>w</sub>*S<sub>A</sub>*K<sub>p</sub>*ABS*ET*EF*ED*CF</u><br>BW*AT |                         |  |                                 |              |  |                                      |             |
|---|---|-------------------------|--|---------------------------------|--------------|--|--------------------------------------|-------------|
| C <sub>w</sub> - Concentration in surface water =                       | mg/L  | see below               |  |                                 |              |  |                                      |             |
| S <sub>A</sub> - Surface area available for exposure =                  | cm <sup>2</sup>   | 3620                    | calculated                             |                                 |              |  |                                      |             |
| S <sub>A<sub>t</sub></sub> - Total skin surface area =                  | cm <sup>2</sup>   | 20000                   | USEPA 1997, EFH                        |                                 |              |  |                                      |             |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = |   | 18.1%                   | USEPA 1997, EFH                        |                                 |              |  |                                      |             |
| K <sub>p</sub> - Dermal permeability constant =                         | cm/hr   | see below               |  |                                 |              |  |                                      |             |
| A <sub>B<sub>bap</sub></sub> - Absorption - B(a)P =                     |   | 0.03                    | USEPA 1995, Region III                 |                                 |              |  |                                      |             |
| A <sub>B<sub>pah</sub></sub> - Absorption - PAHs =                      |   | 0.1                     | USEPA 1995, Region III                 |                                 |              |  |                                      |             |
| E <sub>T</sub> - Exposure time =  | hrs/day   | 1                       | USEPA 1992, Dermal Exposure Assessment |                                 |              |  |                                      |             |
| E <sub>F</sub> - Exposure frequency =                                   | days/year   | 40                      | reasonable assumption                  |                                 |              |  |                                      |             |
| E <sub>D</sub> - Exposure duration =                                    | years   | 24                      | USEPA 1995, Region IV                  |                                 |              |  |                                      |             |
| C <sub>F</sub> - Conversion factor =                                    | L/cm <sup>3</sup>   | 1.00E-03                |  |                                 |              |  |                                      |             |
| B <sub>W</sub> - Body weight =  | kg  | 70                      | USEPA 1995, Region IV                  |                                 |              |  |                                      |             |
| A <sub>T<sub>n</sub></sub> - Averaging time - noncarcinogenic =         | days  | 8760                    | USEPA 1991, HHEM                       |                                 |              |  |                                      |             |
| A <sub>T<sub>c</sub></sub> - Averaging time - carcinogenic =            | days  | 25550                   | USEPA 1991, HHEM                       |                                 |              |  |                                      |             |
| Concentration in Surface Water  |   |                         |  |                                 |              |  |                                      |             |
| Constituent   | mg/L  | K <sub>p</sub><br>cm/hr | Average Daily Intake<br>mg/kg-day      | Dermal Chronic RfD<br>mg/kg-day | Hazard Index | Average Lifetime Daily Intake<br>mg/kg-day | Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>  |   |                         |  |                                 |              |  |                                      |             |
| Benzo(a)anthracene  | 5.00E-04  | 8.10E-01                | 6.89E-08                               | NA                              | NA           | 2.36E-08                                   | 7.30E-01                             | 1.72E-08    |
| Benzo(a)pyrene  | 5.00E-04  | 1.20E+00                | 1.02E-07                               | NA                              | NA           | 3.50E-08                                   | 7.30E+00                             | 2.55E-07    |
| Benzo(b)fluoranthene  | 9.00E-03  | 1.20E+00                | 1.84E-06                               | NA                              | NA           | 6.30E-07                                   | 7.30E-01                             | 4.60E-07    |
| Benzo(k)fluoranthene  | 5.00E-04  | 4.48E+01                | 3.81E-06                               | NA                              | NA           | 1.31E-06                                   | 7.30E-02                             | 9.53E-08    |
| Chrysene  | 5.00E-04  | 8.10E-01                | 6.89E-08                               | NA                              | NA           | 2.36E-08                                   | 7.30E-03                             | 1.72E-10    |
| Dibenz(a,h)anthracene   | 5.00E-04  | 2.70E+00                | 2.30E-07                               | NA                              | NA           | 7.87E-08                                   | 7.30E+00                             | 5.74E-07    |
| Indeno(1,2,3-cd)pyrene  | 5.00E-04  | 1.90E+00                | 1.62E-07                               | NA                              | NA           | 5.54E-08                                   | 7.30E-01                             | 4.04E-08    |

NA - Not Available

Total Cancer Risk = 1.44E-06



**Table 65**

**Dermal Exposure to EU4 Sediment by a Maintenance Worker**

**Preliminary Remediation Goal Calculation**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                |                              |              |   |                                   |             |
|---|---|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| Cs - Concentration in sediment =                            | mg/kg   | chem. spec.                    |                              |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                                | 3620                           | calculated                   |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                                     | 20000                          | USEPA 1997, EFH              |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |   | 18.1%                          | USEPA 1997, EFH              |              |   |                                   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                                  | 0.034                          | USEPA 1997, EFH              |              |   |                                   |             |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |   | 0.03                           | USEPA 1995, Region III       |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year   | 30                             | reasonable assumption        |              |   |                                   |             |
| ED - Exposure duration =                                    | years   | 25                             | USEPA 1995, Region IV        |              |   |                                   |             |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                       |                              |              |   |                                   |             |
| BW - Body weight =  | kg  | 70                             | USEPA 1995, Region IV        |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 9125                           | USEPA 1991, HHEM             |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                          | USEPA 1991, HHEM             |              |   |                                   |             |
|   |   |                                |                              |              |   |                                   |             |
| Constituent   | Sediment mg/kg                                      | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>  |   |                                |                              |              |   |                                   |             |
| Benzo(a)pyrene  | 7.10E+01  | 3.08E-07                       | NA                           | NA           | 1.10E-07                                | 7.30E+00                          | 8.03E-07    |

NA - Not Available

Total Hazard Index =

Total Cancer Risk = 8.03E-07



**Table 66****Dermal Exposure to EU4 Surface Soil (0-6') by a Maintenance Worker****Preliminary Remediation Goal Calculation (3)****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                |                              |              |   |                                   |             |
|---|---|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
| Cs - Concentration in soil =                                | mg/kg   | chem. spec.                    |                              |              |   |                                   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day                                | 3000                           | calculated                   |              |   |                                   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                                     | 20000                          | USEPA 1997, EFH              |              |   |                                   |             |
| Fs - Fraction of skin surface area available for exposure = |   | 15%                            | USEPA 1997, EFH              |              |   |                                   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                                  | 0.038                          | USEPA 1997, EFH              |              |   |                                   |             |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |   | 0.03                           | USEPA 1995, Region III       |              |   |                                   |             |
| EF - Exposure frequency =                                   | days/year   | 30                             | reasonable assumption        |              |   |                                   |             |
| ED - Exposure duration =                                    | years   | 25                             | USEPA 1995, Region IV        |              |   |                                   |             |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                       |                              |              |   |                                   |             |
| BW - Body weight =  | kg  | 70                             | USEPA 1995, Region IV        |              |   |                                   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 9125                           | USEPA 1991, HHEM             |              |   |                                   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                          | USEPA 1991, HHEM             |              |   |                                   |             |
|   |   |                                |                              |              |   |                                   |             |
| Constituent   | Concentration in Soil mg/kg                         | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |   |                                |                              |              |   |                                   |             |
| Benzo(a)pyrene  | 5.60E+01  | 2.25E-07                       | NA                           | NA           | 8.03E-08                                | 7.30E+00                          | 5.86E-07    |

NA - Not Available

Total Hazard Index =

Total Cancer Risk = 5.86E-07



**Table 67****Dermal Exposure to EUS Surface Soil (0-6') by a Maintenance Worker****Preliminary Remediation Goal Calculation****Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) = $\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}$ |                                |                                   |                                 |              |                                    |  |             |
|--|--------------------------------|-----------------------------------|---------------------------------|--------------|------------------------------------|--|-------------|
| Constituent  | Concentration in Soil<br>mg/kg | Average Daily Intake<br>mg/kg-day | Dermal Chronic RfD<br>mg/kg-day | Hazard Index | Lifetime Daily Intake<br>mg/kg-day | Average Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
| Semivolatiles  |                                |                                   |                                 |              |                                    |  |             |
| Benzo(a)pyrene   | 8.00E+00                       | 1.61E-07                          | NA                              | NA           | 5.74E-08                           | 7.30E+00                                     | 4.19E-07    |

NA - Not Available

Total Cancer Risk = 4.19E-07





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**HUMAN HEALTH RISK ASSESSMENT  
FOR THE FORMER GULF STATES CREOSOTING FACILITY,  
HATTIESBURG, MISSISSIPPI**

November 12, 1999

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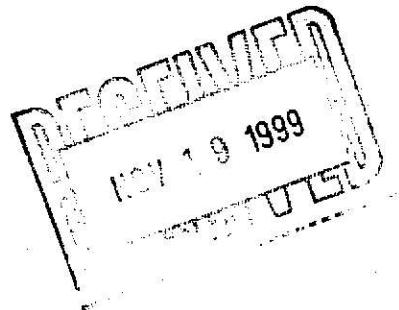
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## Executive Summary

A baseline human health risk assessment (HHRA) was conducted for the Former Gulf States Creosoting facility in Hattiesburg, Mississippi. The HHRA was performed in accordance with: Mississippi Commission on Environmental Quality's (MCEQ's) *Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi* (1999); US EPA's *Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A)* (1989); US EPA Region 4 guidance entitled *Technical Services Supplemental Guidance to RAGS, Region 4 Bulletins* (1995); and other relevant US EPA guidance documents.

Creosoting constituents of potential health concern include polycyclic aromatic hydrocarbons, of which benzo(a)pyrene is the predominant contributor to potential risks. Much of the former creosoting process area is currently covered with asphalt or large building structures. Potential future exposure scenarios included a construction worker, a maintenance worker, and a Site visitor, and assumed bare soils available for direct contact.

Hazards posed by chemical constituents in soil for health effects other than an increased risk of cancer were well below a threshold of possible concern. Cancer risks, due largely to benzo(a)pyrene, were within the US EPA's acceptable target risk range of  $1 \times 10^{-6}$  to  $1 \times 10^{-4}$  (*i.e.*, one in one million to one in ten thousand). The added lifetime cancer risk conservatively estimated for a maintenance worker was  $2 \times 10^{-5}$  (*i.e.*, an upper bound risk of two incidences of cancer out of a population of 100,000 persons so exposed). The potential risk for a construction worker was estimated to be  $3 \times 10^{-6}$ , and the estimated potential risk for a Site visitor was  $7 \times 10^{-7}$ , which is below the *de minimis* acceptable target risk range.

Implementation of a remedy to preclude direct contact with surface soils contaminated with the three highest concentrations of benzo(a)pyrene and other potentially carcinogenic polycyclic aromatic hydrocarbons would reduce estimated cancer risks to below the *de minimis* risk of  $1 \times 10^{-6}$  for all potential exposure scenarios applicable to this site.

## **1.0 Introduction**

Environmental Standards, Inc. (Environmental Standards) was retained by Kerr-McGee Chemical Corporation LLC (Kerr-McGee) to perform a human health risk assessment (HHRA) to evaluate hazards and risks potentially posed by residual levels of chemicals present at the Former Gulf States Creosoting facility (Site). The Site, located near the intersection of US Highways 49 and 11 in Hattiesburg, Mississippi, was formerly a wood treating facility that operated between the early 1900s and 1960. In the early 1960s, the Site was redeveloped for commercial and light industrial uses (Michael Pisani & Assoc., 1997). The land on which the site is located is a portion of the Sixteenth Section land owned by the Hattiesburg Public School District and leased to the current tenants under a 99-year lease, granted on July 7, 1947. At the time of this report, the Site, with the exception of the grassy and wooded areas in the south and southwest, respectively, was primarily used for automobile dealerships. There are no residential or institutional (schools) uses of the Site (Michael Pisani & Assoc., 1997).

Operations at the Site consisted of a small-scale wood preserving process using creosote. The creosoting process was primarily confined to a 2.5-acre area in the northeast corner of the Site; this is known as the Process Area and is currently occupied by Courtesy Ford. During the redevelopment of the Site in the early 1960s, construction debris (e.g., broken concrete, asphalt, etc.) appears to have been relocated to the southwestern corner of the Site along Gordon's Creek. This area is known as the Fill Area and currently remains undeveloped.

This assessment has been conducted as a result of an agreement between Kerr-McGee, the Mississippi Department of Environmental Quality (MDEQ), and the Mississippi Commission on Environmental Quality (MCEQ) pursuant to the Uncontrolled Site Voluntary Evaluation Program. The MDEQ Office of Pollution Control, Uncontrolled Sites Section has been providing oversight and review of investigations and reports relating to the former Gulf States Creosoting facility.

This report will address the potential for on-Site exposures to human receptors. Off-Site areas were considered a separate operable unit and related exposures will be addressed at a later time.

The primary guidance used to develop this risk assessment was the MCEQ *Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi* (1999). US EPA Region 4's *Technical Services Supplemental Guidance to RAGS: Region 4 Bulletins* (1995) were also referred to for guidance. Additional US EPA guidance documents cited herein include:

- *Guidance for Remediation of Uncontrolled Hazardous Substance Sites in Mississippi (MDEQ, 1990);*
- *Risk Assessment Guidance for Superfund, Volume 1, Human Health Evaluation Manual/Part A (RAGS/Part A)* (US EPA, 1989);
- *Human Health Evaluation Manual, Supplemental Guidance: "Standard Default Exposure Factors"* (US EPA, 1991);
- *Exposure Factors Handbook* (US EPA, 1997);
- *Guidelines for Exposure Assessment* (US EPA 1992);
- *Dermal Exposure Assessment: Principles and Applications* (US EPA, 1992);

These documents are not listed in a hierarchical manner; other US EPA guidance documents and peer-reviewed technical papers may have also been referenced in this risk assessment report.

## **2.0 Hazard Identification and Conceptual Site Model**

As a result of the historical wood preservation process, residual levels of creosote-related chemicals are present in soils in the Process Area. Sediment and surface water in a drainage ditch along the southeast border of the Process Area also contain chemical residuals. These site-related chemicals, mostly polycyclic aromatic hydrocarbons (PAHs) are also present in the Fill Area. Residual levels of PAHs have been found in soil in the Fill Area and in Gordon's Creek surface water and sediment.

The PAH residuals have also been detected in shallow groundwater underlying the Site. Currently, there are no private water wells located on-Site that access this shallow groundwater for potable purposes; therefore, groundwater exposures will not be evaluated in this report. Off-Site uses of private wells for potable purposes will be addressed at a later time.

A conceptual site model (CSM) was developed for the Site to aid in determining the potential receptors and exposure units to be evaluated under current and future potential land use (Figure 1). These receptors were identified as infrequent Site visitors, maintenance workers, and construction workers.

Under current land use assumptions, Site visitors may potentially contact residual chemicals in Gordon's Creek surface water and sediment, and/or surface soils in the Fill Area and surrounding woods, the grassy field southeast of the Fill Area, and/or the drainage ditch along side of the Process Area. Visitors may also potentially contact surface soil, surface water, and sediment in the Process Area drainage ditch. The remaining affected areas of the Site are covered with either buildings or pavement precluding casual direct contact with surface soils.

Under both current and future land use assumptions, a maintenance worker may contact surface soils in the Fill Area and surrounding woods, the grassy field southeast of the Fill Area, and/or the Process Area and surrounding affected areas. Although most of the Process Area and vicinity are paved, maintenance activities may involve some shallow digging; therefore, direct

contact with shallow soils in this area was assessed. Maintenance activities are not expected to occur in Gordon's Creek; therefore, maintenance worker exposure to surface water and sediment in Gordon's Creek was not assessed. The remainder of the Site was relatively unaffected by historical creosoting activities.

Although there are currently no major construction activities at the Site, these types of activities may occur at some time in the future. As with the maintenance worker scenario, construction activities could potentially occur in the Fill Area and vicinity, the grassy field southeast of the Fill Area, and the Process Area and vicinity. Construction workers may be exposed to both surface and subsurface soils. Construction activities are not anticipated to occur in Gordon's Creek; therefore, construction worker exposure to surface water and sediment in Gordon's Creek was not assessed. The remainder of the Site was relatively unaffected by historical creosoting activities.

### **3.0 Data Evaluation**

To characterize potential exposures to Site-related chemicals, the former Gulf States Creosoting facility was divided into five exposure units (EUs). Each exposure unit outlines potentially affected areas of the Site and adjacent on-Site locales that may be frequented by individuals accessing the Site for recreational or occupational purposes. The use of EUs is encouraged by the US EPA Region 4 (1995), which defines an EU as "an areal extent of a receptor's movements during a single day..." Each of these exposure units is depicted on Figure 2 and is discussed below.

#### **3.1 Exposure Unit Delineation**

The following EUs were delineated based upon the presence of residual chemicals and the potential for receptors to contact those chemicals. Areas of the Site most affected were included in at least one of the five EUs while areas with relatively low or non-detectable concentrations of residuals were not included in an EU. By limiting Site-wide exposures to the EUs most affected by historical activities at the Site, worst-case scenarios were created.

##### **3.1.1 Exposure Unit 1**

EU1 outlines the on-Site areas in, adjacent to, and downstream of the Fill Area along Gordon's Creek (Figure 2). EU1 addresses exposures to surface water and sediment by an infrequent Site visitor.

##### **3.1.2 Exposure Unit 2**

EU2 delineates the upland areas of the Fill Area and adjacent woody and grassy areas (Figure 2). Surface soils (zero to one foot below ground surface [bgs]) in this area were evaluated for potential visitor and future hypothetical maintenance worker scenarios. Subsurface soils were also evaluated for a hypothetical future construction worker scenario. Available data for subsurface soils for a construction scenario were evaluated to a depth of 16 feet bgs.

### 3.1.3 Exposure Unit 3

In the southwest corner of the Site there exists a grassy field east of West Pine Street between Henson Auto Sales and Eagan Cars and Trucks. This grassy area has been defined as EU3 for purposes of this risk assessment (Figure 2). Similar to EU2, surface soil (zero to one foot bgs) will be evaluated in EU2 for visitor and hypothetical future maintenance worker scenarios. Subsurface soils will also be evaluated for a hypothetical future construction worker scenario. Available data for subsurface soils for a construction scenario were evaluated to a depth of 16 feet bgs.

### 3.1.4 Exposure Unit 4

EU4 encompasses the grassy, drainage ditch area along the fenceline behind Courtesy Ford in the northeast corner of the Site (Figure 2). This area, along the southeast side of the Process Area, has been defined to evaluate Site visitor exposures via casual contact with surface soil, sediment, and surface water in this area. Surface soils in EU4 were defined as zero to one foot bgs.

### 3.1.5 Exposure Unit 5

EU5 outlines the Process Area and the historical drip track areas of the former Gulf States Creosoting facility (Figure 2). This EU encompasses the area delineated in EU4. Surface soils from zero to one foot bgs were evaluated in EU5 for a hypothetical maintenance worker scenario. Available data for subsurface soils down to 16 feet bgs were evaluated in EU5 for a hypothetical future construction worker scenario.

## 3.2 Statistical Evaluation

Environmental samples undergo laboratory analyses that are designed to quantitate the concentrations of constituents in the various environmental media. As a result of the analytical procedures, a constituent may be detected and its concentration measured, detected but not able to be quantitated, or not detected at all in a sample. The data set for the Site contains a number of nondetections for some chemicals of potential concern (COPCs) in various samples. Assuming that the COPC is present in these samples at the achieved detection limit is biased

because the chemical may be absent altogether. Assuming a concentration of zero is also flawed because the chemical could be present at a level below laboratory capabilities to detect and quantify the concentration. Consequently, in the event that an analyte identified at least once in a given medium was not detected in a given sample, it was conservatively assumed for the risk assessment purposes to be present at a concentration equivalent to one-half of the sample quantitation limit (SQL). In addition, samples labeled with an "R" (rejected) qualifier were not included in the data analysis because those data were deemed unreliable and, therefore, unusable. Constituents that were not detected in any sample from a particular medium were eliminated from further consideration in accordance with US EPA guidelines (1989).

Site analytical data used in this assessment were collected during the Phase I (1997) and Phase II (1998) remedial investigations. These data were fully validated by qualified technical professionals using standard data validation protocols, as required by the MCEQ (1999).

In accordance with MCEQ guidance (1999), relative potency factors (RPFs) were applied to determine Site-related concentrations of carcinogenic PAHs. The detected concentration of the carcinogenic PAH (or one-half the detection limit for non-detected samples) was multiplied by an RPF to estimate that PAH's concentration relative to benzo(a)pyrene's (B(a)P) potency. The resulting concentration is a B(a)P potency equivalent. The B(a)P equivalent concentrations for each carcinogenic PAH for each sample were then summed to provide a single B(a)P equivalent concentration per sample. The B(a)P equivalents were then statistically analyzed as discussed below. The RPFs used for this process are as follows (US EPA, 1993):

| <u>Compound</u>      | <u>RPF</u> |
|----------------------|------------|
| Benzo(a)pyrene       | 1.0        |
| Benzo(a)anthracene   | 0.1        |
| Benzo(b)fluoranthene | 0.1        |
| Benzo(k)fluoranthene | 0.01       |
| Chrysene             | 0.001      |

| <u>Compound</u>        | <u>RPF</u> |
|------------------------|------------|
| Dibenz(a,h)anthracene  | 1.0        |
| Indeno(1,2,3-cd)pyrene | 0.1        |

The validated laboratory data were compiled into data sets representing areas of potential exposure (EUs) for each potential receptor. Each data set was analyzed statistically using SiteStat®, a commercially available software package, to calculate the minimum, maximum, arithmetic mean, logarithmic mean, standard error of the mean, and the 95% upper confidence limit of the mean concentration (95% UCL) for each constituent based on distributional analysis of the data (*i.e.*, utilizing goodness-of-fit statistical tests to determine whether the data are distributed normally or lognormally). The data qualifier associated with the *minimum* and *maximum* detected concentrations as well as the location of the maximum detected concentration for each EU were also determined. Results of the quantitative and statistical analyses for each of the EUs discussed above are presented in Tables 1 through 11.

Soils one foot deep were assumed to be representative of surface soils at the Site. This assumption is in accordance with Region 4 guidance (1995). A depth of 16 feet was used to define the extent of subsurface soils at the Site. This depth represents a logical break in the depth intervals from which samples were collected at the Site, and it is the approximate depth that can be reached by construction equipment during the hypothetical excavation of a foundation.

### 3.3 Determination of Exposure-Point Concentrations

Exposure-point concentrations were determined to be the 95% UCL or the maximum concentration of a COPC in an EU, whichever is lower. This methodology is in accordance with US EPA guidance (1989). If the distribution of the concentration data was determined to be lognormal, then the lognormal 95% UCL was compared to the maximum concentration to determine the exposure-point concentration. In the event that the distribution of a chemical in any given medium could not be confidently labeled as normal or lognormal, it is termed either "unknown" or "normal/lognormal." In these cases, the lognormal 95% UCL was compared to

maximum concentration when determining the exposure-point concentration. It should be noted, however, that in cases where the distribution is “unknown,” the normal and lognormal 95% UCLs could not be reliably predicted. Assuming a lognormal distribution of the data increases the uncertainty associated with this step of the risk assessment process; however, hazard and risk estimates are likely to be less uncertain than if the maximum concentrations were used.

Exposure-point concentrations are provided on the statistical summary tables, Tables 1 through 11.

#### 3.4 COPC Selection

Soils (both surface and subsurface) were screened according to MCEQ (1999) guidance. The first tier of the screening process compared maximum concentrations of a constituent in an EU with the Restricted Tier 1 target remediation goal (TRG). Restricted TRGs were used because the site is not currently used for residential purposes and the current commercial/industrial land-use is anticipated to remain into the future. If a maximum concentration of a constituent was less than the Restricted Tier 1 TRG, then that constituent was eliminated from further quantitative assessment.

If the maximum concentration of a constituent in an EU exceeded the Restricted Tier 1 TRG, then the 95% UCL of the constituent was compared to the Restricted Tier 1 TRG. In the event that the concentrations of a chemical were distributed lognormally, the lognormal 95% UCL of that constituent was compared to the Restricted Tier 1 TRG. If the distribution of data of a chemical could not be positively identified as either normal or lognormal, the lognormal 95% UCL was used in the screening process. In these cases, either the maximum concentration or the lognormal 95% UCL can be conservatively used. The US EPA, however, justifies the use of an average concentration as the exposure-point concentration by explaining that toxicity criteria for both carcinogenic and non-carcinogenic effects are based on lifetime average exposures and that the “average concentration is most representative of the concentration that would be contacted at a site over time” (*Supplemental Guidance to RAGS: Calculating the Concentration Term*, 1992). Other US EPA guidance states that “...in most situations, assuming long-term contact with the

maximum concentration is not reasonable" (*Risk Assessment Guidance for Superfund, Part A*, 1989). US EPA Region 4 also states that, generally, it is reasonable to assume that soil data are distributed lognormally (1995). In keeping with these guidances, the lognormal 95% UCL was considered in the screening process where the data distribution for a compound could not be defined as specifically normal or lognormal.

If the 95% UCL (or lognormal 95% UCL where appropriate) of a constituent was less than the Restricted Tier 1 TRG, then that constituent was eliminated from further quantitative analysis. If the 95% UCL (or lognormal 95% UCL where appropriate) of a constituent in soil exceeded the Restricted Tier 1 TRG, then that constituent was retained for quantitative analysis.

MCEQ guidance (1999) does not specify screening levels for constituents in surface water or sediment; therefore, Region 4 was referred to for guidance (1995). In accordance with US EPA Region 4 (1995), maximum constituent concentrations in sediment in an EU were compared to US EPA Region 3 Residential Risk-Based Soil Concentrations (RBCs, 1999). If an RBC was calculated based on non-carcinogenic effects, it was multiplied by 0.1 before comparison to the EU maximum concentration, in accordance with US EPA Region 4 guidance (1995). If a maximum concentration of a constituent in an EU was less than the US EPA Region 3 RBC, then that constituent was eliminated from further quantitative analysis. If the maximum concentration of a constituent in an EU exceeded the US EPA Region 3 RBC, then that constituent was retained for the quantitative risk assessment.

For surface water, the maximum detected concentration of a constituent in an EU was compared to the US EPA Human Health Water Quality Standard (WQS) for consumption of water and organisms in accordance with US EPA Region 4 guidance (1995). If the maximum concentration of a constituent in surface water was less than the WQS, then that constituent was eliminated from quantitative analysis. If the maximum concentration of a constituent in surface water exceeded the WQS, then that constituent was retained for quantitative analysis.

The results of the screening process are presented on the statistical summary tables, Tables 1 through 11. The screening process eliminated all detected COPCs from subsurface soils down to 16 feet bgs in EU3 and surface soils in EU4. For this reason, construction worker exposures to soils in EU3 and maintenance worker and visitor exposures to surface soils in EU4 were not evaluated quantitatively in this assessment.

## **4.0 Exposure Assessment**

Currently, a majority of the Site is used for commercial and light industrial purposes and is paved for roads and parking lots. Unpaved areas are limited to the wooded portion in and around the Fill Area (EU2) and the grassy field outlined by EU3. Since the developed and undeveloped areas of the Site vary considerably with respect to both residual chemical concentrations and land use, the Site was divided into five EUs for the exposure assessment. Chemical data from each EU were combined with EU-specific exposure parameter values and receptor scenarios to determine the chemical intake for each receptor potentially accessing an EU for occupational or recreational purposes.

### **4.1 Receptor Identification**

The following exposures pathways have been selected for this risk assessment as reasonable and realistic scenarios under current and future land-use assumptions:

- Visitor dermal contact with sediment in EU1;
- Visitor ingestion of sediment in EU1;
- Visitor dermal contact with surface water in EU1;
- Visitor dermal contact with surface soil in EU2;
- Visitor ingestion of surface soil in EU2;
- Visitor dermal contact with sediment in EU4;
- Visitor ingestion of sediment in EU4;
- Visitor dermal contact with surface water in EU4;
- Visitor dermal contact with surface soil in EU4;
- Visitor ingestion of surface soil in EU4
- Hypothetical future maintenance worker dermal contact with surface soil in EU2;
- Hypothetical future maintenance worker ingestion of surface soil in EU2;
- Hypothetical future maintenance worker dermal contact with surface soil in EU3;
- Hypothetical future maintenance worker ingestion of surface soil in EU3;

- Hypothetical future maintenance worker dermal contact with surface soil in EU5;
- Hypothetical future maintenance worker ingestion of surface soil in EU5;
- Hypothetical future construction worker dermal exposure to soils in EU2;
- Hypothetical future construction worker ingestion of soils in EU2;
- Hypothetical future construction worker inhalation of fugitive dust in EU2;
- Hypothetical future construction worker ingestion of fugitive dust in EU2;
- Hypothetical future construction worker dermal exposure to soils in EU3;
- Hypothetical future construction worker ingestion of soils in EU3;
- Hypothetical future construction worker inhalation of fugitive dust in EU3;
- Hypothetical future construction worker ingestion of fugitive dust in EU3;
- Hypothetical future construction worker dermal exposure to soils in EU5;
- Hypothetical future construction worker ingestion of soils in EU5;
- Hypothetical future construction worker inhalation of fugitive dust in EU5; and
- Hypothetical future construction worker ingestion of fugitive dust in EU5.

In accordance with MCEQ guidance (1999), intake of carcinogenic PAH compounds via the ingestion route were evaluated qualitatively (see Section 6.1) because the published cancer slope factor for benzo(a)pyrene (IRIS, 1999) cannot be used to quantify carcinogenic risks from ingestion. As a result of the screening process presented in the Data Evaluation section, B(a)P equivalents were the only COPCs selected for the visitor, maintenance worker, and construction worker soil exposures in EU2, the visitor and maintenance worker soil exposures in EU3, and the construction worker and maintenance worker soil exposures in EU5. These soil ingestion scenarios, therefore, were not evaluated quantitatively because of the lack of appropriate published toxicity values for B(a)P. In addition, published oral reference doses were not available at the time of this report to quantify the non-carcinogenic effects of ingestion of PAH compounds in EU1 sediment for the visitor scenario.

Surface water present on-site is either ephemeral or very shallow and is conducive only to wading-type activities. Ingestion of site surface water was considered an insignificant exposure

pathway because the shallow on-site depths preclude complete submersion and subsequent incidental swallowing of the surface water.

In summary, the only ingestion scenario evaluated quantitatively was that of the adolescent visitor in EU4. Appropriate published oral toxicity values for the remainder of the soil exposure scenarios were not available. Surface water ingestion scenarios were considered insignificant in relation to dermal surface water exposures and were not evaluated quantitatively herein.

Each of the potential receptors is discussed below.

#### 4.1.1 Infrequent Site Visitor

Since the Site is not currently fenced or guarded, the general public has access to most areas of the Site at any given time. It is possible that an individual may use some areas of the Site, such as EU1, EU2, or EU3, for recreational purposes. For this reason, sediment and surface water exposures to visitors in EU1, and surface soil exposures in EU2 and EU3 were assessed for the visitor scenario. A strong majority of the remainder of the Site is covered with either buildings or pavement, precluding direct contact with surface soils; however, a small exposed area encompassing a drainage ditch exists along side of the Process Area (EU4). Although this area is not attractive for recreational purposes, it is possible that an individual traversing the Site may contact surface soils, sediment, or surface water in this EU; therefore, these potential exposures were assessed. Sediment exposures in EU1 and EU4 were addressed in accordance with US EPA Region 4 guidance that recommends evaluating sediment exposures in intermittent streams.

#### 4.1.2 Maintenance Worker

Currently, maintenance activities are most likely limited to the developed portions of the Site. Of these, the Process Area and adjacent former drip track area (EU5) were most affected by historical wood preserving processes. Although these areas are mostly paved or built upon, it is possible that maintenance activities may require some shallow digging in unpaved areas; therefore, exposures to surface soils in EU5 were assessed. As a conservative measure, surface soil data from sample locations located in paved areas were evaluated in conjunction with

surface soil data from exposed areas in EU5. If the currently undeveloped portions of the Site (EU2 and EU3) become developed in the future, similar maintenance activities may be required and, therefore, exposures to surface soils in EU2 and EU3 were also assessed.

#### 4.1.3 Construction Worker

Although there are currently no major construction activities at the Site, such activities may hypothetically occur in the future. The three most affected areas of the Site, EU2, EU3, and EU5, were selected to address hypothetical future construction worker scenarios. Construction workers may be exposed to both surface and subsurface soils during activities such as excavating. Subsurface soils, for purposes of this assessment, were defined as those soils 16 feet bgs and shallower. Subsurface soil samples at the Site were collected at various depth intervals, one of which was 14 to 16 feet. Excavating equipment (*e.g.*, backhoe) can generally reach to about 15 feet bgs; consequently, the 14 to 16 foot depth interval presented a logical break in the Site data and was used to define the extent of subsurface soils. Available subsurface soil data were utilized for this exposure scenario.

#### 4.1.4 Future On-Site Residents

The Site is currently zoned for industrial or light-commercial use, and, at the time of this report, there were no plans to develop the Site for residential housing. It is reasonable and realistic to assume that the Site will remain commercial/industrial in the future for the reasons presented below.

The land on which the Site is located is a portion of the Sixteenth Section lands owned by the Hattiesburg Public School District and leased to the current tenants under a 99-year lease, granted on July 7, 1947. The Mississippi Code requires the Board of Education, with the assistance of appropriate public agencies, to survey and classify all Sixteenth Section lands (Miss. Code Ann. §29-3-31 [1972]). Such lands may be classified into one of eight categories, including industrial and commercial. The Site has been classified since approximately 1981 by the Hattiesburg Public School District as either commercial or industrial. Established businesses have been located on the Site since the early 1960s, and later constructions continue in the

commercial or industrial category (car dealerships, used car lots, hardware stores, auto supply, and the like). Current tenants occupy under a lease which does not expire until the year 2046; thus, it can be anticipated that future use will continue as commercial and industrial, the uses to which the current tenants have put the property.

Also, the assumption that future land use will be the same as current land use for exposure assessment purposes is consistent with US EPA's OSWER Directive 9355.7-04 "Land Use in the CERCLA Remedy Selection Process" (1995). This memo states that "...for example, future industrial land use is likely to be a reasonable assumption where a site is currently used for industrial purposes, is located in an area where the surroundings are zoned for industrial use, and the comprehensive plan predicts the site will continue to be used for industrial purposes." The OSWER Directive (1995) goes on to state that "The baseline risk assessment generally needs only to consider the reasonably anticipated future land use..." and "In cases where the future land use is relatively certain, the remedial action objective generally should reflect this land use."

It is reasonable to assume that the future land use is "relatively certain" given the historical land use and the requirements set forth in the Mississippi Code as discussed above. For these reasons, on-Site residential exposures will not be evaluated in this assessment.

#### 4.2 General Intake Equation

Chemical exposure/intake is expressed as the amount of the agent at the exchange boundaries of an organism (*i.e.*, skin, lungs, gut) that is available for systemic absorption. An applied dose is defined as the amount of a chemical at the absorption barriers such as skin, lung, digestive tract, available for absorption and is (usually expressed in milligrams, or mg) absorbed per unit of body weight of the receptor (usually expressed in units of kilogram, or kg). Absorbed dose can be defined as the amount of chemical that penetrates the exchange boundaries. If the exposure occurs over time, the total exposure can be divided by the time period of interest to obtain an average exposure rate (*e.g.*, mg/kg-day). The general equation, as defined by US EPA, for estimating a time-weighted average intake is:

$$\text{Intake (mg/kg - day)} = \frac{C \times IR \times EF \times ED}{BW \times AT} \quad [\text{Equation 1}]$$

where:

C = chemical concentration at the exposure point (e.g., mg/m<sup>3</sup> air);  
IR = intake rate (e.g., m<sup>3</sup>/hr);  
EF = exposure frequency (days/year);  
ED = exposure duration (years);  
BW = body weight of exposed individual (kg); and  
AT = averaging time (period over which exposure is averaged, usually measured in days).

Additional parameters (e.g., skin surface area) were incorporated into the above general equation to evaluate the different potential exposure routes (dermal, oral, inhalation).

Table 12 presents the general and pathway-specific exposure parameters utilized for the intake equations in this assessment.

#### 4.2.1 General Exposure Parameters

Although some of the parameters used to calculate potential exposure are pathway- or route-specific, exposure frequency (EF), exposure duration (ED), averaging time (AT; determined separately for carcinogenic and non-carcinogenic exposures), and body weight (BW) are present in each intake model. These general parameters remain consistent throughout the intake calculations for each specific receptor.

##### 4.2.1.1 Exposure Frequency

The exposure frequency (EF) describes the number of times per year an event is likely to occur. It is most often expressed in units of days/year or events/year, depending on the scenario. Variables such as weather, vacations, sick days, and institutional controls often aid in determining reasonable and realistic exposure frequencies.

The EF for an adolescent visitor was extracted from US EPA *Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A) Interim Final* (1989). This EF

value of 12 days/year per EU is a reasonable estimate that assumes an adolescent would most likely be engaged in outdoor activity on the unpaved areas of the Site for one day a week during the three warmest months of the year.

The EF parameter used for construction workers was 80 days. This is a very conservative assumption in that typical construction projects, especially at industrial complexes, generally involve several phases of activity prior to completion. Examples of these activities include foundation excavation, foundation pouring, building framing, plumbing installation, electrical installation, and roofing. Generally, to complete each of these phases, a different team of specialized contractors is employed to perform the tasks for which they are most qualified. As a result, an individual may only remain at the construction site for a few weeks until his/her task has been completed and the next phase has begun. This is especially true for those activities involving direct contact with soil such as excavating and foundation pouring. Individuals performing these tasks are not usually qualified or employed to continue with the actual building processes. An 80-day EF accounts for an individual to be involved in construction activities for two entire months of the year (assuming five-day work weeks). As explained above, this is a very conservative assumption.

The EF value used for the maintenance worker scenario was 150 days/year. This is also a conservative assumption in that the currently developed areas of the site are covered with buildings or pavement. Maintenance activities in these areas would require little contact with the obscured surface soils. The undeveloped areas of the site currently require little or no maintenance as they are only occasionally mowed or allowed to grow naturally. Should these areas become developed, they will most likely take on the appearance of the remainder of the site, including industrial/commercial buildings and paved roads or parking lots. Once again, extensive direct contact with surface soils would be minimal for a maintenance worker.

#### 4.2.1.2      Exposure Duration

The ED parameter represents the number of years during which an event is likely to occur. Factors affecting this parameter include variables such as age of receptor, population mobility,

and occupational mobility. Exposure durations of less than seven years typically correspond to subchronic exposures while those greater than seven years are typically considered chronic exposures (US EPA, 1989). Toxicity indices are selected based on subchronic or chronic exposure durations.

The future construction worker scenario used an ED of one year because it is highly unlikely that a future construction worker would remain on one site for more than a year. Often, two months is considered the maximum amount of time a construction worker may reasonably remain at the same site.

The future maintenance worker ED, on the other hand, is based on occupational mobility studies. The ED of 25 years was obtained from US EPA (1991) which recommends a 95th percentile value of 25 years based on a study by the Bureau of Labor Statistics as of 1987. US EPA Region 4 also recommends a default value of 25 years for worker scenarios (1995).

The adolescent visitor scenario used an ED of 10 years. An adolescent was defined in this assessment as an individual aged seven to 16 years in accordance with US EPA Region 4 (1995); therefore, an exposure duration of 10 years was most appropriate.

#### 4.2.1.3 Averaging Time

The averaging time (AT) parameter is the time period over which exposure is averaged. For human health cancer risk calculations, the AT<sub>c</sub> value prorates a total cumulative dose over a lifetime. As a conservative approach, the AT<sub>c</sub> value for each receptor is the product of a 365-day year and a 70-year life span, equaling 25,550 days.

The AT<sub>n</sub> used for non-carcinogenic effects is the product of a 365-day year and the exposure duration (*i.e.*, AT<sub>n</sub> = 365 days × ED). Because the ED parameter changes for each receptor, the AT<sub>n</sub> changes as well. The AT<sub>n</sub> values used for each receptor are summarized below:

Future Construction Worker - 365 days  
Maintenance Worker - 9125 days  
Adolescent Visitor - 3650 days

#### 4.2.1.4 Body Weight

The body weight used for the adult exposures (future construction worker, maintenance worker) analyzed in this assessment was the current US EPA default value of 70 kg (US EPA, 1989; US EPA Region 4, 1995). The adolescent body weight used for the visitor scenarios was 45 kg. This value was extracted from US EPA Region 4 guidance (1995).

#### 4.2.2 Route-Specific Exposure Parameters

The general intake equation discussed above (Equation 1) was modified by including route-specific exposure parameters in order to calculate route-specific intake values. For dermal exposures, skin surface area, adherence factor, exposure time (surface water exposures only), and absorption factor parameters were included in the intake equation. For ingestion exposures, an ingestion rate and a matrix effect were included in the intake calculation. For inhalation exposures, an inhalation rate and a retention factor for fugitive dusts were included in the intake equation. Also, for inhalation exposures, an additional paradigm was necessary to convert soil concentrations to concentrations in air available for intake.

##### 4.2.2.1 Dermal Exposure Parameters

###### Skin Surface Area

The total skin surface area used for adult receptors in this assessment was 20,000 cm<sup>2</sup>. This is a US EPA default value extracted from the *Exposure Factors Handbook* (1997). For adolescent exposures, a value of 12,768.3 cm<sup>2</sup> was used for total skin surface area. This was a mean value calculated based on the distributions of total skin surface areas for males and females between the ages of 7 and 16 as presented in *Exposure Factors Handbook* (1997).

For purposes of occupational or recreational exposures, it was assumed that only portions of the body would be exposed to the affected media on the site. For the maintenance worker and construction worker scenarios, it was assumed that the hands and face would be exposed to site

soils. These body parts comprise 9.1% of the total skin surface area, or 1820 cm<sup>2</sup>. The trespasser scenario assumed that the hands, face, and lower legs would be exposed for contact with site media. These body parts comprise 25% of the total skin surface area, or 3192 cm<sup>2</sup>.

#### Soil Adherence Factor

Until recently, the US EPA-recommended default for soil adherence on skin ranged from 0.2 to 1.0 mg/cm<sup>2</sup> for the entire exposed surface area, without consideration of the type of activity (US EPA, 1992). However, the data from which that range was derived were primarily the result of indirect measurements, artificial activities, and sampling of hands only. A more recent study has presented the results of direct measurement of soil loading on skin surfaces before and after normal occupational and recreational activities that might result in soil contact (Kissel *et al.*, 1996). A five-order of magnitude range (roughly 10<sup>3</sup> to 10<sup>+2</sup> mg/cm<sup>2</sup>) was reported for observed activity-related hand loadings. That report indicated that hand loadings within the range of 0.2 to 1.0 mg/cm<sup>2</sup> were produced by activities in which there was vigorous soil contact (e.g., rugby, farming); but for activities in which there was less soil contact (e.g., soccer, professional grounds maintenance), loadings substantially less than 0.2 mg/cm<sup>2</sup> were found on hands and other body parts. Kissel *et al.* (1996) concluded that, because non-hand loadings attributable to higher contact activities exceeded hand loadings resulting from lower contact activities, hand data from limited activities cannot be used as a conservative predictor of loadings that might occur on other body surfaces without regard to activity. Furthermore, because exposures are activity-dependent, dermal exposure to soil should be quantified using data describing human behavior (e.g., type of activity, frequency, duration, including interval before bathing, clothing worn, etc.).

The most recent version of the *Exposure Factors Handbook* (1997) states:

In consideration, of these general observations and the recent data from Kissel *et al.* (1996, 1997), this document recommends a new approach for estimating soil adherence to skin. First use Table 6-12 [Summary of Field Studies, Kissel *et al.*, 1996a] to select the activity which best approximates the exposure scenario of concern. Next, use Table 6-13 [Mean Soil Adherence by Activity and Body Region, Kissel *et al.*, 1996a] to select soil loadings on exposed skin surfaces

which correspond to the activity of interest. This table contains soil loading estimates for various body parts. The estimates were derived from soil adherence measurements of body parts of individuals engaged in specific activities described in Table 6-12. These results provide the best estimate of central loadings, but are based on limited data. Therefore, they have a high degree of uncertainty such that considerable judgment must be used when selecting them for an assessment.

In another study that assessed the percentage of skin coverage in several soil contact trials in a greenhouse and an irrigation pipe laying trial, Kissel *et al.* (1996) concluded that adjusted loadings may be two to three orders of magnitude larger than average loadings if average loadings are small.

The activity-specific soil adherence factor for exposures to a maintenance worker was calculated based on data presented by Kissel *et al.* (1996) for grounds keepers, as presented below:

|                    |                         | Soil Adherence Factor by Body Part (mg/cm <sup>2</sup> ) |                |                 |               |
|--------------------|-------------------------|--|----------------|-----------------|---------------|
| Receptor           | Representative Activity | Hands  | Arms           | Lower Legs      | Face          |
| Maintenance Worker | Grounds Keepers         | 0.030 - 0.15   | 0.0021 - 0.023 | 0.0008 - 0.0012 | 0.0021 - 0.01 |

Data for the grounds keepers were used for the maintenance worker estimates because the activities of a grounds keeper best mimic those of a maintenance worker.

Soil adherence factors were calculated by normalizing each body part-specific soil adherence value (using the mid-points of the ranges tabulated above) with regard to the percentage of total body surface area represented by the respective body part (extracted from the US EPA *Dermal Exposure Assessment: Principles and Applications* [US EPA, 1992]). The maintenance worker adherence factor was calculated based upon exposure to the hands and face. Surface area percentages for hands and face are 5.2 and 3.9 percent, respectively (US EPA, 1992). Those body parts comprise 9.1 percent of the total body surface area. The normalized values for all body parts of interest were added, and the sum was divided by the total percentage of body

surface area occupied by the parts. For example, the soil adherence factor for the maintenance worker ( $0.054 \text{ mg/cm}^2$ ) was calculated as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.09 \times 0.052) + (0.006 \times 0.039)}{0.091} = 0.054$$

The construction worker adherence factor was also calculated in this fashion. This exposure scenario, however, assumed that the hands, arms, legs, and face would be exposed to site soils. Soil loadings for the upper torso (chest and back) were not measured by Kissel *et al.* (1996) for construction workers because this body area is generally covered. However, to account for exposure to the upper torso during the very hot months of the year, the total area of the arms, legs, hands, and face were assumed to be completely exposed. The hands, arms, legs, and face comprise 5.2%, 5.9%, 12.8%, and 2.6% of the total skin surface area, respectively (with the face comprising one-third the surface area of the head), for a total of 26.5% exposed surface area. The construction worker soil adherence factor was based on data from Kissel *et al.* (1996) for construction workers as follows:

|                     |                         | Soil Adherence Factor by Body Part ( $\text{mg/cm}^2$ ) |       |            |       |
|---------------------|-------------------------|---|-------|------------|-------|
| Receptor            | Representative Activity | Hands   | Arms  | Lower Legs | Face  |
| Construction Worker | Construction Worker     | 0.24  | 0.098 | 0.066      | 0.029 |

The soil adherence factor for the construction worker scenario was calculated as follows:

$$AF(\text{mg/cm}^2) = \frac{(0.052 \times 0.24) + (0.059 \times 0.098) + (0.128 \times 0.066) + (0.026 \times 0.029)}{0.265} = 0.104$$

The adherence factor for visitor exposures to soil and sediment assumed that the arms, hands, and lower legs would be exposed to soil or sediment. The data used in these calculation were

based on data by Kissel *et al.* (1996) for soccer players (exposed to a playing field of roughly one-half grass and one-half bare earth in a light mist) as presented below:

| Receptor | Representative Activity | Soil Adherence Factor by Body Part (mg/cm <sup>2</sup> ) |              |                |
|----------|-------------------------|--|--------------|----------------|
|          |                         | Arms   | Hands        | Lower Legs     |
| Visitor  | Soccer Players          | 0.0029 – 0.011   | 0.019 – 0.11 | 0.0081 – 0.031 |

The arms, hands, and lower legs comprise 5.9%, 5.2%, and 12.8% of the total skin surface area, respectively, for a total of 23.9% (US EPA *Exposure Factors Handbook*, 1997). The adherence factor was then calculated for visitor dermal exposures to soil and sediment as follows:

$$AF \text{ (mg/cm}^2\text{)} = \frac{(0.00695 \times 0.059) + (0.0645 \times 0.052) + (0.0196 \times 0.128)}{0.239} = 0.026$$

A value of 0.026 mg/cm<sup>2</sup> was used as the soil adherence factor for visitors to the Site.

#### Exposure Time

To estimate intakes as a result of dermal exposure to surface water, an exposure time (ET) parameter was included in the intake formula for Site visitors. The parameter value of 1.0 hour/day was estimated using best professional judgement. This value represents the amount of time a site visitor may spend exposed to surface water in any one EU during the course of his/her visit to the site.

#### Dermal Permeability Constant

The permeability constant, K<sub>p</sub>, accounts for the movement of a constituent dissolved in water through the skin, across the stratum corneum, and into the blood stream. K<sub>p</sub> values for the constituents examined in this assessment for surface water exposures were obtained from US EPA *Dermal Exposure Assessment: Principles and Applications* (1992). For values not available in

US EPA *Dermal Exposure Assessment* (1992), the K<sub>p</sub> value can be calculated using the equations provided by the US EPA in the same document.

#### Dermal Absorption Factor

The final parameter included in the dermal intake paradigm was a dermal absorption factor. In general, the skin provides an effective barrier to environmental toxins. For example, certain hair-coloring formulations which are vigorously rubbed onto the scalp on a daily basis contain lead acetate at concentrations up to 200,000 ppm, yet lead toxicity does not appear to result. Moore *et al.* (1980) determined that the rate of lead absorption from 203<sup>Pb</sup> labeled lead acetate in cosmetic preparations containing six mmol Pb acetate/L in male volunteers over 12 hours was 0.06% during normal use of such preparations. For most inorganic salts, percutaneous (skin) absorption is considered insignificant relative to incidental ingestion (for example, US EPA, 1986). On the other hand, some drugs (*e.g.*, nicotine) are effectively administered and absorbed into the blood stream from dermal "patches."

Most dermal bioavailability data for impacted soil have been obtained in laboratory animals or in vitro test systems. This introduces a significant source of uncertainty for predicting the human response. Safety factors have sometimes been applied to dermal absorption data obtained in animals to conservatively estimate the upper-bound of likely human percutaneous uptake of a certain constituent from skin exposure. This is usually unnecessary because human skin has generally been shown, for a diverse group of constituents, to be about 10-fold less permeable than the skin of typical animal species, such as rabbits and rats (Bartek and LaBudde, 1975; Shu *et al.*, 1988).

US EPA Region III evaluated available data concerning the dermal absorption of specific constituents and classes of constituents and provided several recommendations (US EPA Region 3, 1995). For semivolatile compounds, such as *bis(2-ethylhexyl)phthalate*, the US EPA recommends a range of 1% to 10% (US EPA, 1995). Kao *et al.* (1985) reported 2.7 percent for absorption of topically applied pure benzo(a)pyrene by human skin *in vitro*. The US EPA

Region 3 recommends using 10% as a conservative assumption based on the Ryan *et al.* study (1987). For the purpose of this risk assessment, an ABS of 3% for benzo(a)pyrene and of 10% for other semivolatile organic compounds were conservatively assumed for dermal absorption, in keeping with US EPA Region 3's recommendations.

#### 4.2.2.2 Ingestion Exposure Parameters

##### Ingestion Rate

The ingestion rate used for the adolescent visitor scenario was 100 mg/day. The US EPA *Exposure Factors Handbook* (1997) recommends a value of 100 mg/day as a mean ingestion rate for children under six years of age. This value was conservatively used in this assessment to estimate sediment ingestion exposures for an adolescent visitor aged 7 to 16 years.

##### Gastrointestinal Matrix Effects of Soil

Incidental ingestion incorporates the matrix effect (ME; sometimes called the absorption adjustment factor [AAF]) into the general intake equation. When constituents are administered in solid vehicles such as food and soil, only a fraction of the ingested dose is extracted from the vehicle and subsequently absorbed through the gastrointestinal tract (US EPA *Estimated Exposure to Dioxin-like Compounds*, 1992). Gastrointestinal absorption of constituents sorbed onto such a medium is inhibited by physical-constituent bonding to the matrix (Hawley, 1985). This phenomenon is referred to as the gastrointestinal matrix effect of soil. Several studies referenced in the US EPA's *Estimated Exposure to Dioxin-like Compounds* (1992) have been performed to estimate the oral absorption factors of constituents from soil.

For PAHs, a gastrointestinal matrix effect value of 0.29 was utilized as recommended by Magee *et al.*, 1996. This value is the average of 12 oral soil AAF studies performed by Rozett *et al.*, Weyand *et al.*, and Goonet (Magee, 1996).

#### 4.2.2.3 Inhalation Exposure Parameters and Paradigms

##### Inhalation Rate

The inhalation rate used for the construction worker scenario was 20 m<sup>3</sup>/day. This is a common US EPA default value and was recommended by US EPA Region 4 (1995).

##### Retention Factor

According to the International Commission on Radiological Protection (ICRP), 75 percent of respirable dust particles (PM<sub>10</sub>, or particles less than 10 microns in aerodynamic diameter) are retained when inhaled, the vast majority of which is potentially subsequently swallowed (ICRP, 1968). This 75% was included in the inhalation intake equation as the retention factor parameter (RF). This parameter applies only to non-VOC constituents entrained onto dust particles.

##### Concentration in Air

To estimate airborne dust levels during hypothetical construction activities, an emission rate of suspendible particles of less than 15 microns in aerodynamic diameter (PM<sub>15</sub>) was calculated (grams/second); particles less than 10 microns were considered to be respirable. Considering particles of 15 microns or less in diameter in the emission rate calculation is a conservative assumption, inasmuch as only particles with an aerodynamic diameter of less than five to seven microns are inhaled into the lung.

The two types of construction activities at the Site that have the potential to emit fugitive dusts are vehicular movement over bare (unpaved or unvegetated) surfaces and the excavation of soil. Estimation of fugitive dust emissions caused by each activity were examined separately, as follows, and were derived from existing estimates of general construction exposure. The sum of the emissions from these two activities was multiplied by the concentration of constituent in the soil (Cs) in order to derive the total emission rate (Ei) for non-VOCs as follows:

$$Ei = C_s \times (PERv + PERe) \quad [Equation 2]$$

where:

- E<sub>i</sub> = Emission rate (mg/sec);  
C<sub>s</sub> = Concentration in soil (mg/kg);  
PERv = Particulate emission rate for vehicular movement (lb/vehicle mile);  
and  
PERe = Particulate emission rate for excavation (lb/vehicle mile).

The following empirical expression (US EPA, 1988) was used to estimate the fugitive dust generated by vehicles during construction activities:

$$\text{PERv (lbs/vehicle mile)} = k \times 5.9 \times (s/12)(S/30) \times (\text{mvw}/3)^{0.7} \times (\text{ww}/4)^{0.5} \times ((365 - p)/365)$$

[Equation 3]

where:

- PERv = Vehicle particle emission rate (lb/vehicle mile traveled);  
s = Percent silt content (unitless);  
k = Particle size multiplier (unitless);  
S = Mean vehicle speed (mph);  
mvw = Mean vehicle weight (ton);  
ww = Mean number of wheels per vehicle (unitless); and  
p = Mean number of days with  $\geq 0.01$  inches of precipitation per year (unitless).

It was assumed that the vehicle travels during 40% of the 80-day exposure duration and 0.5 miles per day. The result is a value of 16 miles per construction event. Percent silt content was estimated to have a mean value of 50%, based on geotechnical data provided in the *Remedial Investigation Report* (Pisani & Assoc., 1997). US EPA default values were utilized and referenced for all other parameters. The particle size multiplier was assumed to be 0.50, corresponding to particles less than 15 microns (US EPA, 1996). Vehicle characteristics consist of the following: mean vehicle speed was assumed to be 15 mph, with mean vehicle weight assumed to be approximately 12.5 tons, for 8-wheeled vehicles (US EPA, 1988). The estimated mean number of days with precipitation equal to or greater than 0.01 inches per year is 110 (US EPA, 1988). Total resultant dust emissions for constituents during vehicular movement activities were estimated to be approximately 16.5 lbs/vehicle mile traveled, or 0.0001 kg/sec. Calculations are summarized in Table 13.

Future excavation may be performed by bulldozers, a backhoe, or other heavy construction equipment. The following estimate of particulate emissions, less than 15  $\mu\text{m}$  in diameter resulting from bulldozing activity, was based on the approach described in the US EPA *Compilation of Air Pollution Emission Factors* (1996), as developed from studies of emissions from uncontrolled open dust sources resulting from bulldozing at western surface coal mines.

$$\text{PERe (lb/hour)} = \frac{1.0 \times s^{1.5}}{M^{1.4}} \quad [\text{Equation 4}]$$

where:

PERe = Excavation particle emission rate (lb/hr);  
s = Percent silt content (unitless); and  
M = Soil moisture content (unitless).

Percent soil moisture content was assumed to be 15.1%, an average of site-specific soil moisture data and percent silt content 50%, as described above.

The resultant fugitive dust emission rate during excavation activities was 7.9 lbs/hr or 0.001 kg/sec. Table 13 summarizes these calculations.

Once the emission rate (Ei in Equation 2) was calculated, it was converted to a concentration in ambient air. Gaussian models are conventionally used to determine downwind ambient air concentrations, Ca, from the emission rate, Ei, estimated. However, in this scenario, such models have limited applicability when the receptor(s) is at or very near the source of emission. In this case, a bulldozer operator, for example, is situated directly within the area of ground emissions of vapors and dusts. Average ambient air concentrations in this circumstance are best estimated by use of a near-field box model (US EPA, 1988).

The near-field box model assumes uniform wind speed and uniform mixing throughout the box. The release and mixing of VOCs or respirable dusts in ambient air is estimated as follows:

$$C_a \text{ (mg/m}^3\text{)} = \frac{E_i}{W_b \times H_b \times V} \quad [\text{Equation 5}]$$

where:

$C_a$  = concentration of constituent in ambient air ( $\text{mg/m}^3$ );  
 $E_i$  = emission rate of constituent ( $\text{mg/sec}$ );  
 $W_b$  = width of box in crosswind dimension within the area of residual constituent in soil (m);  
 $H_b$  = downwind height of box (m); and  
 $V$  = average wind speed through the box ( $\text{m/sec}$ ).

The value of  $H_b$  in this calculation is determined by the downwind distance and the atmospheric turbulence at ground level, which determines the trajectory of a release from the upwind edge of the source of vapor or dust emissions. For neutral atmospheric conditions, the height at the downwind boundary ( $H_b$ ) may be expressed by the following function (Pasquill 1975, Horst 1979):

$$z = 6.25 r [H_b/r \times \ln(H_b/r) - 1.58 H_b/r + 1.58] \quad [\text{Equation 6}]$$

where:

$H_b$  = Downwind height of box (m);  
 $z$  = Downwind distance to boundary (m); and  
 $r$  = A terrain-dependent roughness height (m)

$H_b$  (defined in Equation 5) is adjusted until the  $z$  parameter is equal to  $W_b$  (defined in Equation 5). The resulting  $H_b$  value is the height of the box. On any given workday, it is estimated that grading or excavation activities occur over the entire "workable" site area (exposure unit) from which dusts are generated. This area is estimated to be  $2,500 \text{ m}^2$ , with

length of the box estimated to be 50 meters (downwind distance) and the width of the box (W) estimated to be 50 meters. The greater the roughness height, the greater the wind turbulence and constituent dilution (*i.e.*, the height of the box increases). For the purposes of this risk assessment, it is conservatively assumed that the roughness height is 0.20 meters, which corresponds to a terrain with grass, some small bushes, and occasional trees (US EPA *Rapid Assessment of Exposure to Particulate Emission from Surface Contamination Sites*, 1985). This assumption is appropriate for the actual Site conditions. An annual average wind speed (4.69 m/sec) is obtained from the STAR data set, accessed through the Personal Computer Graphical Exposure Modeling System (PCGEMS), for STAR station 03940, Jackson/Thompson, MS for the period 1974-1978 (Table 14).

## **5.0      Toxicity Assessment**

The toxicity assessment involves the evaluation of available toxicity information to be utilized in the risk assessment process. Toxicity values derived from a dose-response relationship can be used to estimate the potential for the occurrence of adverse effects in individuals exposed to various constituent levels.

Exposure to a constituent does not necessarily result in adverse effects. The relationship between dose and response defines the quantitative indices of toxicity required to evaluate the potential health risks associated with a given level of exposure. If the nature of the dose-response relationship is such that no effects can be demonstrated below a certain level of exposure, a threshold can be defined and an acceptable exposure level derived. Humans are routinely exposed to naturally-occurring constituents and man-made constituents through the typical diet, air, and water, with no apparent adverse effects. However, the potential for adverse effects may occur if the exposure level exceeds the threshold in a variably sensitive population. This threshold applies primarily to constituents which produce non-carcinogenic (systemic) effects, although there is a growing body of scientific evidence which suggests that exposure thresholds may exist for certain carcinogenic constituents as well.

Adverse effects can be caused by acute exposure, which is a single or short-term exposure to a toxic substance, or by chronic exposure on a continuous or repeated basis over an extended period of time. "Acceptable" acute or chronic levels of exposure are considered to be without any anticipated adverse effects. Such exposure levels are commonly expressed as reference doses (RfDs), health advisories, etc. An acceptable exposure level is calculated to provide an "adequate margin of safety."

Chronic RfDs, which have been derived by the US EPA for a large number of constituents, were utilized to evaluate exposures lasting seven to 70 years (US EPA, 1989). Activities involving exposures of shorter duration to COPCs at the Site are anticipated to result in hazard and risk estimates that are lower than those associated with the long-term exposures. Identification of

subchronic toxicity values corresponding to shorter-term exposure scenarios (*i.e.*, less than seven years) are included in the risk assessment to ensure that both short-term and long-term risks can be addressed.

Currently, the US EPA has not developed toxicity values to be utilized in dermal exposure scenarios; however, the US EPA does provide the following guidance for dermal exposure:

No RfDs or slope factors are available for the dermal route of exposure. In some cases, however, non-carcinogenic or carcinogenic risks associated with dermal exposure can be evaluated using an oral RfD or oral slope factor, respectively. (US EPA, 1989).

Provisional dermal toxicity values were developed and utilized in the dermal exposure pathways considered in the human health risk assessment to provide a more accurate Site-specific risk assessment. These dermal RfD values were developed by multiplying the published oral RfD for a given constituent by the fraction of that constituent that can be absorbed through the gastrointestinal tract (stomach/intestine lining). The absorption fraction utilized was 50% for semivolatiles as extracted from US EPA Region 4 guidance (1995).

A number of sources of toxicity information exists, and these sources vary with regard to the availability and strength of supporting evidence. The following protocol has been established for the determination of toxicity indices; it defines a hierarchy of sources to be consulted and the methodology for the determination of toxicity values. This protocol has been developed in accordance with current US EPA methodology. Toxicity values for the COPCs at the Site were obtained with reference to the following hierarchy of sources developed in accordance with MCEQ guidance (1999):

- 1) Toxicity values were obtained from the *Integrated Risk Information System* (IRIS, 1999) database. This database contains the RfDs and Cancer Slope Factors (CSFs), which have been verified by the US EPA's RfD and Carcinogen Risk Assessment Verification Endeavor (CRAVE) workgroups, and is, thus, the

agency's preferred source for toxicity values. IRIS supersedes all other information sources.

- 2) For toxicity values which are unavailable on IRIS, the most current source of information is the Health Effects Assessment Summary Tables (HEAST, US EPA, 1997), published by the US EPA. HEAST contains interim, as well as verified RfDs and CSFs. Supporting toxicity information for verified values is provided in an extensive reference section of HEAST.
- 3) In cases where IRIS or HEAST could not provide toxicity values, US EPA Region III's Risk-Based Concentration (RBC) Tables were visited. These tables often provide toxicity values generated by reliable sources other than IRIS or HEAST. For example, in response to specific requests from risk assessors, the US EPA National Center for Environmental Assessment (NCEA) develops provisional RfDs or CSFs for chemicals not listed in IRIS or HEAST. Region III's RBC tables will list such provisional values. Also, RfDs or CSFs that have since been withdrawn from IRIS or HEAST may still be listed on the Region III RBC tables, although they are flagged with a "W." These toxicity values were no longer agreed upon by US EPA scientists; however, the Region III RBC tables continue to publish such values because risk assessors still need to quantify exposures to these chemicals. Lastly, the Region III RBC tables will list toxicity indices found in "other" US EPA documents. These values are flagged with an "O" on the tables.

The US EPA has derived carcinogenic slope factors for both oral and inhalation pathways, and these are utilized to quantitatively estimate risks. In the first step of the US EPA's evaluation, the available data are analyzed to determine the likelihood that the agent is a human carcinogen. The evidence is characterized separately for human studies and animal studies as sufficient, limited, inadequate, no data, or evidence of no effect. The characterizations of these two types of data are combined, and based on the extent to which the agent has been shown to be a carcinogen in experimental animals or humans, or both, the agent is given a provisional weight-of-evidence classification. The US EPA scientists then adjust the provisional classification upward or downward, based on other supporting evidence of carcinogenicity (see Section 7.1.3, US EPA, 1989). For a further description of the role of supporting evidence, see the US EPA guidelines (US EPA, 1986).

The US EPA classification system for weight of evidence is shown in the table below. This system is adapted from the approach taken by the International Agency for Research on Cancer.

| US EPA WEIGHT-OF-EVIDENCE<br>CLASSIFICATION SYSTEM FOR<br>CARCINOGENICITY |  |
|---|--|
| Group   | Description  |
| A   | Human carcinogen   |
| B1 or<br>B2   | Probable human carcinogen<br><br>B1 indicates that limited human data are available<br><br>B2 indicates sufficient evidence in animals and inadequate or no evidence in humans |
| C   | Possible human carcinogen  |
| D   | Not classifiable as to human carcinogenicity   |
| E   | Evidence of non-carcinogenicity for humans   |

(US EPA, 1989)

Table 15 summarizes the available toxicity values for the identified COPCs. COPCs lacking published toxicity values were not able to be quantitatively evaluated in this assessment in accordance with MCEQ guidance (1999). The MCEQ limits the use of toxicity values to those that have been published in IRIS, HEAST, ATSDR toxicity profiles, or other peer-reviewed reference sources or literature approved by the MCEQ (1999).

## **6.0 Risk Characterization**

The objective of the risk characterization is to determine potential risk to receptors by combining the results of the exposure and toxicity assessments. Non-carcinogenic effects and carcinogenic risks are summarized in Table 16. Tables 17 through 30 provide algorithms and parameters for each pathway.

The estimated intakes calculated for each exposure pathway considered and each COPC were compared to RfDs for non-carcinogenic effects. The following formula was used to estimate the potential for non-carcinogenic health effects for each COPC.

$$HQ = ADI/RfD \quad [Equation 6]$$

where:

HQ = Hazard quotient - potential for noncancer health effects (unitless);  
ADI = Average daily intake of COPC (mg/kg-day); and  
RfD = Reference dose (mg/kg-day).

RfDs have been developed by the US EPA for chronic (*e.g.*, lifetime) and/or subchronic exposure to constituents based on the most sensitive non-carcinogenic effects. The chronic RfD for a constituent is an estimate of a lifetime daily exposure level for the human population, including sensitive subpopulations, that is likely to be without an appreciable risk of deleterious effects. The potential for noncancer health effects was evaluated by comparing the Site-specific exposure level with the RfD derived by the US EPA for a similar exposure period. This ratio of exposure to toxicity is called the hazard quotient (HQ). If the Site-specific exposure level exceeds the threshold (*i.e.*, the HQ exceeds a value greater than 1.0), there may be concern for potential noncancer effects.

To assess the overall potential for noncancer effects posed by multiple constituents, a hazard index (HI) is derived by summing the individual HQs. This approach assumes additivity of critical effects of multiple constituents. This is appropriate only for compounds that induce the

same effect by the same mechanism of action. This conservative approach significantly overestimates the actual potential for adverse health impacts.

In cancer risk assessment, the US EPA has required the use of the upper limit which produces an estimate of potential risk that has a 95% probability of exceeding the actual risk, which may, in fact, be zero. The following formula was utilized to estimate the upper bound excess cancer risk for each carcinogen (note that not all COPCs are carcinogens):

$$TR = CLDI \times SF \quad [\text{Equation 7}]$$

where:

- TR = Target risk - excess probability of an individual developing cancer (unitless);  
CLDI = Calculated lifetime average daily intake of carcinogenic COPC (mg/kg-day); and  
SF = Cancer slope factor (mg/kg-day)<sup>-1</sup>.

For exposures to multiple carcinogens, the upper limits of cancer risks are summed to derive a total cancer risk. The US EPA recognizes that it is not technically appropriate to sum upper confidence limits of the risk to produce a realistic total probability, but requires this approach be used.

Carcinogenic risk refers to the probability of developing cancer as a result of exposure to known or suspected carcinogens. The National Contingency Plan (NCP) endorses an acceptable risk range of  $10^{-4}$  to  $10^{-6}$  for exposure to multiple carcinogens. This range represents an incremental increase of 1 in 10,000 to 1 in 1,000,000 in the chance of developing cancer over a lifetime.

Table 16 provides a summary of the non-carcinogenic effects and carcinogenic risks associated with each of the pathways evaluated in this assessment.

The overall hazard index across the assessed pathways and EU's was 0.02 for the Site visitor scenario. This value is below the acceptable benchmark of 1.0. The highest hazard index

associated with the Site visitor scenario was 0.01 corresponding to oral exposure to sediment in EU4. The overall cancer risk for exposures to Site visitors was estimated to be  $7 \times 10^{-7}$ . This value is below the *de minimis* benchmark of  $1 \times 10^{-6}$ .

Hazard indices for the maintenance worker scenario were not quantified because published reference doses were not available for the selected COPCs. The overall cancer risk for the maintenance worker scenario was  $2 \times 10^{-5}$ . This value is within the acceptable range of  $1 \times 10^{-4}$  to  $1 \times 10^{-6}$ . The cancer risk level for the maintenance worker was primarily attributable to dermal exposure to B(a)P equivalents in surface soils in EU5.

Hazard indices for the hypothetical future construction worker were not quantified because published reference doses were not available for the selected COPCs. The overall cancer risk for the hypothetical future construction worker scenario was  $3 \times 10^{-6}$ . The cancer risk level for the construction worker was primarily attributable to dermal exposure to B(a)P equivalents in soils in EU5. The construction worker scenario assumed that a significant skin surface area was available for soil adherence. During the cooler times of the year when more clothing is probably worn, risks posed to construction workers would likely fall below the *de minimis* target risk value of  $1 \times 10^{-6}$ .

The scenarios that drove the cancer risk levels at the Site were the maintenance worker and construction worker dermal exposures to B(a)P equivalents in EU5. These scenarios exceeded a *de minimis* risk level of  $1 \times 10^{-6}$ , but were less than an upper-bound target risk level of  $1 \times 10^{-4}$ , as specified by US EPA.

## 6.1 Assessment of the Ingestion Path of Carcinogenic PAHs

### 6.1.1 Background

Under Subpart II of the Risk Evaluation Procedures, MCEQ states that US EPA's published cancer slope factor utilized to quantitatively estimate risks due to dermal absorption of

carcinogenic PAHs cannot be used to quantify carcinogenic risks from incidental ingestion of PAHs. Rather, the guidance states that such potential ingestion risk shall be described qualitatively. This section addresses the qualitative added risk from this exposure route and also provides useful perspective concerning typical daily exposures to PAH compounds by the general populace.

PAHs are ubiquitous in the environment and occur from both natural sources and human-related activities. Microbial synthesis, higher plant synthesis, volcanic activity, and prairie and forest fires are all major contributors to the natural background levels of PAHs (ATSDR, 1999; Suess, 1976; Andelman *et al.*, 1980; and Youngblood *et al.*, 1975). However, the majority of PAHs in the environment is formed by anthropogenic sources. PAH mixtures are found in tar, soot, petroleum, tobacco smoke, automotive exhaust, engine lubricant wastes, asphalt, fried and broiled foods, smoked meats, etc., and are generally formed by the incomplete combustion of organic material. PAHs are present in ambient air, water, sediments, soils, and foods (ATSDR, 1999).

In the environment, PAHs almost always occur as complex mixtures of many compounds made up of three or more fused benzene rings in linear, angular, or cluster arrangements and contain only carbon and hydrogen (Edwards, 1983). The most studied member of the class is B(a)P. The use of B(a)P as a model PAH is so prevalent that PAHs and B(a)P are often thought of synonymously. B(a)P may be as much as three orders of magnitude more potent than certain other carcinogenic PAHs (Chu and Chen, 1984, as cited in ATSDR, 1999). It should be emphasized that B(a)P constitutes only 1-20% of the total PAHs usually found in a complex environmental mixture (Blumer, 1961). Menzie *et al.* (1992), report that B(a)P represents about 10 to 30% of carcinogenic PAHs in a mixture. Nevertheless, because B(a)P is well-characterized and many PAHs in a mixture may not be identified in typical gas chromatography/mass spectrometry (GC/MS) analysis, some authors still report only B(a)P concentrations as representative of a complex mixture. Typically, total PAHs (*i.e.*, the five to 20 PAHs quantified in published reports) in soils are about 8 to 10 times the value of B(a)P alone (Edwards, 1983).

Typical concentrations of B(a)P in various materials or media are summarized in the following table:

#### **Polycyclic Aromatic Hydrocarbon Content Of Various Materials And Media**

| Material                                | Benzo(a)pyrene ( $\mu\text{g}/\text{kg}$ ) | Reference                     |
|---|--|-------------------------------|
| Coal                                    | 4,000                                      | Coomes – 1981                 |
| Asphalt                                 | 10,000 – 100,000                           | Coomes – 1981                 |
| Raw shale oil                           | 3,200                                      | Coomes – 1981                 |
| Crude oil                               | 400 – 2,800                                | Pancirov & Brown – 1975       |
| Coal tar pitch                          | up to 12,500,000                           | Wallcave – 1971               |
| Eastern US soils<br>(rural)             | 40 – 1,300                                 | Blumer – 1961                 |
| Soil near Swiss<br>highway              | 2,100 – 30,000                             | Blumer <i>et al.</i> – 1977   |
| Fresh waters ( $\mu\text{g}/\text{L}$ ) | 0.001 – 0.100                              | Andelman & Suess – 1980       |
| Urban air (ng/m)                        | 0.03 – 104                                 | IARC – 1973                   |
| Marine sediments                        | nil – 5,000                                | Zobell – 1971                 |
| Margarine                               | 0.2 – 6.8                                  | Swallow – 1976                |
| Sunflower oil                           | 8.0  | Swallow – 1976                |
| Barbecued ribs                          | 10.5                                       | Libinsky & Shubik – 1965      |
| Liquid smoke<br>flavoring               | 9,400,000                                  | Youngblood & Blumer –<br>1975 |
| Barbecued beef                          | 3.5  | Frethein – 1976               |
| Tea leaves                              | 3.9 – 21.3                                 | IARC – 1973                   |
| Lettuce<br>(industrial area)            | 8.6 – 150                                  | Edwards – 1983                |

The main source of contamination of the soil by PAHs is by deposition from air. It has been estimated that global emissions of B(a)P during the time period between 1966 and 1969 were about 4.6 million kg (about 10 million pounds) per year, mostly from fossil fuel burning, with motor vehicles contributing only about 1% of the total (Suess, 1976, as cited in ATSDR, 1988). However, background concentrations of PAHs in surface soils near major highways tend to be

relatively high. Blumer and coworkers (1975) showed that the major source of PAHs found in soils of a Swiss mountain town located in a deep valley with frequent temperature inversions was automobile exhaust (Blumer *et al.*, 1977); concentrations of PAHs in soils close to a major highway were found to be much greater than those in soils of the same locality but nearer to industries.

The PAHs emitted from jet aircraft exhaust provide clear illustration of the influence of emission source locations to background soil levels of PAHs. Shabad reported that a jet airplane engine emits 2000 to 4000 µg B(a)P/min when cruising at high altitude, but emits amazingly up to 40,000,000 µg B(a)P/min during take-off (Shabad, 1980, as cited in ATSDR, 1988). Not surprisingly then, soil, snow, and vegetation near airports were found to contain B(a)P (and PAHs) at levels an order of magnitude or more higher than at a control site (Smirnov, 1970 and Audere *et al.* 1973, as cited in ATSDR, 1988). On a global scale, however, amounts of PAHs emitted from jet engine exhausts are very small compared to emissions from fossil fuel combustion and other sources (Edwards, 1983).

Carcinogenic PAHs are found in all surface soils. Evidence of the global distribution of PAHs was given by Thomas (1986, as cited in ASTDR, 1999) who detected carcinogenic PAHs in arctic soils above 150 µg/kg.

Although background levels of PAHs will vary depending on locality and proximity to emission sources, some generalizations regarding typical levels in soils, sediments, water, air, etc., have been reported. Blumer and coworkers examined surface soils in many areas of the eastern United States and noted that levels between 40 and 1,300 µg/kg in soils from relatively rural areas (Blumer, 1961; 1976). Typical concentrations of B(a)P in soils of the world according to Edwards range from about 100 µg/kg to 1000 µg/kg, with total PAHs about 10 times the value of B(a)P alone (Edwards, 1983).

In source dominated areas soils may contain much higher concentrations. Butler (1984) examined soils in Switzerland and demonstrated that much higher levels of PAHs, up to 300,000 µg/kg, are found near complex road interchanges than in areas more distant (Blumer *et al.*, 1977; Butler *et al.*, 1984). B(a)P concentrations as high as 650,000 µg/kg have been reported in soil samples collected 10 m from a German soot plant and as high as 120,000 µg/kg 500 m from a tar plant (Freitz, 1971).

As anticipated, urban areas have higher soil concentrations than do more remote areas because of the proximity to sources of fossil fuel combustion. The majority of urban soil concentrations of carcinogenic PAH concentrations falls in the 600 to 3,000 µg/kg range (Menzie *et al.*, 1992). The highest ambient concentrations of carcinogenic PAHs in soils have been reported for road dust, which can contain levels of 8,000 to 336,000 µg/kg (Menzie *et al.*, 1992).

Sediments are major sinks for PAHs, primarily because of the low solubility of these compounds and their strong affinity for organic carbon in particulate matter. For example, sediment samples taken from the Detroit River in 1982 contained B(a)P at concentrations ranging from 120 µg/kg to 17,640 µg/kg (Health & Welfare Canada, 1979,).

#### 6.1.2 Qualitative Assessment of Incidental Ingestion of PAHs in Soil

Carcinogenic PAHs are thought to increase the risk of cancer by all routes of exposure. Accordingly, the risks estimated by dermal absorption and inhalation exposure may underestimate the upper bound of cancer risk from additional exposure via incidental ingestion of soil at the Site. However, the risks estimated for dermal absorption of PAHs adsorbed into soils adhering to skin may overestimate risks for a host of reasons.

Early studies conducted by Falk and coworkers indicated that the carcinogenic effect of B(a)P on subcutaneous injection in mice could be markedly inhibited by the simultaneous administration of various non-carcinogenic PAHs (Falk *et al.*, 1964, as cited in ATSDR, 1988. In other subcutaneous injection and skin-painting studies with mice, it was shown that a combination of

several non-carcinogenic PAH compounds, mixed according to the proportion occurring in auto exhaust, did not enhance or inhibit the action of two potent PAH carcinogens, B(a)P and dibenz(a,h)anthracene.(ATSDR, 1988).

The carcinogenic potency of B(a)P and other carcinogenic PAHs is generally determined by injecting solutions under the skin, painting the skin with the carcinogenic PAH dissolved in a solvent, or dissolved in corn oil in feeding studies. This vehicle or matrix affords a high level of bioavailability of the carcinogenic PAH compound. Recently, Krueger *et al.* (1999) conducted *in vitro* percutaneous absorption studies with contaminated soils and organic solvent extracts of contaminated soils collected at former manufactured gas plant (MGP) sites. The MGP tar-contaminated soils contained PAHs at levels ranging from 10 to 2400 mg/kg. The dermal penetration rates of PAH from the MGP tar-contaminated soils and soil solvent extracts were determined experimentally through human skin using tritium-labelled B(a)P as a surrogate. Results showed reductions of two to three orders of magnitude in PAH absorption through human skin from the most contaminated soils in comparison to the soil extracts. Reduction in PAH penetration was attributed to soil matrix properties. That is, PAH compounds adsorbed to organic carbon in a soil matrix are far less bioavailable for dermal flux than PAH compounds dissolved in a solvent. [No correction for such a profound soil matrix effect was applied in quantitatively estimating cancer risks due to dermal absorption of B(a)P and other carcinogenic PAHs in this assessment.]

Because PAHs are ubiquitous, humans are exposed to these chemicals as part of everyday living. PAHs have been detected in many food products, particularly leafy vegetables, especially those grown in urban areas or near roadways. Even smoking "one pack of cigarettes a day," which has been estimated to result in a B(a)P exposure of 0.4 µg/kg (ATSDR, 1988), is less than the typical intake via the diet. Santodonato *et al.* (1981) reviewed the literature on human exposure including dietary doses of PAHs and summarized the findings as follows.

| Source | B(a)P Concentration | Carcinogenic PAH Concentration | Total PAH Concentration |
|--------|---------------------|--------------------------------|-------------------------|
| Air    | 0.0095 – 0.0435 µg  | 0.038 µg                       | 0.207 µg                |
| Water  | 0.0011 µg           | 0.0042 µg                      | 0.0270 µg               |
| Food   | 0.16 – 1.6 µg       | no data                        | 1.6 – 16 µg             |

Statistics available from the United States Department of Agriculture (Pao *et al.*, and United States Department of Agriculture, 1986) were used to determine an average diet for US males. For the average American diet, the intake of carcinogenic PAHs was estimated to be between 1 and 5 µg/day. Menzie *et al.* (1992) reported comparable results for total potential dose of carcinogenic PAHs for nonsmoking adult males, and estimated a median of 3 µg/day and a maximum of 15 µg/day. According to these authors, this would translate to a daily dietary dose of B(a)P of 0.6 to 3.0 µg.

It should be noted that the available monitoring data are very limited with regard to dairy products, such as cheese and milk (Santodonato *et al.*, 1981), which make up approximately one-third of the daily diet of 1600 g. In one study conducted in Canada, B(a)P concentrations of 7.6 to 387 µg/L, with a mean value of 129.5 µg/L were found in samples of milk from ten nursing mothers (HSDB, 1999). Assuming ingestion of one liter per day of mother's milk by a nursing baby, the intake of B(a)P alone would, on average, be about 130 µg.

To provide qualitative perspective with regard to the incidental ingestion of soil at the Hattiesburg Site, if a maintenance worker were to ingest 100 mg of soil per day, 150 days per year, with an average B(a)P concentration of 39.25 mg/kg in the former Process Area, the average daily intake of B(a)P would be about 1.6 µg. This appears to be comparable to or less than the typical dietary intake from food. A charcoal-broiled T-bone steak reportedly contains about 50 µg/kg of B(a)P, as well as various levels of other PAHs (Lijinsky and Ross, 1977, as cited in HSDB, 1999). For a 10-ounce portion, this translates to a dose of 14 µg B(a)P from this

food item alone. It may be noted here also that, with few exceptions, PAHs in soil in the Process Area are covered with asphalt, and therefore, not readily available for direct contact and incidental ingestion.

## **7.0 Uncertainty Analysis**

Risk assessment uses a wide array of information sources and techniques. Even in those rare circumstances where constituent intake for an exposed individual may be measured relatively precisely, assumptions will still be required to evaluate the associated risk. Generally, data are not available for critical aspects of the risk assessment, and the use of professional judgment, inferences based on analogy, the use of default values, model estimation techniques, etc., result in uncertainty of varying degrees.

The expressions of risk in this assessment are not probabilistic; the expressions of risk are conditional, based on the conditions represented by the single-point values selected for the analysis. This section is intended to identify and qualitatively evaluate the more salient Site-specific uncertainties and their potential influence on the credibility of the estimated Site risks.

### **7.1 Uncertainty of Data Evaluation Factors**

Uncertainties in data analysis include analytical error, selection of COPCs, adequacy of sampling design, etc. Generally, there is far less uncertainty in this phase of the risk assessment process than other aspects contribute.

Laboratory analysis is extremely accurate relative to the potential error of "professional judgment" in exposure assessments. The uncertainty of analytical data is likely to be less than 25 percent, most of the time.

The adequacy of the sampling strategies to characterize Site conditions is a potentially large source of uncertainty. Because of the limited availability of resources, sample collection is generally limited. However, sampling (especially in multiple surveys) is not random, but is designed to locate the areas with the highest levels of constituents. Thus, test data are biased toward overestimation of average constituent levels. In addition, in most instances, the upper 95-percent confidence limit of the average concentration is utilized as an exposure-point

concentration in the risk assessment. The use of this value likely will result in an overestimation of risk, as the 95% UCL represents a value that will be greater than the true average 95% of the time.

Oftentimes, only a portion of detected constituents are carried through the risk assessment process because constituents are eliminated through COPC screening procedures (US EPA, 1989). This could result in an underestimation of risk, although the COPC selection process is intended to identify those constituents which account for the vast majority of potential risk. COPCs lacking published RfD values were not quantitatively evaluated and this may result in an underestimation of potential hazards (non-carcinogenic effects).

## 7.2 Uncertainty of Toxicity Values

The US EPA's IRIS states that the uncertainty associated with RfD values for non-carcinogenic endpoints of toxicity "span perhaps an order of magnitude." In fact, the uncertainty of extrapolating dose-response data from animals to humans with the application of multiple safety factors (100 to 10,000 or more) is likely to be several orders of magnitude. Current policies for deriving RfD values will often result in an overestimation of risk.

The uncertainty associated with the estimation of cancer risk contributes, by far, the major source of potential error and uncertainty. It is beyond the scope of this analysis to explore this toxicity assessment factor in any detail. However, a few salient points are noted below.

Some constituents classified as carcinogens have been shown to produce an increased incidence of cancer in mice but not rats, for example. If the mouse is not an adequate model for the rat, it may be wondered how reliable a model it is for human beings. The assumption of linearity and a non-threshold phenomenon in the dose versus risk relationship may not be valid and could result in a very large overestimation of actual cancer risk, if any even exist at low doses in humans.

The US EPA evaluated the uncertainty of cancer risk estimates from exposures to trichloroethene and several other related VOCs in public drinking water supplies (Cothorn *et al.*, 1984). These US EPA scientists concluded the following:

- The largest uncertainty in the calculations is due to the choice of the model [Multistage, Weibull, Logit, Probit, etc.] used in extrapolating risk to low doses in humans, and is 5 to 6 orders of magnitude;
- If a single model were chosen [assumed to be valid], the overall uncertainty in risk estimates would be 2 to 3 orders of magnitude;
- The exposure estimates contribute, at most, an order of magnitude to the uncertainty; and
- It would appear that until a particular compound's mechanisms of cancer are better known, it is likely that the uncertainty in the toxicity will not be improved.

### 7.3 Uncertainties in Assessing Potential Exposure

Ideally, Site-specific exposure values should be used when assessing potential intakes of chemicals at a Site. Oftentimes, however, Site-specific data are not available; therefore, the risk assessor must estimate values that most accurately reflect Site conditions. In doing so, US EPA or other regulatory default values were utilized in place of Site-specific data. These values may over- or under-estimate risks, depending on Site conditions and the percentile range in which the default values fall (*e.g.*, 50<sup>th</sup>, 95<sup>th</sup>).

Although a considerable amount of published data is available on the most common exposure parameters (*e.g.*, body weight, skin surface area), even these data contain uncertainties. Studies conducted by different scientists often provide differing levels of detail, statistics, and accuracy based on sample size, study design, geographic area, etc. Such discrepancies can increase uncertainty when the data are combined to derive a single-point default value. These data may be the best available; however, the reflection of reality may still be imprecise.

Where published exposure parameters were not available, best professional judgment had to be used, thereby increasing uncertainty. The default or estimated exposure parameters used in this assessment likely resulted in a moderate over-estimation of risk.

MCEQ (1999) guidance recommends a qualitative evaluation of carcinogenic risks resulting from ingestion of carcinogenic PAH compounds. By not including carcinogenic PAH ingestion in the quantitative evaluation, potential risks are clearly underestimated. However, by evaluating carcinogenic PAHs as a B(a)P equivalent for dermal and inhalation exposures, the hazard and risk estimates are actually calculated for the seven carcinogenic PAHs. Therefore, hazard and risk levels calculated using B(a)P equivalents actually account for the intakes of seven PAH compounds where acceptable risk levels (*e.g.*, one in one million) are provided for individual compounds (MCEQ, 1999).

## 8.0 Summary of Findings

The results of the baseline human health risk assessment indicate potentially unacceptable risk values for the maintenance worker and construction worker exposed to soils in EU5 as a result of the presence of residual B(a)P equivalents. To determine the extent of remediation necessary to reduce these risks to acceptable levels, soil data in EU5 were closely examined.

In EU5, the surface sample locations contributing most to elevated risk levels for the maintenance worker scenario were GEO-19/0-1', GEO-21/0-1', and GEO-33/0-1' (see Figure 2). B(a)P equivalent concentrations at these locations contributed most to the maintenance worker risk estimates. Sample location GEO-33/0-1' is located within a paved area in a parcel of land southwest of Courtesy Ford (Figure 2). Pavement in this area precludes direct contact with surface soils; therefore, it is not anticipated that current or future maintenance workers will have access to surface soils in or around sample location GEO-33/0-1'. Sample locations GEO-19/0-1' and GEO-21/0-1' are located between the fenced Courtesy Ford property and the railroad tracks to the southeast of Courtesy Ford (Figure 2). These sample locations are adjacent to a drainage ditch that is mostly vegetated and accessible to the general public.

For the maintenance worker scenario, elimination of contact with the three maximum concentrations of B(a)P equivalents would reduce risk levels in EU5 to below the acceptable risk range of  $1 \times 10^{-6}$  to  $1 \times 10^{-4}$ . Existing pavement already precludes direct contact with surface soils at location GEO-33/0-1'. Implementing a remedy to preclude contact with surface soils along the drainage ditch in the vicinity of EU4 would significantly reduce current risk levels. The risk level associated with B(a)P residuals was recalculated excluding the three maximum B(a)P concentration sample locations in EU5. A reevaluation of risks excluding contact with the most affected soils is presented in Tables 31 and 32. Results of this evaluation indicate risk levels below the *de minimis* lower range of the target risk level.

B(a)P equivalents in soil sample locations GEO-21/0-1' and GEO-21/2-3' contributed most to the elevated risk levels for the construction worker scenario in EU5. Implementing a remedy to

preclude contact with soils down to three feet bgs at sample location GEO-21 would significantly reduce current risk levels. The remedy for the construction worker scenario in EU5 can be implemented in conjunction with the remedy for the maintenance worker scenario in EU5 since sampling location GEO-21 is a risk driver for both scenarios. By implementing a remedy that would preclude contact with sample locations GEO-21/0-1' and GEO-21/2-3', cancer risk levels for the construction worker scenario in EU5 would be reduced to below the *de minimis* risk level of  $1\times10^{-6}$  (see Tables 31 and 33). Implementation of a remedy precluding contact with GEO-19/0-1', as a result of the EU5 maintenance worker exposures, would further reduce construction worker risk levels.

## Bibliography

- Andelman, J. B., and M. J. Suess. 1980. Polynuclear aromatic hydrocarbons in the water environment. Bull. WHO 43:479-508.
- ATSDR (Agency for Toxic Substances and Disease Registry). Toxicological Profile for Benzo(a)Pyrene. Oak Ridge National Laboratory. 1988.
- ATSDR (Agency for Toxic Substances and Disease Registry). ATSDR's Toxicological Profiles on CD-ROM. Polycyclic Aromatic Hydrocarbons (PAHs), Update. CRC Press, 1999.
- Audere, A. K., Z. Y. Lindberg, G. A. Smirnov, and L. M. Shabad. 1973. Experiment in studying the influence of an airport located within the limits of a city on the level of environmental pollution by benzo(a)pyrene. Gig. Sanit. 38(9): 90-92.
- Bartek, M.J. and J.A. LaBudde. Percutaneous Absorption *in vitro*, in Animal Models in Dermatology. Ed. H.I. Maibach. New York: Churchill Livingstone, 1975. p. 103.
- Blumer, M. 1961. Benzpyrenes in soil. Science 134, 474-475.
- Blumer, M., W. Blumer, and T. Relch. 1977. Polycyclic aromatic hydrocarbons in soils of a mountain valley; correlation with highway traffic and cancer incidence. Environ. Sci. Technol. 11 (12), 1082-1084.
- Butler, J. D., V. Butterworth, C. Kellow, and H. G. Robinson. 1984. Some observations on the polycyclic aromatic hydrocarbon (PAH) content of surface soils in urban areas. Sci. Total Environ. 38, 75-85.
- Chu, M. M. L. and G. W. Chem. 1984. Evaluation and estimation of potential carcinogenic risks of polynuclear aromatic hydrocarbons. Paper presented at the Pacific Rim Risk Conference.
- Coomes, R. M. 1981. Carcinogenic testing of oil shale materials. Twelfth Oil Shale Symposium Proceedings. Colorado School of Mines Pres.
- Cothorn, C. R., W. Conniglio, W. Marcus. Techniques for the Assessment of the Carcinogenic Risk to the US Population due to Exposure from Selected Volatile Organic Compounds from Drinking Water via the Ingestion, Inhalation and Dermal Routes. NTIS PB84-213941. Office of Drinking Water. Washington DC: Environmental Protection Agency, 1984.
- Edwards, N. T. 1983. Polycyclic aromatic hydrocarbons (PAHs) in the terrestrial environment – a review. J. Environ. Qual. 12 (4), 427-441.

- Falk, H. L., and P. T. S. Kotin. Inhibition of carcinogenesis: The effects of polycyclic hydrocarbons and related compounds. *Arch. Environ. Health* Vol. 9 (1964):169-179.
- Fritz, W. 1971. Extent and sources of contamination of our food with carcinogenic hydrocarbons. *Ernaehrungsforschung* 16(4), 547-557.
- Health & Welfare Canada. Polycyclic Aromatic Hydrocarbons, Report No. 80-EHD-50, (1979) p. 38.
- Horst, T. W. Langrangian Similarity Modeling of Vertical Diffusion for a Ground Level Source. *Int. Applied Met.*, Vol. 18 (1979): 733-740.
- HSDB (Hazardous Substances Data Bank), 1999. National Library of Medicine (NLM) On-Line Toxicological Network (TOXNET). Bethesda, MD.
- ICRP (International Commission on Radiological Protection). Report of Committee IV on Evaluation of Radiation Doses to Body Tissues from Internal Contamination due to Occupational Exposure. ICRP Publication 10. New York: Pergamon Press, 1968.
- Kao, J.K., F.K. Patterson, and J. Hall. Skin Penetration and Metabolism of Topically Applied Chemicals in Six Mammalian Species, Including Man: an in vitro Study with Benzo(a)pyrene and Testcaterone. *Toxicol. Appl. Pharmacol.*, Vol. 81 (1985): 502-516.
- Kissel, J., K. Richter, and R. Fenske. Field Measurements of Dermal Soil Loading Attributable to Various Activities: Implications for Exposure Assessment. *Risk Analysis*, Vol. 16, No. 1 (1996): 115-125.
- Magee, B., P. Anderson, and D. Burmaster. Absorption Adjustment Factor (AAF) Distributions for Polycyclic Aromatic Hydrocarbons (PAHs). *Human and Ecological Assessment: An International Journal*. Vol. 2, No. 4 (December 1996): 841-873.
- Menzi, C.A., B.B. Potocki and J. Santodonato. Exposure to Carcinogenic PAHs in the Environment. *Environ. Sci. Technol.* Vol. 26, No. 7, 1992.
- Michael Pisani & Associates. Remedial Investigation Report, Former Gulf States Creosoting Site, Hattiesburg, Mississippi. New Orleans, Louisiana. 1997.
- Michael Pisani & Associates. Phase II Remedial Investigation Report, Former Gulf States Creosoting Site, Hattiesburg, Mississippi. New Orleans, Louisiana. 1998.
- Mississippi Commission on Environmental Quality (MCEQ). Final Regulations Governing Brownfields Voluntary Cleanup and Redevelopment in Mississippi. 1999.

Mississippi Department of Environmental Quality (MDEQ). Guidance for Remediation of Uncontrolled Hazardous Substance Sites in Mississippi. Office of Pollution Control. 1990.

Moore, M.R., P.A. Meredith, W.S. Watson, D.J. Sumner, M.K. Taylor, and A. Goldberg. "The Percutaneous Absorption of Lead-203 in Humans From Cosmetic Preparations Containing Lead Acetate, as Assessed by Whole-Body Counting and Other Techniques." *Food Cosmet. Toxicol.* 18. (1980): 399.

Pancirov, R. J. and R. A. Brown. 1975. Analytical methods for polynuclear aromatic hydrocarbons in crude oil, heating oils, and marine tissues. In: Conference on prevention and control of oil pollution, San Francisco, CA, March, 1975. American Petroleum Institute, Wash., DC. pp 103-13.

Pao, E. M. *et al.* Home Economics Research Report No. 44. United States Department of Agriculture, Washington, DC. 1982.

Pasquill, I.. The Dispersion of Material in the Atmospheric Boundary Layer - The Basis for Generalization. In: *Lectures on Air Pollution and Environmental Impact Analysis*. Boston, MA: American Meteorological Society, 1975.

Ryan, E.A., E.T. Hawkins *et al.* "Assessing Risk From Dermal Exposure at Hazardous Waste Sites. in Bennett." Ed. G. and J. Bennett. Superfund '87: Proceedings of the Eighth National Conference. Washington, DC, 16-18 November 1987. The Hazardous Material Control Research Institute. p.166-168.

Santodonato, J., P. Howard, and D. Basu. Health and Ecological Assessment of Polynuclear Aromatic Hydrocarbons. Pathotox Publishers, Inc., Park Forest South, IL. 1981.

Shabad, L. M. 1980. Circulation of carcinogenic polycyclic aromatic hydrocarbons in the human environment and cancer prevention. *J. Natl. Cancer Inst.* 64(3): 405-410.

Shu, H.P., P. Teitelbaum, A.S. Webb, L. Marple, B. Brunck, D. Dei Rossi, F.J. Murray, and D.J. Paustenbach. "Bioavailability of Soil Bound TCDD: Dermal Bioavailability in the Rat." *Fundam. Appl. Toxicol.*, Vol. 10 (1988): 648-654.

Smirnov, G. A. 1970. The study of benz(a)pyrene content in soil and vegetation in the airfield region. *Vopr. Onkol.* 16(5): 83-86.

State of Mississippi. Mississippi Code 1972 Annotated. Title 29 Public Lands, Buildings, and Property, Chapter 3, Sixteenth Section and Lieu Lands in General. 1998.

Suess, M. J. 1976. The environmental load and cycle of polycyclic aromatic hydrocarbons. *Sci. Total Environ.* 6:239-250.

Ta, Roy *et al.*, Studies Estimating the Dermal Bioavailability of Polynuclear Aromatic Hydrocarbons from Manufactured Gas Plan Tar-Contaminated Soils. Env. Sci. Tech. 32(20). 1998. 3113-3117.

Thomas, J.F., M. Mukai, and B.D. Teggens. Fate of airbone benzo(a)pyrene. Environ. Sci. Technol. 2:33-39, 1968.

United States Department of Agriculture (USDA). Nationwide Food Consumption Survey: Continuing Survey of Food Intakes by Individuals, Men 19-50 years, 1 Day, 1985; United States Department of Agriculture. Human Nutrition Information Service. Nutrition Monitoring Division; Washington, DC, Report No. 86-1. 1986.

US EPA (United States Environmental Protection Agency). Rapid Assessment of Exposure to Particulate Emission from Surface Contamination Sites. EPA/OHEA/EPA. 600/8-85/002. Cowherd, C., Jr., G.E. Muleski, P.J. Engelhart and D.A. Gillett, Ed. Washington DC: Midwest Research Inst.. 1985.

US EPA (United States Environmental Protection Agency). "Guidelines for Carcinogenic Risk Assessment." Federal Register 51:33992-34003. 1986.

US EPA (United States Environmental Protection Agency). Superfund Exposure Assessment Manual (SEAM). EPA 540/1-88/001. Office of Remedial Response. Washington, DC, 1988.

US EPA (United States Environmental Protection Agency). Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (Part A). Interim Final. EPA/540/1-89/002. Office of Emergency and Remedial Response. Washington, DC, 1989.

US EPA (United States Environmental Protection Agency). Human Health Evaluation Manual, Supplemental Guidance: Standard Default Exposure Factors. OSWER Directive. 9285.6-03. Office of Solid Waste and Emergency Response. Washington, DC, 1991.

US EPA (United States Environmental Protection Agency). Dermal Exposure Assessment: Principles and Applications. Office of Research and Development. EPA/600/8-91/011B. Washington, DC, 1992.

US EPA (United States Environmental Protection Agency). Estimating Exposure to Dioxin-Like Compounds. Office of Research and Development. EPA/600/6-88/005B. Washington, DC, 1992.

US EPA (United States Environmental Protection Agency). Provisional Guidance for Quantitative Risk Assessment of Polynuclear Aromatic Hydrocarbons. Office of Solid Waste and Environmental Remediation. EPA/600/R-93/089, July 1993

US EPA (United States Environmental Protection Agency). Land Use Directive in the CERCLA Remedy Selection Process. OSWER Directive 9355.7-04. Office of Solid Waste and Emergency Response. Washington, DC, May 1995.

US EPA (United States Environmental Protection Agency) Region 3. Technical Guidance Manual: Risk Assessment, Assessing Dermal Exposure From Soil. EPA/903-K-95-003. Office of Superfund Programs, Hazardous Waste Management Division. Washington, DC, 1995.

US EPA (United States Environmental Protection Agency) Region 4. Technical Services Supplemental Guidance to RAGS: Region 4 Bulletins. Waste Management Division, Atlanta, GA. 1995.

US EPA (United States Environmental Protection Agency). Supplement B to Compilation of Air Pollutant Emission Factors, Volume I: Stationary Point and Area Sources. AP-42, Fifth Edition, Supplement B. Office of Air Quality Planning and Standards, Office of Air and Radiation. Research Triangle Park, NC, 1996

US EPA (United States Environmental Protection Agency). Exposure Factors Handbook., EPA/600/P-95/002F. Office of Research and Development, Washington, DC, August 1997.

US EPA (United States Environmental Protection Agency). Health Effects Assessment Summary Tables (HEAST). Office of Health and Environmental Assessment, Environmental Criteria and Assessment Office (ECAO). Cincinnati OH, 1997.

US EPA (United States Environmental Protection Agency). IRIS (Integrated Risk Information System). A Continuously Updated Electronic Database Maintained by the US Environmental Protection Agency. Bethesda, Maryland: National Library of Medicine, 1999.

US EPA (United States Environmental Protection Agency) Region 3. Updated Risk-Based Concentration Tables. Office of RCRA Technical & Program Support Branch. Philadelphia, PA, April, 1999.

Wallcave, L., H. Garcia, R. Fedlman, W. Linjinsky, and P. Shubik. 1971. Skin tumorigenesis in mice by petroleum asphalts and coal-tar pitches of known polynuclear aromatic hydrocarbon content. Toxicol. Appl. Pharmacol. 18, 41-52.

Youngblood, W. W., and M. Blumer. 1975. Polycyclic aromatic hydrocarbons in the environment: homologous series in soils and recent marine sediments. *Geochim. Cosmochim. Acta* 39:1303-1315.

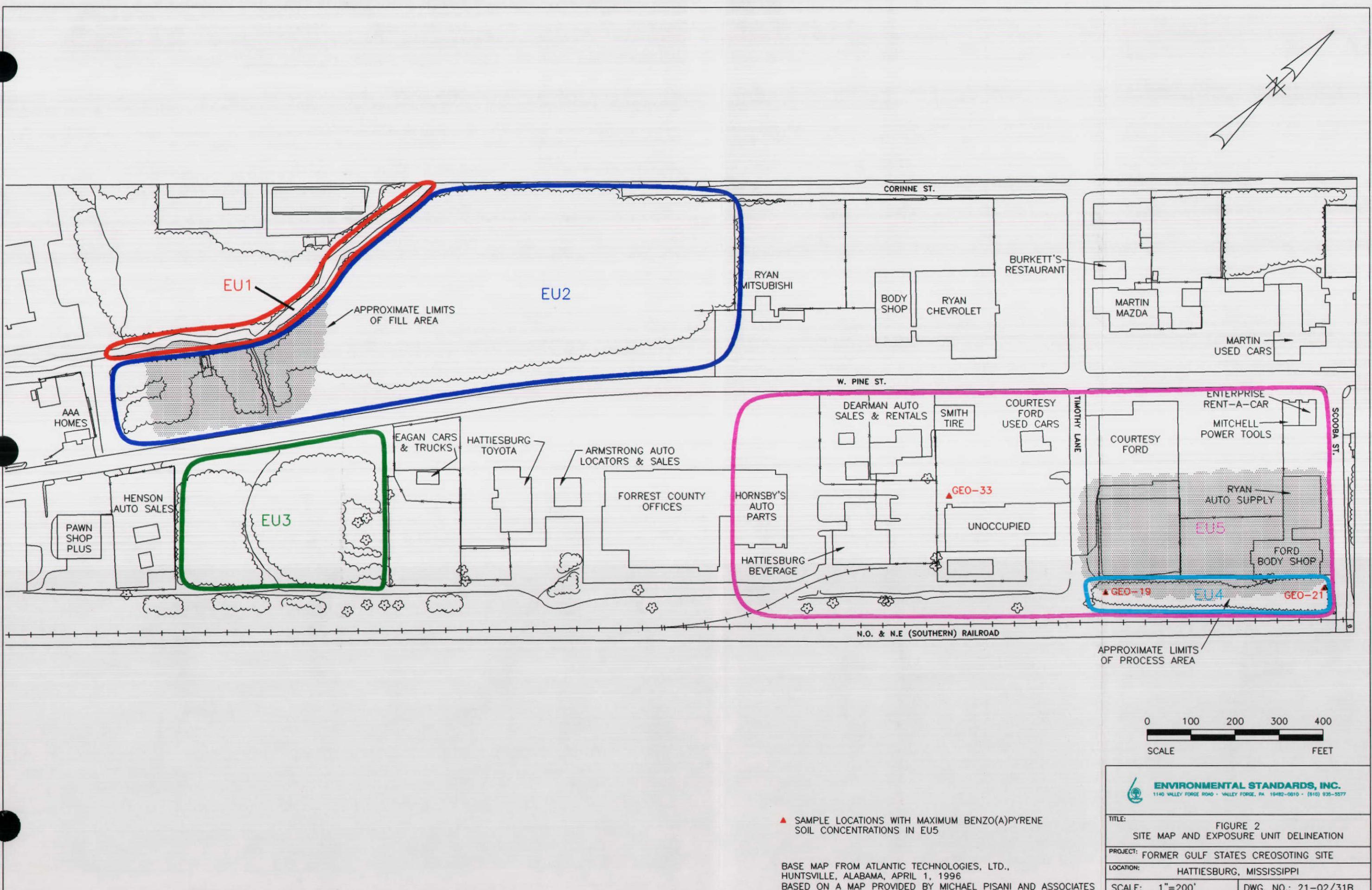
**Figure 1**  
**Site Conceptual Model and Selection of Exposure Pathways**  
**Kerr McGee, Hattiesburg, MS**

| Scenario      | Medium        | Exposure Medium | Exposure Point     | Receptor Population | Receptor Age | Exposure Route | On-Site/ Off-Site        | Type of Analysis  | Rationale for Selection or Exclusion of Exposure Pathway   |
|---------------|---------------|-----------------|--------------------|---------------------|--------------|----------------|--------------------------|---|--|
| Current       | Surface Soil  | Surface Soil    | Exposure Unit 1    | Visitor             | Adolescent   | Dermal Oral    | On-Site                  | None  | Surface Soils not included in EU1 data   |
|               |               |                 | Exposure Unit 2    | Visitor             | Adolescent   | Dermal Oral    | On-Site                  | None  | Surface Soils not included in EU1 data   |
|               |               |                 | Exposure Unit 3    | Visitor             | Adolescent   | Dermal Oral    | On-Site                  | Quantitative Qualitative  | Area potentially attractive for occasional recreational use PAHs to be addressed qualitatively according to MDDEQ guidance |
|               |               |                 | Exposure Unit 4    | Visitor             | Adolescent   | Dermal Oral    | On-Site                  | None  | VOCs not present at levels of concern  |
|               |               |                 | Exposure Unit 5    | Visitor             | Adolescent   | Dermal Oral    | On-Site                  | Quantitative Qualitative  | Area potentially attractive for occasional recreational use PAHs to be addressed qualitatively according to MDDEQ guidance |
|               |               |                 | Exposure Unit 5    | Maintenance Worker  | Adult        | Dermal Oral    | On-Site                  | None  | VOCs not present at levels of concern  |
| Sediment      | Sediment      | Exposure Unit 1 | Visitor            | Adolescent          | Dermal Oral  | On-Site        | Quantitative Qualitative | Exposed ground around drainage ditch potentially contacted by a visitor COPCs eliminated during screening process |  |
|               |               | Exposure Unit 4 | Visitor            | Adolescent          | Dermal Oral  | On-Site        | None                     | None  | Surface Soil exposures addressed in EU4, remainder of EU paved   |
|               |               | Exposure Unit 4 | Maintenance Worker | Adult               | Dermal Oral  | On-Site        | None                     | None  | Surface Soil exposures addressed in EU4, remainder of EU paved   |
|               |               | Exposure Unit 4 | Visitor            | Adolescent          | Dermal Oral  | On-Site        | Quantitative Qualitative | Potential exposure during digging-type activities PAHs to be addressed qualitatively according to MDDEQ guidance  |  |
|               |               | Exposure Unit 4 | Maintenance Worker | Adult               | Dermal Oral  | On-Site        | None                     | None  | VOCs not present at levels of concern  |
|               |               | Exposure Unit 4 | Visitor            | Adolescent          | Dermal Oral  | On-Site        | Quantitative Qualitative | Visitor may potentially wade in Gordon's Creek Published RfD values not available                                 |  |
|               |               | Exposure Unit 4 | Maintenance Worker | Adult               | Dermal Oral  | On-Site        | Quantitative Qualitative | Visitor may potentially walk through drainage ditch VOCs not present at levels of concern                         |  |
|               |               | Exposure Unit 4 | Visitor            | Adolescent          | Dermal Oral  | On-Site        | Quantitative Qualitative | Visitor may potentially walk through drainage ditch VOCs not present at levels of concern                         |  |
| Surface Water | Surface Water | Exposure Unit 1 | Visitor            | Adolescent          | Dermal Oral  | On-Site        | Quantitative Qualitative | Outside of maintained, fenced areas Exposure pathway insignificant  |  |
|               |               | Exposure Unit 4 | Visitor            | Adolescent          | Dermal Oral  | On-Site        | Quantitative Qualitative | Visitor may potentially walk through drainage ditch VOCs not present at levels of concern                         |  |
|               |               | Exposure Unit 4 | Maintenance Worker | Adult               | Dermal Oral  | On-Site        | None                     | Outside of maintained, fenced areas   |  |
|               |               | Exposure Unit 4 | Maintenance Worker | Adult               | Dermal Oral  | On-Site        | None                     | Outside of maintained, fenced areas   |  |
|               |               | Exposure Unit 4 | Visitor            | Adolescent          | Dermal Oral  | On-Site        | Quantitative Qualitative | Visitor may potentially wade in Gordon's Creek Exposure pathway insignificant                                     |  |

Figure 1

**Site Conceptual Model and Selection of Exposure Pathways**  
**Kerr McGee, Hattiesburg, MS**

| Scenario<br>Timeframe | Medium             | Exposure<br>Medium | Exposure<br>Point      | Receptor<br>Population | Receptor<br>Age              | Exposure<br>Route            | On-Site/<br>Off-Site                | Type of<br>Analysis  | Rationale for Selection or Exclusion<br>of Exposure Pathway  |
|-----------------------|--------------------|--------------------|------------------------|------------------------|------------------------------|------------------------------|-------------------------------------|--|--|
| Future                | Surface<br>Soil    | Surface<br>Soil    | Exposure Unit 1        | Maintenance<br>Worker  | Adult                        | Dermal<br>Oral<br>Inhalation | On-Site                             | None   | Surface Soils not included in EU1 data   |
|                       |                    |                    | Exposure Unit 2        | Maintenance<br>Worker  | Adult                        | Dermal<br>Oral<br>Inhalation | On-Site                             | None   | Surface Soils not included in EU1 data   |
|                       |                    |                    | Exposure Unit 3        | Maintenance<br>Worker  | Adult                        | Dermal<br>Oral<br>Inhalation | On-Site                             | Quantitative<br>Qualitative<br>None  | May potentially become a maintained area<br>PAHs to be addressed qualitatively according to MDEQ guidance<br>VOCs not present at levels of concern   |
|                       |                    |                    | Exposure Unit 4        | Maintenance<br>Worker  | Adult                        | Dermal<br>Oral<br>Inhalation | On-Site                             | Quantitative<br>Qualitative<br>None  | May potentially become a maintained area<br>PAHs to be addressed qualitatively according to MDEQ guidance<br>VOCs not present at levels of concern   |
|                       | Subsurface<br>Soil | Exposure Unit 1    | Construction<br>Worker | Adult                  | Dermal<br>Oral<br>Inhalation | On-Site                      | None                                | Surface Soil exposures addressed in EU5 under a current scenario   | Surface Soil exposures addressed in EU5 under a current scenario   |
|                       |                    | Exposure Unit 2    | Construction<br>Worker | Adult                  | Dermal<br>Oral<br>Inhalation | On-Site                      | None                                | Surface Soil exposures addressed in EU5 under a current scenario   | Surface Soil exposures addressed in EU5 under a current scenario   |
|                       |                    | Exposure Unit 3    | Construction<br>Worker | Adult                  | Dermal<br>Oral<br>Inhalation | On-Site                      | None                                | Potentially constructable area not constructable   | Potentially constructable area not constructable   |
|                       |                    | Exposure Unit 4    | Construction<br>Worker | Adult                  | Dermal<br>Oral<br>Inhalation | On-Site                      | Quantitative<br>Qualitative<br>None | Potentially constructable area in the future<br>PAHs to be addressed qualitatively according to MDEQ guidance<br>VOCs not present at levels of concern | Potentially constructable area in the future<br>PAHs to be addressed qualitatively according to MDEQ guidance<br>VOCs not present at levels of concern   |
|                       |                    | Exposure Unit 5    | Construction<br>Worker | Adult                  | Dermal<br>Oral<br>Inhalation | On-Site                      | Quantitative<br>Qualitative<br>None | Potentially constructable area in the future<br>COPCs eliminated during screening process  | Potentially constructable area in the future<br>COPCs eliminated during screening process  |
|                       |                    |                    |                        |                        |                              |                              |                                     | Quantitative<br>Qualitative<br>None  | Potential for non-VOC entrained fugitive dust generation during potential construction activities  |
|                       |                    |                    |                        |                        |                              |                              |                                     | None   | Soil exposures addressed in EU5  |
|                       |                    |                    |                        |                        |                              |                              |                                     | None   | Soil exposures addressed in EU5  |
|                       |                    |                    |                        |                        |                              |                              |                                     | None   | Soil exposures addressed in EU5  |
|                       |                    |                    |                        |                        |                              |                              |                                     | Quantitative<br>Qualitative<br>Quantitative  | Potentially constructable area in the future<br>PAHs to be addressed qualitatively according to MDEQ guidance<br>Potential for non-VOC entrained fugitive dust generation during potential construction activities |
|                       |                    |                    |                        |                        |                              |                              |                                     | Quantitative<br>Qualitative<br>Quantitative  | Potentially constructable area in the future<br>PAHs to be addressed qualitatively according to MDEQ guidance<br>Potential for non-VOC entrained fugitive dust generation during potential construction activities |



## Table

Statistical Summary and Selection of COPCs in EU1 Sediment  
Kerr McGee, Hattiesburg, MS

| Constituent                 | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Minimum Detected mg/kg | Detected Qualifier | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Detected Qualifier | Maximum Concentration mg/kg | Standard Deviation mg/kg |
|-----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|------------------------|--------------------|------------|------------------------|------------------------|--------------------|-----------------------------|--------------------------|
| <b>Semivolatiles</b>        |            |               |                |                 |               |               |                       |                        |                    |            |                        |                        |                    |                             |                          |
| 2-Methylnaphthalene         | 91-37-6    | 2             | 2              | 100             | NA            | NA            | 7.40E-02              | J                      | 2.92E-01           | 1.94E-01   | 5.10E-01               | SD-07                  | 3.08E-01           |                             |                          |
| Acenaphthene                | 83-32-9    | 2             | 2              | 100             | NA            | NA            | 1.80E-01              | J                      | 3.15E-01           | 2.85E-01   | 4.50E-01               | SD-07                  | 1.91E-01           |                             |                          |
| Acenaphthylene              | 208-96-8   | 2             | 1              | 50              | 4.00E-02      | 4.00E-02      | 7.80E-02              | J                      | 4.90E-02           | 3.95E-02   | 7.80E-02               | J                      | SD-07              | 4.10E-02                    |                          |
| Anthracene                  | 120-12-7   | 2             | 2              | 100             | NA            | NA            | 2.60E-01              | J                      | 3.60E-01           | 3.46E-01   | 4.60E-01               | J                      | SD-07              | 1.41E-01                    |                          |
| Benzo (a) Pyrene Equivalent | -          | 2             | 2              | 100             | NA            | NA            | 1.83E-01              |                        | 3.88E-01           | 3.29E-01   | 5.93E-01               |                        | SD-07              | 2.90E-01                    |                          |
| Benzo(ghi)perylene          | 191-24-2   | 2             | 2              | 100             | NA            | NA            | 6.50E-02              | J                      | 1.23E-01           | 1.08E-01   | 1.80E-01               | J                      | SD-07              | 8.13E-02                    |                          |
| Carbazole                   | 86-74-8    | 2             | 2              | 100             | NA            | NA            | 1.60E-01              | J                      | 3.65E-01           | 3.02E-01   | 5.70E-01               | J                      | SD-07              | 2.90E-01                    |                          |
| Dibenzofuran                | 132-64-9   | 2             | 2              | 100             | NA            | NA            | 1.50E-01              | J                      | 2.80E-01           | 2.48E-01   | 4.10E-01               | J                      | SD-07              | 1.84E-01                    |                          |
| Fluoranthene                | 206-44-0   | 2             | 2              | 100             | NA            | NA            | 6.80E-01              |                        | 1.19E+00           | 1.08E+00   | 1.70E+00               |                        | SD-07              | 7.21E-01                    |                          |
| Fluorene                    | 86-73-7    | 2             | 2              | 100             | NA            | NA            | 2.30E-01              | J                      | 4.25E-01           | 3.78E-01   | 6.20E-01               | J                      | SD-07              | 2.76E-01                    |                          |
| Naphthalene                 | 91-20-3    | 2             | 2              | 100             | NA            | NA            | 1.80E-01              | J                      | 6.40E-01           | 4.45E-01   | 1.10E+00               | J                      | SD-07              | 6.51E-01                    |                          |
| Phenanthrene                | 85-01-8    | 2             | 2              | 100             | NA            | NA            | 7.20E-01              | 1.21E+00               | 1.11E+00           | 1.70E+00   | 1.70E+00               | SD-07                  | 6.93E-01           |                             |                          |
| Pyrene                      | 129-00-0   | 2             | 2              | 100             | NA            | NA            | 4.80E-01              | 9.40E-01               | 8.20E-01           | 1.40E+00   | 1.40E+00               | SD-07                  | 6.51E-01           |                             |                          |

NA - Not Available

**Table 1**  
**Statistical Summary and Selection of COPCs in EU1 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent          | 95% UCL          |                  | Lognormal<br>95% UCL<br>mg/kg |                  | Distribution<br>99%<br>Confidence | Exposure Point<br>Concentration<br>mg/kg | Region III<br>Residential Soil |              | Is the Maximum<br>Detected > RBC? |
|----------------------|------------------|------------------|-------------------------------|------------------|-----------------------------------|--|--------------------------------|--------------|-----------------------------------|
|                      | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg | 95% UCL<br>mg/kg              | 95% UCL<br>mg/kg |                                   |  | 95% UCL<br>mg/kg               | RBC<br>mg/kg |                                   |
| <b>Semivolatiles</b> |                  |                  |                               |                  |                                   |  |                                |              |                                   |
| 2-Methylnaphthalene  | 1.67E+00         | 1.60E+22         | Unknown                       | 5.10E-01         | 1.60E+02                          | no                                       |                                |              |                                   |
| Acenaphthene         | 1.17E+00         | 3.23E+04         | Unknown                       | 4.50E-01         | 4.70E+02                          | no                                       |                                |              |                                   |
| Acenaphthylene*      | 2.32E-01         | 8.34E+09         | Unknown                       | 7.80E-02         | NA                                | no                                       |                                |              |                                   |
| Anthracene           | 9.91E-01         | 2.23E+01         | Unknown                       | 4.60E-01         | 2.30E+03                          | no                                       |                                |              |                                   |
| Benz(a)Pyrene Equiv. | 1.68E+00         | 9.30E+07         | Unknown                       | 5.93E-01         | 8.70E-02                          | YES - COPC                               |                                |              |                                   |
| Benz(ghi)perylene*   | 4.86E-01         | 2.08E+05         | Unknown                       | 1.80E-01         | NA                                | no                                       |                                |              |                                   |
| Carbazole            | 1.66E+00         | 2.15E+09         | Unknown                       | 5.70E-01         | 3.20E+01                          | no                                       |                                |              |                                   |
| Dibenzofuran         | 1.10E+00         | 3.27E+05         | Unknown                       | 4.10E-01         | 3.10E+01                          | no                                       |                                |              |                                   |
| Fluoranthene         | 4.41E+00         | 1.22E+05         | Unknown                       | 1.70E+00         | 3.10E+02                          | no                                       |                                |              |                                   |
| Fluorene             | 1.66E+00         | 3.35E+05         | Unknown                       | 6.20E-01         | 3.10E+02                          | no                                       |                                |              |                                   |
| Naphthalene          | 3.54E+00         | 6.10E+19         | Unknown                       | 1.10E+00         | 1.60E+02                          | no                                       |                                |              |                                   |
| Phenanthrene*        | 4.30E+00         | 2.92E+04         | Unknown                       | 1.70E+00         | NA                                | no                                       |                                |              |                                   |
| Pyrene               | 3.84E+00         | 7.40E+06         | Unknown                       | 1.40E+00         | 2.30E+02                          | no                                       |                                |              |                                   |

NA - Not Available

\* Constituent will be retained as a COPC due to lack of published screening criteria.

**Table 2**  
**Statistical Summary and Selection of COPCs in EUI Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent             | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum Detection Limit mg/L | Maximum Detection Limit mg/L | Minimum Detected Qualifier mg/L | Maximum Detected Qualifier mg/L | Mean mg/L | Logarithmic Mean mg/L | Maximum Detected Qualifier mg/L | Maximum Concentration | Location Maximum Concentration | Standard Deviation mg/L |
|-------------------------|------------|---------------|----------------|-----------------|------------------------------|------------------------------|---------------------------------|---------------------------------|-----------|-----------------------|---------------------------------|-----------------------|--------------------------------|-------------------------|
|                         |            |               |                |                 |                              |                              |                                 |                                 |           |                       |                                 |                       |                                |                         |
| <b>Semi挥发性</b>          |            |               |                |                 |                              |                              |                                 |                                 |           |                       |                                 |                       |                                |                         |
| Benzo (a) Pyrene Equiv. | -          | 2             | 1              | 50              | NA                           | NA                           | 1.00E-03                        | 1.21E-03                        | 1.18E-03  | 1.21E-03              | 1.18E-03                        | SW-08                 | 3.54E-05                       |                         |
| Fluoranthene            | 206-44-0   | 2             | 1              | 50              | 1.00E-03                     | 1.00E-03                     | 7.50E-03                        | 7.50E-03                        | 1.94E-03  | 7.50E-03              | 1.94E-03                        | SW-08                 | 4.95E-03                       |                         |
| Pyrene                  | 129-09-0   | 2             | 1              | 50              | 1.00E-03                     | 1.00E-03                     | 1.00E-03                        | 1.00E-03                        | J         | 7.50E-04              | 7.07E-04                        | 1.00E-03              | J                              | 3.54E-04                |

NA - Not Available



**Table 2**  
**Statistical Summary and Selection of COPCs in EU1 Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/L | Lognormal<br>95% UCL<br>mg/L | Distribution<br>95%<br>Confidence | Exposure Point<br>Concentration<br>mg/L | Human Health<br>Organisms AWQC<br>mg/L | Consumption of Water &<br>Organisms AWQC<br>mg/L | Is the Maximum<br>Concentration > the<br>AWQC? |
|------------------------|-----------------|------------------------------|-----------------------------------|---|--|--|--|
|                        |                 |                              |                                   |   |  |  |  |
| Semivolatiles          |                 |                              |                                   |   |  |  |  |
| Benz (a) Pyrene Equiv. | 1.34E-03        | 1.31E-03                     | Unknown                           | 1.21E-03                                | 4.40E-06                               | YES - COPC                                       |  |
| Fluoranthene           | 2.61E-02        | 2.90E+42                     | Unknown                           | 7.50E-03                                | 3.00E-01                               | no   |  |
| Pyrene*                | 2.33E-03        | 4.37E-01                     | Unknown                           | 1.00E-03                                | NA                                     | no   |  |

NA - Not Available

\* Constituent will be retained as a COPC due to lack of published screening criteria.

**Table**

**Statistical Summary and Selection of COPCs in EU4 Sediment  
Kerr McGee, Hattiesburg, MS**

| Constituent             | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum Detection Limit mg/kg | Maximum Detection Limit mg/kg | Minimum Detected Qualifier mg/kg | Maximum Detected Qualifier mg/kg | Logarithmic Mean mg/kg | Maximum Detected Qualifier mg/kg | Maximum Detected Qualifier mg/kg | Location of Concentration on |
|-------------------------|------------|---------------|----------------|-----------------|-------------------------------|-------------------------------|----------------------------------|----------------------------------|------------------------|----------------------------------|----------------------------------|------------------------------|
| <b>Semivolatiles</b>    |            |               |                |                 |                               |                               |                                  |                                  |                        |                                  |                                  |                              |
| 2,4-Dimethylphenol      | 105-67-9   | 1             | 1              | 100             | NA                            | NA                            | 1.50E+00                         | J                                | 1.50E+00               | 1.50E+00                         | J                                | SD-02                        |
| 2-Methylnaphthalene     | 91-57-6    | 1             | 1              | 100             | NA                            | NA                            | 1.50E+03                         | 1.50E+03                         | 1.50E+03               | 1.50E+03                         | SD-02                            | 0.00E+00                     |
| Acenaphthylene          | 208-96-8   | 1             | 1              | 100             | NA                            | NA                            | 3.50E+01                         | 3.50E+01                         | 3.50E+01               | 3.50E+01                         | SD-02                            | 0.00E+00                     |
| Anthracene              | 120-12-7   | 1             | 1              | 100             | NA                            | NA                            | 1.90E+03                         | 1.90E+03                         | 1.90E+03               | 1.90E+03                         | SD-02                            | 0.00E+00                     |
| Benzo (a) Pyrene Equiv. | -          | 1             | 1              | 100             | NA                            | NA                            | 1.99E+02                         | 1.99E+02                         | 1.99E+02               | 1.99E+02                         | SD-02                            | 0.00E+00                     |
| Benzo(g,h,i)perylene    | 191-24-2   | 1             | 1              | 100             | NA                            | NA                            | 3.60E+01                         | 3.60E+01                         | 3.60E+01               | 3.60E+01                         | SD-02                            | 0.00E+00                     |
| Carbazole               | 86-74-8    | 1             | 1              | 100             | NA                            | NA                            | 5.90E+02                         | 5.90E+02                         | 5.90E+02               | 5.90E+02                         | SD-02                            | 0.00E+00                     |
| Dibenzofuran            | 132-64-9   | 1             | 1              | 100             | NA                            | NA                            | 9.40E+02                         | 9.40E+02                         | 9.40E+02               | 9.40E+02                         | SD-02                            | 0.00E+00                     |
| Fluoranthene            | 206-44-0   | 1             | 1              | 100             | NA                            | NA                            | 1.60E+03                         | 1.60E+03                         | 1.60E+03               | 1.60E+03                         | SD-02                            | 0.00E+00                     |
| Fluorene                | 86-73-7    | 1             | 1              | 100             | NA                            | NA                            | 1.20E+03                         | 1.20E+03                         | 1.20E+03               | 1.20E+03                         | SD-02                            | 0.00E+00                     |
| Naphthalene             | 91-20-3    | 1             | 1              | 100             | NA                            | NA                            | 3.00E+03                         | 3.00E+03                         | 3.00E+03               | 3.00E+03                         | SD-02                            | 0.00E+00                     |
| Phenanthrene            | 85-01-8    | 1             | 1              | 100             | NA                            | NA                            | 3.20E+03                         | 3.20E+03                         | 3.20E+03               | 3.20E+03                         | SD-02                            | 0.00E+00                     |
| Pyrene                  | 129-00-0   | 1             | 1              | 100             | NA                            | NA                            | 1.00E+03                         | 1.00E+03                         | 1.00E+03               | 1.00E+03                         | SD-02                            | 0.00E+00                     |

NA - Not Available

**Table 3**  
**Statistical Summary and Selection of COPCs in EU4 Sediment**  
**Kerr McGee, Hattiesburg, MS**

| Constituent             | 95% UCL |       | Lognormal 95% UCL |         | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Residential Soil<br>RBC<br>mg/kg | Is the Maximum<br>Detected > RBC? |
|-------------------------|---------|-------|-------------------|---------|--------------------------------|--|----------------------------------|-----------------------------------|
|                         | mg/kg   | mg/kg | mg/kg             | mg/kg   |                                |  |                                  |                                   |
| <b>Semivolatile</b>     |         |       |                   |         |                                |  |                                  |                                   |
| 2,4-Dimethylphenol      | NA      | NA    | NA                | Unknown | 1.50E+00                       | 1.60E+02                                 | no                               |                                   |
| 2-Methylnaphthalene     | NA      | NA    | NA                | Unknown | 1.50E+03                       | 1.60E+02                                 | YES - COPC                       |                                   |
| Acenaphthylene*         | NA      | NA    | NA                | Unknown | 3.50E+01                       | NA                                       | no                               |                                   |
| Anthracene              | NA      | NA    | NA                | Unknown | 1.90E+03                       | 2.30E+03                                 | no                               |                                   |
| Benzo (a) Pyrene Equiv. | NA      | NA    | NA                | Unknown | 1.99E+02                       | 8.70E-02                                 | YES - COPC                       |                                   |
| Benzo(g,h,i)perylene*   | NA      | NA    | NA                | Unknown | 3.60E+01                       | NA                                       | no                               |                                   |
| Carbazole               | NA      | NA    | NA                | Unknown | 5.90E+02                       | 3.20E+01                                 | YES - COPC                       |                                   |
| Dibenzofuran            | NA      | NA    | NA                | Unknown | 9.40E+02                       | 3.10E+01                                 | YES - COPC                       |                                   |
| Fluoranthene            | NA      | NA    | NA                | Unknown | 1.60E+03                       | 3.10E+02                                 | YES - COPC                       |                                   |
| Fluorene                | NA      | NA    | NA                | Unknown | 1.20E+03                       | 3.10E+02                                 | YES - COPC                       |                                   |
| Naphthalene             | NA      | NA    | NA                | Unknown | 3.00E+03                       | 1.60E+02                                 | YES - COPC                       |                                   |
| Phenanthrene*           | NA      | NA    | NA                | Unknown | 3.20E+03                       | NA                                       | no                               |                                   |
| Pyrene                  | NA      | NA    | NA                | Unknown | 1.00E+03                       | 2.30E+02                                 | YES - COPC                       |                                   |

NA - Not Available

\* Constituent will be retained as a COPC due to lack of published screening criteria.

Table 4

Statistical Summary and Selection of COPCs in EU4 Surface Water  
Kerr McGee, Hattiesburg, MS

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/L | Maximum mg/L | Detection Limit mg/L | Detected Qualifier | Minimum mg/L | Maximum mg/L | Detected Qualifier | Logarithmic Mean mg/L | Logarithmic Mean mg/L | Maximum Detected Qualifier | Maximum Detected Qualifier | Location Maximum Concentration mg/L | Standard Deviation mg/L |
|----------------------------|------------|---------------|----------------|-----------------|--------------|--------------|----------------------|--------------------|--------------|--------------|--------------------|-----------------------|-----------------------|----------------------------|----------------------------|-------------------------------------|-------------------------|
| <b>Semivolatiles</b>       |            |               |                |                 |              |              |                      |                    |              |              |                    |                       |                       |                            |                            |                                     |                         |
| Acenaphthene               | 83-32-9    | 1             | 1              | 100             | NA           | NA           | 1.40E-02             | 1.40E-02           | 1.40E-02     | 1.40E-02     | 1.40E-02           | SW-02                 | 0.00E+00              |                            |                            |                                     |                         |
| Antracene                  | 120-12-7   | 1             | 1              | 100             | NA           | NA           | 1.30E-02             | 1.30E-02           | 1.30E-02     | 1.30E-02     | 1.30E-02           | SW-02                 | 0.00E+00              |                            |                            |                                     |                         |
| Benzo (a) Pyrene Equiv.    | -          | 1             | 1              | 100             | NA           | NA           | 2.78E-03             | 2.78E-03           | 2.78E-03     | 2.78E-03     | 2.78E-03           | SW-02                 | 0.00E+00              |                            |                            |                                     |                         |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 1             | 1              | 100             | NA           | NA           | 3.00E-03             | J                  | 3.00E-03     | 3.00E-03     | 3.00E-03           | J                     | SW-02                 | 0.00E+00                   |                            |                                     |                         |
| Carbazole                  | 86-74-8    | 1             | 1              | 100             | NA           | NA           | 1.00E-02             | J                  | 1.00E-02     | 1.00E-02     | 1.00E-02           | J                     | SW-02                 | 0.00E+00                   |                            |                                     |                         |
| Dibenzofuran               | 132-64-9   | 1             | 1              | 100             | NA           | NA           | 1.10E-02             | 1.10E-02           | 1.10E-02     | 1.10E-02     | 1.10E-02           | SW-02                 | 0.00E+00              |                            |                            |                                     |                         |
| Fluoranthene               | 206-44-0   | 1             | 1              | 100             | NA           | NA           | 3.90E-02             | 3.90E-02           | 3.90E-02     | 3.90E-02     | 3.90E-02           | SW-02                 | 0.00E+00              |                            |                            |                                     |                         |
| Fluorene                   | 86-73-7    | 1             | 1              | 100             | NA           | NA           | 1.20E-02             | 1.20E-02           | 1.20E-02     | 1.20E-02     | 1.20E-02           | SW-02                 | 0.00E+00              |                            |                            |                                     |                         |
| Phenanthrene               | 85-01-8    | 1             | 1              | 100             | NA           | NA           | 1.70E-02             | 1.70E-02           | 1.70E-02     | 1.70E-02     | 1.70E-02           | SW-02                 | 0.00E+00              |                            |                            |                                     |                         |
| Pyrene                     | 129-00-0   | 1             | 1              | 100             | NA           | NA           | 2.10E-02             | 2.10E-02           | 2.10E-02     | 2.10E-02     | 2.10E-02           | SW-02                 | 0.00E+00              |                            |                            |                                     |                         |

NA - Not Available

**Table 4**  
**Statistical Summary and Selection of COPCs in EU4 Surface Water**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | 95% UCL<br>mg/L | Lognormal<br>95% UCL<br>mg/L | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/L | Human Health Consumption of<br>Water & Organisms AWQC<br>mg/L | Is the Maximum<br>Concentration > the<br>AWQC? |
|----------------------------|-----------------|------------------------------|--------------------------------|---|---|--|
| <b>SemiVolatiles</b>       |                 |                              |                                |   |   |  |
| Acenaphthene               | NA              | NA                           | Unknown                        | 1.40E-02                                | 1.20E+00  | no   |
| Anthracene                 | NA              | NA                           | Unknown                        | 1.30E-02                                | 9.60E+00  | no   |
| Benzo (a) Pyrene Equiv.    | NA              | NA                           | Unknown                        | 2.78E-03                                | 4.40E-06  | YES - COPC                                     |
| bis(2-ethylhexyl)phthalate | NA              | NA                           | Unknown                        | 3.00E-03                                | 1.80E-03  | YES - COPC                                     |
| Carbazole*                 | NA              | NA                           | Unknown                        | 1.00E-02                                | NA  | no   |
| Dibenzofuran*              | NA              | NA                           | Unknown                        | 1.10E-02                                | NA  | no   |
| Fluoranthene               | NA              | NA                           | Unknown                        | 3.90E-02                                | 3.00E-01  | no   |
| Fluorene                   | NA              | NA                           | Unknown                        | 1.20E-02                                | 1.30E+00  | no   |
| Phenanthrene*              | NA              | NA                           | Unknown                        | 1.70E-02                                | NA  | no   |
| Pyrene                     | NA              | NA                           | Unknown                        | 2.10E-02                                | 9.60E-01  | no   |

NA - Not Available

\* Constituent will be retained as a COPC due to lack of published screening criteria.

**Table 5**  
**Statistical Summary and Selection of COPCs in EU2 Surface Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent             | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected mg/kg | Minimum Detected Qualifier | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Detected Qualifier | Maximum Concentration | Location of Maximum Concentration | Standard Deviation mg/kg |
|-------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------------|----------------------------|------------|------------------------|------------------------|--------------------|-----------------------|-----------------------------------|--------------------------|
| <b>Semi-volatiles</b>   |            |               |                |                 |               |               |                       |                |                            |            |                        |                        |                    |                       |                                   |                          |
| 2-Methylnaphthalene     | 91-57-6    | 14            | 2              | 14.29           | 3.30E-02      | 3.30E-02      | 7.00E-02              | J              | 3.06E-02                   | 2.15E-02   | 1.60E-01               | J                      | SS-10              | 3.99E-02              |                                   |                          |
| Acenaphthene            | 83-32-9    | 14            | 1              | 7.14            | 3.30E-02      | 3.30E-02      | 4.90E-02              | J              | 1.88E-02                   | 1.78E-02   | 4.90E-02               | J                      | GEO-13/0-1'        | 8.69E-03              |                                   |                          |
| Acenaphthylene          | 208-96-8   | 14            | 6              | 42.86           | 3.30E-02      | 3.30E-02      | 3.70E-02              | J              | 1.59E-01                   | 4.29E-02   | 1.30E+00               | J                      | GEO-13/0-1'        | 3.52E-01              |                                   |                          |
| Anthracene              | 120-12-7   | 14            | 7              | 50              | 3.30E-02      | 3.30E-02      | 4.10E-02              | J              | 1.89E-01                   | 5.00E-02   | 1.60E+00               | J                      | GEO-13/0-1'        | 4.28E-01              |                                   |                          |
| Benzo (a) Pyrene Equiv. | -          | 14            | 12             | 85.71           | NA            | NA            | 8.85E-02              | J              | 1.36E+00                   | 5.11E-01   | 8.11E+00               | J                      | GEO-13/0-1'        | 2.23E+00              |                                   |                          |
| Benzo(ghi)perylene      | 191-24-2   | 14            | 10             | 71.43           | 6.70E-02      | 6.70E-02      | 1.70E-01              | J              | 5.17E-01                   | 2.20E-01   | 2.30E+00               | J                      | SS-10              | 6.93E-01              |                                   |                          |
| Carbazole               | 86-74-8    | 14            | 4              | 28.57           | 3.30E-02      | 3.30E-02      | 4.30E-02              | J              | 6.28E-02                   | 2.94E-02   | 3.50E-01               | J                      | GEO-13/0-1'        | 1.03E-01              |                                   |                          |
| Di-n-butylphthalate     | 84-74-2    | 14            | 9              | 64.29           | 3.30E-02      | 3.30E-02      | 7.20E-02              | J              | 3.60E-02                   | 3.68E-02   | 1.10E-01               | J                      | SS-10              | 2.50E-02              |                                   |                          |
| Dibenzofuran            | 132-64-9   | 14            | 2              | 14.29           | 3.30E-02      | 3.30E-02      | 7.20E-02              | J              | 2.63E-02                   | 2.08E-02   | 9.80E-02               | J                      | SS-10              | 2.54E-02              |                                   |                          |
| Fluoranthene            | 206-44-0   | 14            | 12             | 85.71           | 3.30E-02      | 3.30E-02      | 6.60E-02              | J              | 1.40E+00                   | 3.00E-01   | 1.20E+01               | J                      | GEO-13/0-1'        | 3.16E+00              |                                   |                          |
| Fluorene                | 86-73-7    | 14            | 2              | 14.29           | 3.30E-02      | 3.30E-02      | 4.50E-02              | J              | 4.38E-02                   | 2.21E-02   | 3.70E-01               | J                      | GEO-13/0-1'        | 9.42E-02              |                                   |                          |
| Naphthalene             | 91-20-3    | 14            | 2              | 14.29           | 3.30E-02      | 3.30E-02      | 8.80E-02              | J              | 3.26E-02                   | 2.20E-02   | 1.70E-01               | J                      | SS-10              | 4.39E-02              |                                   |                          |
| Phenanthrene            | 85-01-8    | 14            | 8              | 57.14           | 3.30E-02      | 3.30E-02      | 3.70E-02              | J              | 1.28E-01                   | 5.30E-02   | 7.40E-01               | J                      | GEO-13/0-1'        | 2.08E-01              |                                   |                          |
| Pyrene                  | 129-00-0   | 14            | 12             | 85.71           | 6.70E-02      | 6.70E-02      | 9.80E-02              | J              | 1.70E+00                   | 4.60E-01   | 1.40E+01               | J                      | GEO-13/0-1'        | 3.66E+00              |                                   |                          |

NA - Not Available

Table 5

**Statistical Summary and Selection of COPCs in EU2 Surface Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent             | 95% UCL<br>mg/kg | Lognormal 95%<br>UCL<br>mg/kg | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG?<br>no | Is the 95% UCL ><br>TRG? no |
|-------------------------|------------------|-------------------------------|--------------------------------|--|--|---|-----------------------------|
| <b>Semi-volatiles</b>   |                  |                               |                                |  |  |   |                             |
| 2-Methylnaphthalene     | 4.95E-02         | 4.29E-02                      | Unknown                        | 4.29E-02                                 | 8.18E-04                               | no                                      | no                          |
| Acenaphthene            | 2.29E-02         | 2.17E-02                      | Unknown                        | 2.17E-02                                 | 1.23E-05                               | no                                      | no                          |
| Acenaphthylene          | 3.26E-01         | 4.99E-01                      | Unknown                        | 4.99E-01                                 | 1.23E-05                               | no                                      | no                          |
| Anthracene              | 3.91E-01         | 6.29E-01                      | Unknown                        | 6.29E-01                                 | 6.13E-05                               | no                                      | no                          |
| Benzo (a) Pyrene Equiv. | 2.42E+00         | 6.37E+00                      | Lognormal                      | 6.37E+00                                 | 7.84E-01                               | YES                                     | YES - COPC                  |
| Benzo (g,h,i)perylene   | 8.46E-01         | 2.74E+00                      | Lognormal                      | 2.30E+00                                 | 6.13E-04                               | no                                      | no                          |
| Carbazole               | 1.12E-01         | 1.24E-01                      | Unknown                        | 1.24E-01                                 | 2.86E-02                               | no                                      | no                          |
| Di-n-butylphthalate     | 5.48E-02         | 6.30E-02                      | Normal/Lognormal               | 6.30E-02                                 | 2.28E-03                               | no                                      | no                          |
| Dibenzofuran            | 3.83E-02         | 3.57E-02                      | Unknown                        | 3.57E-02                                 | 8.18E-03                               | no                                      | no                          |
| Fluoranthene            | 2.89E+00         | 1.66E+01                      | Lognormal                      | 1.20E+01                                 | 8.17E-04                               | no                                      | no                          |
| Fluorene                | 8.84E-02         | 5.84E-02                      | Unknown                        | 5.84E-02                                 | 8.17E-04                               | no                                      | no                          |
| Naphthalene             | 5.34E-02         | 4.71E-02                      | Unknown                        | 4.71E-02                                 | 8.24E-02                               | no                                      | no                          |
| Phenanthrene            | 2.26E-01         | 3.96E-01                      | Lognormal                      | 3.96E-01                                 | 6.13E-04                               | no                                      | no                          |
| Pyrene                  | 3.43E+00         | 1.23E+01                      | Lognormal                      | 1.25E+01                                 | 6.13E-04                               | no                                      | no                          |

NA - Not Available

Table 6

Statistical Summary and Selection of COPCs in EU3 Surface Soil (0-1' bgs)  
Kerr McGee, Hattiesburg, MS

| Constituent             | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum Detection Limit mg/kg | Maximum Detection Limit mg/kg | Detected mg/kg | Minimum Detected Qualifier mg/kg | Maximum Detected Qualifier mg/kg | Logarithmic Mean mg/kg | Maximum Mean mg/kg | Detected mg/kg | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|-------------------------|------------|---------------|----------------|-----------------|-------------------------------|-------------------------------|----------------|----------------------------------|----------------------------------|------------------------|--------------------|----------------|-----------------------------|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>    |            |               |                |                 |                               |                               |                |                                  |                                  |                        |                    |                |                             |                                   |                          |
| 2-Methylphthalene       | 91-57-6    | 3             | 1              | 33.33           | 3.30E-02                      | 3.30E-02                      | J              | 8.77E-02                         | 3.97E-02                         | 2.30E-01               | J                  | SS-16          | 1.23E-01                    |                                   |                          |
| Benzo (a) Pyrene Equiv. | 208-95-8   | 3             | 3              | 100             | NA                            | 1.01E-01                      |                | 7.17E-01                         | 4.72E-01                         | 1.12E+00               |                    | SS-16          | 5.43E-01                    |                                   |                          |
| Acenaphthylene          | 120-12-7   | 3             | 2              | 66.67           | 3.30E-02                      | 3.30E-02                      | J              | 1.20E-01                         | 1.02E-01                         | 6.96E-02               | 1.70E-01           | J              | SS-16                       | 7.83E-02                          |                          |
| Anthracene              | 191-24-2   | 3             | 3              | 100             | NA                            | NA                            |                | 1.20E-01                         | J                                | 1.02E-01               | 6.96E-02           | J              | SS-16                       | 7.83E-02                          |                          |
| Benz(ghi)perylene       | 86-74-8    | 3             | 2              | 66.67           | 3.30E-02                      | 3.30E-02                      | J              | 8.00E-02                         | J                                | 6.53E-01               | 4.03E-01           | 1.20E+00       | SS-16                       | 5.60E-01                          |                          |
| Carbazole               | 84-74-2    | 3             | 3              | 100             | NA                            | NA                            |                | 4.60E-02                         | J                                | 5.75E-02               | 4.37E-02           | 1.10E-01       | J                           | SS-16                             | 4.78E-02                 |
| Di-n-butylphthalate     | 132-64-9   | 3             | 2              | 66.67           | 3.30E-02                      | 3.30E-02                      | J              | 4.00E-02                         | J                                | 8.30E-02               | 7.58E-02           | 1.10E-01       | J                           | SS-16                             | 3.76E-02                 |
| Dibenzofuran            | 206-44-0   | 3             | 3              | 100             | NA                            | NA                            |                | 3.60E-02                         | J                                | 4.85E-02               | 3.81E-02           | 9.30E-02       | J                           | SS-16                             | 3.98E-02                 |
| Fluoranthene            | 91-20-3    | 3             | 1              | 33.33           | 3.30E-02                      | 3.30E-02                      | J              | 1.20E-01                         | J                                | 5.27E-01               | 3.99E-01           | 7.80E-01       | J                           | SS-16                             | 3.56E-01                 |
| Naphthalene             | 85-01-8    | 3             | 2              | 66.67           | 3.30E-02                      | 3.30E-02                      | J              | 1.60E-01                         | J                                | 6.43E-02               | 3.52E-02           | 1.60E-01       | J                           | SS-16                             | 8.28E-02                 |
| Phenanthrene            | 129-00-0   | 3             | 3              | 100             | NA                            | NA                            |                | 1.30E-01                         | J                                | 1.32E-01               | 8.12E-02           | 2.50E-01       | J                           | SS-16                             | 1.17E-01                 |
| Pyrene                  |            |               |                |                 |                               |                               |                | 1.20E-01                         | J                                | 6.90E-01               | 4.85E-01           | 1.00E+00       |                             | SS-17                             | 4.94E-01                 |

NA - Not Available

**Table 6**  
**Statistical Summary and Selection of COPCs in EU3 Surface Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent           | 95% UCL<br>mg/kg | Lognormal<br>95% UCL<br>mg/kg | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRC<br>mg/kg | Is the Maximum<br>Detected > TRC? | Is the 95% UCL ><br>TRC? |
|-----------------------|------------------|-------------------------------|--------------------------------|--|--|-----------------------------------|--------------------------|
| Semivolatiles         |                  |                               |                                |  |  |                                   |                          |
| 2-Methylnaphthalene   | 2.95E-01         | 2.43E+08                      | Unknown                        | 2.30E-01                                 | 8.18E-04                               | no                                | no                       |
| Benz(a) Pyrene Equiv. | 1.63E+00         | 1.82E+07                      | Normal/Lognormal               | 1.12E+00                                 | 7.84E-01                               | no                                | YES - COPC               |
| Acenaphthylene        | 2.34E-01         | 3.45E+05                      | Normal/Lognormal               | 1.70E-01                                 | 1.23E-05                               | no                                | no                       |
| Anthracene            | 2.34E-01         | 3.45E+05                      | Normal/Lognormal               | 1.70E-01                                 | 6.13E-05                               | no                                | no                       |
| Benzo(ghi)perylene    | 1.60E+00         | 1.70E+08                      | Normal/Lognormal               | 1.20E+00                                 | 6.13E+04                               | no                                | no                       |
| Carbazole             | 1.38E-01         | 2.81E+02                      | Normal/Lognormal               | 1.10E-01                                 | 2.86E+02                               | no                                | no                       |
| Di-n-butylphthalate   | 1.46E-01         | 1.55E+00                      | Normal/Lognormal               | 1.10E-01                                 | 2.28E+03                               | no                                | no                       |
| Dibenzofuran          | 1.16E-01         | 5.59E+01                      | Normal/Lognormal               | 9.30E-02                                 | 8.18E-03                               | no                                | no                       |
| Fluoranthene          | 1.13E+00         | 1.59E+04                      | Normal/Lognormal               | 7.80E-01                                 | 8.17E+04                               | no                                | no                       |
| Naphthalene           | 2.04E-01         | 6.64E+05                      | Unknown                        | 1.60E-01                                 | 8.24E+02                               | no                                | no                       |
| Phenanthrene          | 3.29E-01         | 2.65E+07                      | Normal/Lognormal               | 2.50E-01                                 | 6.13E+04                               | no                                | no                       |
| Pyrene                | 1.52E+00         | 7.45E+05                      | Normal/Lognormal               | 1.00E+00                                 | 6.13E+04                               | no                                | no                       |

**Table 7**  
**Statistical Summary and Selection of COPCs in EU4 Surface Soils (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Detection Limit | Maximum mg/kg | Detected | Minimum mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected   | Maximum mg/kg | Location of Maximum Concentration Qualifier | Standard Deviation mg/kg |
|------------------------|------------|---------------|----------------|-----------------|---------------|-----------------|---------------|----------|---------------|------------|------------------------|---------------|------------|---------------|---|--------------------------|
| <b>Semi挥发物</b>         |            |               |                |                 |               |                 |               |          |               |            |                        |               |            |               |   |                          |
| 2,4-Dimethylphenol     | 105-67-9   | 3             | 1              | 33.33           | 4.10E-01      | 9.90E+00        | 2.50E-01      | J        | 1.80E+00      | 6.33E-01   | 2.50E-01               | J             | GEO-19/0-1 | 2.73E+00      |   |                          |
| 2-Methylphthalene      | 91-57-6    | 3             | 3              | 100             | NA            | NA              | 2.70E-01      | 9.36E-01 | 3.42E+00      | 2.80E+02   | 3.42E+00               | J             | GEO-21/0-1 | 1.61E+02      |   |                          |
| 2-Methylphenol         | 95-48-7    | 3             | 1              | 33.33           | 2.00E-01      | 5.00E+00        | 7.30E-02      | J        | 8.91E-01      | 2.62E-01   | 7.30E-02               | J             | GEO-19/0-1 | 1.39E+00      |   |                          |
| 3- and 4-Methylphenol  | 106-44-5   | 3             | 1              | 33.33           | 4.10E-01      | 9.90E+00        | 2.10E-01      | J        | 1.79E-00      | 5.97E-01   | 2.10E-01               | J             | GEO-19/0-1 | 2.74E+00      |   |                          |
| Acenaphthene           | 83-32-9    | 3             | 2              | 66.67           | 2.00E-01      | 2.00E-01        | 1.00E+00      | J        | 6.37E-01      | 2.67E+00   | 1.90E+02               | J             | GEO-21/0-1 | 1.09E+02      |   |                          |
| Acenaphthylene         | 208-96-8   | 3             | 3              | 100             | NA            | NA              | 1.40E+00      | J        | 2.08E-01      | 9.73E+00   | 4.70E+01               | J             | GEO-21/0-1 | 2.35E+01      |   |                          |
| Anthracene             | 120-12-7   | 3             | 3              | 100             | NA            | NA              | 2.10E+00      | J        | 2.62E-02      | 3.42E+01   | 7.60E+02               | J             | GEO-21/0-1 | 4.31E+02      |   |                          |
| Benz(a) Pyrene Equiv.  | -          | 3             | 3              | 100             | NA            | NA              | 5.15E+00      | J        | 1.47E-02      | 5.33E+01   | 3.52E+02               | J             | GEO-21/0-1 | 1.82E+02      |   |                          |
| Benzoglypterylene      | 191-24-2   | 3             | 3              | 100             | NA            | NA              | 2.70E+00      | J        | 3.89E-01      | 1.80E+01   | 9.00E+01               | J             | GEO-21/0-1 | 4.55E+01      |   |                          |
| Carbazole              | 86-74-8    | 3             | 3              | 100             | NA            | NA              | 6.00E-01      | J        | 7.83E+01      | 9.34E+00   | 2.30E+02               | J             | GEO-21/0-1 | 1.31E+02      |   |                          |
| Dibenzofuran           | 132-64-9   | 3             | 3              | 100             | NA            | NA              | 3.40E-01      | J        | 6.37E+01      | 3.65E+00   | 1.90E+02               | J             | GEO-21/0-1 | 1.09E+02      |   |                          |
| Fluoranthene           | 206-44-0   | 3             | 3              | 100             | NA            | NA              | 5.70E+00      | J        | 2.62E+02      | 7.49E+01   | 6.70E+02               | J             | GEO-21/0-1 | 3.57E+02      |   |                          |
| Fluorene               | 86-73-7    | 3             | 2              | 66.67           | 2.00E-01      | 2.00E-01        | 1.40E+00      | J        | 8.72E+01      | 3.31E+00   | 2.60E+02               | J             | GEO-21/0-1 | 1.50E+02      |   |                          |
| Naphthalene            | 91-20-3    | 3             | 3              | 100             | NA            | NA              | 6.80E-01      | J        | 1.64E-02      | 6.35E+00   | 4.90E+02               | J             | GEO-21/0-1 | 2.82E+02      |   |                          |
| N-nitrosodiphenylamine | 86-30-6    | 3             | 1              | 33.33           | 3.70E-02      | 5.00E+00        | 2.00E-01      | J        | 9.06E-01      | 2.10E-01   | 2.00E-01               | J             | GEO-20/0-1 | 1.38E+00      |   |                          |
| Phenanthrene           | 85-01-8    | 3             | 3              | 100             | NA            | NA              | 1.70E+00      | J        | 2.53E-02      | 2.14E+01   | 7.50E+02               | J             | GEO-21/0-1 | 4.30E+02      |   |                          |
| Pyrene                 | 129-00-0   | 3             | 3              | 100             | NA            | NA              | 5.30E+00      | J        | 2.65E+02      | 7.84E+01   | 6.50E+02               | J             | GEO-21/0-1 | 3.40E+02      |   |                          |

NA - Not Available

**Table 7**  
**Statistical Summary and Selection of COPCs in EU4 Surface Soils (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent             | 95% UCL<br>mg/kg | Lognormal<br>95% UCL<br>mg/kg | 99% Confidence   | Distribution | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|-------------------------|------------------|-------------------------------|------------------|--------------|--|--|-----------------------------------|--------------------------|
| <b>Semi-volatiles</b>   |                  |                               |                  |              |  |  |                                   |                          |
| 2,4-Dimethylphenol      | 6.40E+00         | 1.83E+13                      | Normal/Lognormal | 2.50E-01     | 4.08E+04                                 | no                                     | no                                | no                       |
| 2-Methylnaphthalene     | 3.66E+02         | 4.06E+62                      | Lognormal        | 2.80E+02     | 8.18E+04                                 | no                                     | no                                | no                       |
| 2-Methylphenol          | 3.24E+00         | 3.99E+15                      | Normal/Lognormal | 7.30E-02     | 1.02E+05                                 | no                                     | no                                | no                       |
| 3- and 4-Methylphenol   | 6.40E+00         | 9.25E+13                      | Lognormal        | 2.10E-01     | 1.02E+04                                 | no                                     | no                                | no                       |
| Acenaphthene            | 2.48E+02         | 6.33E+63                      | Normal/Lognormal | 1.90E+02     | 1.23E+05                                 | no                                     | no                                | no                       |
| Acenaphthylene          | 6.05E+01         | 2.92E+14                      | Normal/Lognormal | 4.70E+01     | 1.23E+05                                 | no                                     | no                                | no                       |
| Anthracene              | 9.89E+02         | 3.74E+38                      | Normal/Lognormal | 7.60E+01     | 6.13E+05                                 | no                                     | no                                | no                       |
| Benzo (a) Pyrene Equiv. | 4.53E+02         | 1.75E+21                      | Normal/Lognormal | 3.52E+02     | 7.84E-01                                 | YES                                    | no                                | no                       |
| Benzoguaiaculylene      | 1.16E+02         | 3.33E+14                      | Normal/Lognormal | 9.00E+01     | 6.13E+04                                 | no                                     | no                                | no                       |
| Carbazole               | 3.00E+02         | 1.23E+39                      | Normal/Lognormal | 2.30E+02     | 2.86E+02                                 | no                                     | no                                | no                       |
| Dibenzofuran            | 2.48E+02         | 7.02E+50                      | Lognormal        | 1.90E+02     | 8.18E+03                                 | no                                     | no                                | no                       |
| Fluoranthene            | 8.64E+02         | 2.43E+26                      | Normal/Lognormal | 6.70E+02     | 8.17E+04                                 | no                                     | no                                | no                       |
| Fluorene                | 3.40E+02         | 2.06E+68                      | Normal/Lognormal | 2.60E+02     | 8.17E+04                                 | no                                     | no                                | no                       |
| Naphthalene             | 6.40E+02         | 5.82E+60                      | Lognormal        | 4.90E+02     | 8.24E+02                                 | no                                     | no                                | no                       |
| N-nitrosodiphenylamine  | 3.24E+00         | 6.32E+24                      | Normal/Lognormal | 2.00E-01     | 1.17E+03                                 | no                                     | no                                | no                       |
| Phenanthrene            | 9.79E+02         | 7.80E+43                      | Normal/Lognormal | 7.50E+02     | 6.13E+04                                 | no                                     | no                                | no                       |
| Pyrene                  | 8.38E+02         | 2.72E+27                      | Normal/Lognormal | 6.50E+02     | 6.13E+04                                 | no                                     | no                                | no                       |

Table 1

**Statistical Summary and Selection of COPCs in EU5 Surface Soil (0-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent             | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected | Minimum mg/kg | Detected | Qualifier | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Detected | Maximum mg/kg | Detected | Qualifier | Mean mg/kg | Logarithmic Mean mg/kg | Maximum mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|-------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------|---------------|----------|-----------|------------|------------------------|---------------|----------|---------------|----------|-----------|------------|------------------------|---------------|-----------------------------------|--------------------------|
| <b>SemiVolatiles</b>    |            |               |                |                 |               |               |                       |          |               |          |           |            |                        |               |          |               |          |           |            |                        |               |                                   |                          |
| 2,4-Dimethylphenol      | 105-67-9   | 9             | 2              | 22.22           | 7.60E-02      | 9.90E+00      | 1.10E-01              | J        | 6.72E-01      | 1.60E-01 | 2.50E-01  | J          | GEO-19/0-1             | 1.61E+00      |          |               |          |           |            |                        |               |                                   |                          |
| 2-Methylnaphthalene     | 91-57-6    | 9             | 7              | 77.78           | 3.80E-02      | 3.90E-02      | 5.10E-02              | J        | 3.23E+01      | 4.65E-01 | 2.80E+02  | J          | GEO-21/0-1             | 9.29E+01      |          |               |          |           |            |                        |               |                                   |                          |
| 2-Methylphenol          | 95-48-7    | 9             | 2              | 22.22           | 3.80E-02      | 5.00E+00      | 4.20E-02              | J        | 3.30E-01      | 7.20E-02 | 7.30E-02  | J          | GEO-19/0-1             | 8.15E-01      |          |               |          |           |            |                        |               |                                   |                          |
| 3- and 4-Methylphenol   | 106-44-5   | 9             | 2              | 22.22           | 7.60E-02      | 9.90E+00      | 1.40E-01              | J        | 6.68E-01      | 1.59E-01 | 2.10E-01  | J          | GEO-19/0-1             | 1.61E+00      |          |               |          |           |            |                        |               |                                   |                          |
| Benzo (a) Pyrene Equiv. | -          | 9             | 8              | 88.89           | NA            | NA            | 5.09E-01              | 5.99E+01 | 6.47E-00      | 3.52E+02 |           |            | GEO-21/0-1             | 1.15E+02      |          |               |          |           |            |                        |               |                                   |                          |
| N-nitrosodiphenylamine  | 86-30-6    | 9             | 1              | 11.11           | 3.70E-02      | 5.00E+00      | 2.00E-01              | 3.33E-01 | 6.13E-02      | 2.00E-01 |           |            | GEO-20/0-1             | 8.15E-01      |          |               |          |           |            |                        |               |                                   |                          |
| Acenaphthene            | 83-32-9    | 9             | 4              | 44.44           | 3.80E-02      | 2.00E-01      | 1.60E-01              | J        | 2.72E-01      | 3.62E-01 | 1.90E+02  |            | GEO-21/0-1             | 6.35E-01      |          |               |          |           |            |                        |               |                                   |                          |
| Acenaphthylene          | 208-96-8   | 9             | 8              | 88.89           | 3.80E-02      | 3.80E-02      | 4.80E-02              | J        | 9.09E+00      | 1.13E+00 | 4.70E+01  | J          | GEO-21/0-1             | 1.55E-01      |          |               |          |           |            |                        |               |                                   |                          |
| Anthracene              | 120-12-7   | 9             | 7              | 77.78           | 3.80E-02      | 3.90E-02      | 1.30E-01              | J        | 9.70E+01      | 2.14E+00 | 7.60E+02  | J          | GEO-21/0-1             | 2.50E+02      |          |               |          |           |            |                        |               |                                   |                          |
| Benzo(ghi)perylene      | 191-24-2   | 9             | 7              | 77.78           | 3.80E-02      | 3.80E-02      | 2.30E-01              | J        | 1.64E+01      | 1.53E+00 | 9.00E+01  | J          | GEO-21/0-1             | 2.95E+01      |          |               |          |           |            |                        |               |                                   |                          |
| Carbazole               | 86-74-8    | 9             | 6              | 66.67           | 3.80E-02      | 3.90E-02      | 5.30E-01              | J        | 2.80E+01      | 7.44E+01 | 2.30E+02  | J          | GEO-21/0-1             | 7.59E+01      |          |               |          |           |            |                        |               |                                   |                          |
| Dibenzofuran            | 132-64-9   | 9             | 7              | 77.78           | 3.80E-02      | 3.90E-02      | 3.90E-02              | J        | 2.45E+01      | 4.85E+01 | 1.90E+02  |            | GEO-21/0-1             | 6.28E+01      |          |               |          |           |            |                        |               |                                   |                          |
| Fluoranthene            | 206-44-0   | 9             | 8              | 88.89           | 3.80E-02      | 3.80E-02      | 1.30E-01              | J        | 1.30E+02      | 7.24E+00 | 6.70E+02  |            | GEO-21/0-1             | 2.33E+02      |          |               |          |           |            |                        |               |                                   |                          |
| Fluorene                | 86-73-7    | 9             | 5              | 55.56           | 3.80E-02      | 2.00E-01      | 3.30E-01              | J        | 3.62E+01      | 5.14E+01 | 2.60E+02  |            | GEO-21/0-1             | 8.65E+01      |          |               |          |           |            |                        |               |                                   |                          |
| Naphthalene             | 91-20-3    | 9             | 7              | 77.78           | 3.80E-02      | 3.90E-02      | 7.50E-02              | J        | 5.55E+01      | 6.10E+01 | 4.90E+02  |            | GEO-21/0-1             | 1.63E+02      |          |               |          |           |            |                        |               |                                   |                          |
| Phenanthrene            | 85-01-8    | 9             | 7              | 77.78           | 3.80E-02      | 3.90E-02      | 1.80E-01              | J        | 1.12E+02      | 2.07E+00 | 7.50E+02  |            | GEO-21/0-1             | 2.52E+02      |          |               |          |           |            |                        |               |                                   |                          |
| Phenol                  | 108-95-2   | 9             | 4              | 44.44           | 7.40E-02      | 9.90E+00      | 1.40E-01              | J        | 7.10E-01      | 2.08E-01 | 3.80E+01  | J          | GEO-29/0-1             | 1.59E+00      |          |               |          |           |            |                        |               |                                   |                          |
| Pyrene                  | 129-00-0   | 9             | 8              | 88.89           | 3.80E-02      | 3.80E-02      | 2.50E-01              | J        | 1.20E+02      | 7.66E+00 | 6.50E+02  | J          | GEO-21/0-1             | 2.18E+02      |          |               |          |           |            |                        |               |                                   |                          |

**Table 8**  
**Statistical Summary and Selection of COPCs in EU5 Surface Soil (0'-1' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent             | 95% UCL<br>mg/kg | Lognormal<br>95% UCL<br>mg/kg | Distribution<br>95% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|-------------------------|------------------|-------------------------------|--------------------------------|--|--|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>    |                  |                               |                                |  |  |                                   |                          |
| 2,4-Dimethylphenol      | 1.67E+00         | 5.38E+00                      | Lognormal                      | 2.50E-01                                 | 4.08E-04                               | no                                | no                       |
| 2-Methylnaphthalene     | 8.99E-01         | 5.63E+05                      | Lognormal                      | 2.80E-02                                 | 8.18E+04                               | no                                | no                       |
| 2-Methylphenol          | 8.35E-01         | 2.52E+00                      | Lognormal                      | 7.30E-02                                 | 1.02E+05                               | no                                | no                       |
| 3- and 4-Methylphenol   | 1.66E-00         | 5.25E+00                      | Lognormal                      | 2.10E-01                                 | 1.02E+04                               | no                                | no                       |
| Benzo (a) Pyrene Equiv. | 1.31E-02         | 7.89E+05                      | Lognormal                      | 3.52E-02                                 | 7.84E-01                               | YES                               | YES . COPC               |
| N-nitrosodiphenylamine  | 8.38E-01         | 4.44E+00                      | Unknown                        | 2.00E-01                                 | 1.17E+03                               | no                                | no                       |
| Acenaphthene            | 6.66E+01         | 1.62E+07                      | Lognormal                      | 1.90E-02                                 | 1.23E+05                               | no                                | no                       |
| Acenaphthylene          | 1.87E+01         | 7.53E+04                      | Lognormal                      | 4.70E-01                                 | 1.23E+05                               | no                                | no                       |
| Anthracene              | 2.52E+02         | 7.31E+08                      | Lognormal                      | 7.60E-02                                 | 6.13E+05                               | no                                | no                       |
| Benzofuran              | 3.46E+01         | 1.83E+06                      | Lognormal                      | 9.00E-01                                 | 6.13E+04                               | no                                | no                       |
| Carbazole               | 7.51E+01         | 6.76E+06                      | Lognormal                      | 2.30E-02                                 | 2.86E+02                               | no                                | no                       |
| Dibenzofuran            | 6.34E+01         | 1.50E+06                      | Lognormal                      | 1.90E-02                                 | 8.18E+03                               | no                                | no                       |
| Fluoranthene            | 2.75E+02         | 7.48E+08                      | Lognormal                      | 6.70E-02                                 | 8.17E+04                               | no                                | no                       |
| Fluorene                | 8.98E-01         | 3.94E+07                      | Lognormal                      | 2.60E-02                                 | 8.17E+04                               | no                                | no                       |
| Naphthalene             | 1.57E+02         | 1.46E+06                      | Lognormal                      | 4.90E-02                                 | 8.24E+02                               | no                                | no                       |
| Phenanthrene            | 2.69E+02         | 1.17E+09                      | Lognormal                      | 7.50E-02                                 | 6.13E+04                               | no                                | no                       |
| Phenol                  | 1.70E+00         | 5.03E+00                      | Lognormal                      | 3.80E-01                                 | 1.23E+05                               | no                                | no                       |
| Pyrene                  | 2.55E+02         | 1.89E+08                      | Lognormal                      | 6.50E-02                                 | 6.13E+04                               | no                                | no                       |

**Table 9**  
**Statistical Summary and Selection of COPCs in EU2 Soils (0-16' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum mg/kg | Maximum mg/kg | Detection Limit mg/kg | Detected mg/kg | Minimum Detected mg/kg | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Detected Qualifier | Maximum Qualifier mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|----------------------------|------------|---------------|----------------|-----------------|---------------|---------------|-----------------------|----------------|------------------------|------------|------------------------|------------------------|--------------------|-------------------------|-----------------------------------|--------------------------|
| Pesticides                 |            |               |                |                 |               |               |                       |                |                        |            |                        |                        |                    |                         |                                   |                          |
| Endosulfan I               | 959-98-8   | 1             | 1              | 100             | NA            | NA            | 4.00E-03              | J              | 4.00E-03               | 4.00E-03   | 4.00E-03               | J                      | SB-05/4-9          | 0.00E+00                |                                   |                          |
| Heptachlor                 | 76-44-8    | 1             | 1              | 100             | NA            | NA            | 1.00E-02              | 1.00E-02       | 1.00E-02               | 1.00E-02   | 1.00E-02               | J                      | SB-05/4-9          | 0.00E+00                |                                   |                          |
| Semi-volatiles             |            |               |                |                 |               |               |                       |                |                        |            |                        |                        |                    |                         |                                   |                          |
| 2,4-Dimethylphenol         | 105-67-9   | 27            | 2              | 7.41            | 6.70E-02      | 3.30E-01      | 1.10E+00              | J              | 1.26E-01               | 4.77E-02   | 1.30E+00               | J                      | SB-06/10-12        | 3.12E-01                |                                   |                          |
| 2-Methylnaphthalene        | 91-57-6    | 27            | 5              | 18.52           | 3.30E-02      | 3.90E-02      | 7.00E-02              | J              | 2.30E-01               | 5.57E-02   | 2.30E+02               | J                      | SB-07/5-7          | 6.64E+01                |                                   |                          |
| Benzo (a) Pyrene Equiv.    | -          | 27            | 18             | 66.67           | NA            | NA            | 8.41E-02              | NA             | 3.99E+00               | 3.77E-01   | 3.58E+01               | NA                     | SB-07/5-7          | 9.54E+00                |                                   |                          |
| Acenaphthene               | 83-32-9    | 27            | 4              | 14.81           | 3.30E-02      | 3.90E-02      | 4.90E-02              | J              | 1.66E-01               | 4.85E-02   | 2.00E+02               | J                      | SB-07/5-7          | 4.98E+01                |                                   |                          |
| Acenaphthylene             | 208-96-8   | 27            | 9              | 33.33           | 3.30E-02      | 3.90E-02      | 3.70E-02              | J              | 8.23E-01               | 5.42E-02   | 7.70E+00               | J                      | SB-07/5-7          | 2.13E+00                |                                   |                          |
| Anthracene                 | 120-12-7   | 27            | 10             | 37.04           | 3.30E-02      | 3.90E-02      | 4.10E-02              | J              | 9.85E+00               | 7.80E-02   | 1.20E+02               | J                      | SB-07/5-7          | 2.95E+01                |                                   |                          |
| Benzoguaiaculene           | 191-24-2   | 27            | 13             | 48.15           | 3.70E-02      | 6.70E-02      | 1.70E-01              | J              | 8.11E-01               | 1.35E-01   | 6.40E+00               | J                      | SB-07/5-7          | 1.65E+00                |                                   |                          |
| Bis(2-ethylhexyl)phthalate | 117-81-7   | 27            | 1              | 3.7             | 6.70E-02      | 5.00E-01      | 3.70E-01              | J              | 6.46E-02               | 4.55E-02   | 3.70E-01               | J                      | GEO-13/5-6         | 8.05E-02                |                                   |                          |
| Carbazole                  | 86-74-8    | 27            | 7              | 25.93           | 3.30E-02      | 3.90E-02      | 4.30E-02              | J              | 4.30E+00               | 5.41E-02   | 5.00E+01               | J                      | SB-05/4-9          | 1.27E-01                |                                   |                          |
| Di-n-butylphthalate        | 84-74-2    | 27            | 9              | 33.33           | 3.30E-02      | 2.50E-01      | 3.60E-02              | J              | 4.41E-02               | 3.66E-02   | 1.10E-01               | J                      | SB-10              | 2.86E-02                |                                   |                          |
| Dibenzofuran               | 132-64-9   | 27            | 5              | 18.52           | 3.30E-02      | 3.90E-02      | 7.20E-02              | J              | 1.61E+01               | 5.25E-02   | 1.80E+02               | J                      | SB-07/5-7          | 4.72E-01                |                                   |                          |
| Fluoranthene               | 206-44-0   | 27            | 16             | 59.26           | 3.30E-02      | 3.80E-02      | 5.00E-02              | J              | 2.30E+01               | 2.25E-01   | 2.50E+02               | J                      | SB-07/5-7          | 6.50E-01                |                                   |                          |
| Fluorene                   | 86-73-7    | 27            | 6              | 22.22           | 3.30E-02      | 3.80E-02      | 4.50E-02              | J              | 2.02E+01               | 6.16E-02   | 2.50E+02               | J                      | SB-07/5-7          | 6.05E+01                |                                   |                          |
| Naphthalene                | 91-20-3    | 27            | 7              | 25.93           | 3.30E-02      | 3.90E-02      | 8.80E-02              | J              | 3.78E-01               | 6.94E-02   | 4.20E+02               | J                      | SB-06/10-12        | 1.13E+02                |                                   |                          |
| Phenanthrene               | 85-01-8    | 27            | 11             | 40.74           | 3.30E-02      | 3.90E-02      | 3.70E-02              | J              | 4.58E-01               | 9.58E-02   | 5.10E+02               | J                      | SB-07/5-7          | 1.34E+02                |                                   |                          |
| Phenol                     | 108-95-2   | 27            | 2              | 7.41            | 3.30E-02      | 2.50E-01      | 1.10E-01              | J              | 3.94E-02               | 2.75E-02   | 1.90E-01               | J                      | GEO-3/5-6          | 4.30E-02                |                                   |                          |
| Pyrene                     | 129-00-0   | 27            | 16             | 59.26           | 3.70E-02      | 6.70E-02      | 6.80E-02              | J              | 1.98E+01               | 3.09E-01   | 2.30E+02               | J                      | SB-07/5-7          | 5.63E+01                |                                   |                          |
| Volatiles                  |            |               |                |                 |               |               |                       |                |                        |            |                        |                        |                    |                         |                                   |                          |
| Acetone                    | 67-64-1    | 7             | 2              | 28.57           | 7.00E-03      | 3.50E-02      | 4.40E-02              | J              | 1.98E-02               | 9.56E-03   | 6.30E-02               | J                      | SB-05/4-9          | 2.42E-02                |                                   |                          |
| Benzene                    | 71-43-2    | 7             | 1              | 14.29           | 1.00E-03      | 5.00E-03      | 8.00E-03              | J              | 2.14E-03               | 1.18E-03   | 8.00E-03               | J                      | SB-06/10-12        | 2.75E-03                |                                   |                          |
| Ethylbenzene               | 100-41-4   | 7             | 3              | 42.86           | 1.00E-03      | 1.00E-03      | 6.80E-02              | J              | 1.10E-01               | 6.42E-03   | 4.80E-01               | J                      | SB-06/10-12        | 1.82E-01                |                                   |                          |
| Toluene                    | 108-88-3   | 7             | 3              | 42.86           | 1.00E-03      | 1.00E-03      | 1.40E-02              | J              | 2.26E-02               | 3.26E-03   | 9.30E-02               | J                      | SB-06/10-12        | 3.63E-02                |                                   |                          |
| Xylene (total)             | 1330-20-7  | 7             | 3              | 42.86           | 1.00E-03      | 4.90E-01      | 4.85E-01              | 1.30E-02       | 1.70E+00               | 4.85E-01   | 1.70E+00               | J                      | SB-06/10-12        | 6.98E-01                |                                   |                          |

NA - Not Available

**Table 9**  
**Statistical Summary and Selection of COPCs in EU2 Soils (0-16' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | 95% UCL<br>mg/kg | Lognormal 95%<br>UCL<br>mg/kg | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier 1 Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|----------------------------|------------------|-------------------------------|--------------------------------|--|--|-----------------------------------|--------------------------|
| <b>Pesticides</b>          |                  |                               |                                |  |  |                                   |                          |
| Endosulfan 1               | NA               | NA                            | Unknown                        | 4.00E-03                                 | 1.23E-03                               | no                                | NA                       |
| Heptachlor                 | NA               | NA                            | Unknown                        | 1.09E-02                                 | 1.95E-01                               | no                                | NA                       |
| <b>Semi-volatiles</b>      |                  |                               |                                |  |  |                                   |                          |
| 2,4-Dimethylphenol         | 2.28E-01         | 1.23E-01                      | Unknown                        | 1.23E-01                                 | 4.38E-04                               | no                                | no                       |
| 2-Methylnaphthalene        | 4.48E+01         | 1.33E+02                      | Unknown                        | 1.33E+02                                 | 8.18E-04                               | no                                | no                       |
| Benz (a) Pyrene Equiv.     | 7.13E+00         | 1.91E+01                      | Unknown                        | 1.91E+01                                 | 7.84E-01                               | YES                               | YES - COPC               |
| Acenaphthene               | 3.29E+01         | 6.50E+01                      | Unknown                        | 6.50E+01                                 | 1.23E+05                               | no                                | no                       |
| Acenaphthylene             | 1.52E+00         | 2.17E+00                      | Unknown                        | 2.17E+00                                 | 1.23E+05                               | no                                | no                       |
| Anthracene                 | 1.95E+01         | 5.78E+01                      | Unknown                        | 5.78E+01                                 | 6.13E+05                               | no                                | no                       |
| Benz(ghi)perylene          | 1.35E+00         | 3.92E+00                      | Unknown                        | 3.92E+00                                 | 6.13E+04                               | no                                | no                       |
| Bis(2-ethylhexyl)phthalate | 9.11E-02         | 7.71E-02                      | Unknown                        | 7.71E-02                                 | 4.09E-02                               | no                                | no                       |
| Carbazole                  | 8.46E+00         | 1.17E+01                      | Unknown                        | 1.17E+01                                 | 2.86E+02                               | no                                | no                       |
| Di-n-butylphthalate        | 5.35E-02         | 5.71E-02                      | Unknown                        | 5.71E-02                                 | 2.28E-03                               | no                                | no                       |
| Dibenzofuran               | 3.16E+01         | 6.88E+01                      | Unknown                        | 6.88E+01                                 | 8.18E-03                               | no                                | no                       |
| Fluoranthene               | 4.43E+01         | 7.45E+02                      | Unknown                        | 2.50E+02                                 | 8.17E+04                               | no                                | no                       |
| Fluorene                   | 4.01E+01         | 1.31E+02                      | Unknown                        | 1.31E+02                                 | 8.17E+04                               | no                                | no                       |
| Naphthalene                | 7.50E+01         | 3.31E+02                      | Unknown                        | 3.31E+02                                 | 8.24E+02                               | no                                | no                       |
| Phenanthrene               | 8.98E+01         | 6.00E+02                      | Unknown                        | 5.10E+02                                 | 6.13E+04                               | no                                | no                       |
| Phenol                     | 5.35E-02         | 5.15E-02                      | Unknown                        | 5.15E-02                                 | 1.23E+05                               | no                                | no                       |
| Pyrene                     | 3.83E+01         | 4.41E+02                      | Unknown                        | 2.30E+02                                 | 6.13E+04                               | no                                | no                       |
| <b>Volatiles</b>           |                  |                               |                                |  |  |                                   |                          |
| Acetone                    | 3.76E-02         | 2.64E-01                      | Normal/Lognormal               | 6.30E-02                                 | 1.04E+05                               | no                                | no                       |
| Benzene                    | 4.16E-03         | 1.49E-02                      | Lognormal                      | 8.00E-03                                 | 1.36E+00                               | no                                | no                       |
| Ethylbenzene               | 2.44E-01         | 1.63E+05                      | Lognormal                      | 4.80E-01                                 | 3.95E-02                               | no                                | no                       |
| Toluene                    | 4.93E-02         | 1.58E+02                      | Lognormal                      | 9.50E-12                                 | 3.80E-01                               | no                                | no                       |
| Xylene (total)             | 9.97E-01         | 2.80E+11                      | Normal                         | 1.70E+00                                 | 3.18E+02                               | no                                | no                       |

NA = Not Available

Table

Statistical Summary and Selection of COPCs in EU3 Soils (0-16' bgs)  
 Kerr McGee, Hattiesburg, MS

| Constituent             | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum Detection mg/kg | Maximum Detection mg/kg | Minimum Limit mg/kg | Maximum Limit mg/kg | Detected mg/kg | Detected Qualifier | Mean mg/kg | Logarithmic Mean mg/kg | Maximum Detected mg/kg | Maximum Qualifier mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|-------------------------|------------|---------------|----------------|-----------------|-------------------------|-------------------------|---------------------|---------------------|----------------|--------------------|------------|------------------------|------------------------|-------------------------|-----------------------------------|--------------------------|
| <b>Semivolatiles</b>    |            |               |                |                 |                         |                         |                     |                     |                |                    |            |                        |                        |                         |                                   |                          |
| 2-Methylnaphthalene     | 91-57-6    | 7             | 1              | 14.29           | 3.30E-02                | 4.00E-02                | 2.30E-01            | J                   | 4.85E-02       | 2.62E-02           | 2.30E-01   | J                      | SS-16                  | 8.00E-02                |                                   |                          |
| Acenaphthylene          | 208-96-8   | 7             | 2              | 28.57           | 3.30E-02                | 4.00E-02                | 1.20E-01            | J                   | 5.47E-02       | 3.33E-02           | 1.70E-01   | J                      | SS-16                  | 6.34E-02                |                                   |                          |
| Anthracene              | 120-12-7   | 7             | 2              | 28.57           | 3.30E-02                | 4.00E-02                | 1.20E-01            | J                   | 5.47E-02       | 3.33E-02           | 1.70E-01   | J                      | SS-16                  | 6.34E-02                |                                   |                          |
| Benzo (a) Pyrene Equiv. | -          | 7             | 3              | 42.86           | NA                      | NA                      | 1.01E-01            |                     | 3.33E-01       | 1.22E-01           | 1.12E-00   |                        | SS-16                  | 4.77E-01                |                                   |                          |
| Benzofluoranthene       | 191-24-2   | 7             | 3              | 42.86           | 3.70E-02                | 4.00E-02                | 8.00E-02            | J                   | 2.91E-01       | 7.06E-02           | 1.20E-00   |                        | SS-16                  | 4.69E-01                |                                   |                          |
| Carbazole               | 86-74-8    | 7             | 2              | 28.57           | 3.30E-02                | 4.00E-02                | 4.60E-02            | J                   | 3.56E-02       | 2.72E-02           | 1.10E-01   | J                      | SS-16                  | 3.44E-02                |                                   |                          |
| Di-n-butylphthalate     | 84-74-2    | 7             | 3              | 42.86           | 7.50E-02                | 7.90E-02                | 4.00E-02            | J                   | 5.74E-02       | 5.13E-02           | 1.10E-01   | J                      | SS-16                  | 3.23E-02                |                                   |                          |
| Dibenzofuran            | 132-64-9   | 7             | 2              | 28.57           | 3.30E-02                | 4.00E-02                | 3.60E-02            | J                   | 3.17E-02       | 2.57E-02           | 9.30E-02   | J                      | SS-16                  | 2.78E-02                |                                   |                          |
| Fluoranthene            | 206-44-0   | 7             | 3              | 42.86           | 3.70E-02                | 4.00E-02                | 1.20E-01            | J                   | 2.37E-01       | 7.03E-02           | 7.80E-01   |                        | SS-16                  | 3.40E-01                |                                   |                          |
| Naphthalene             | 91-20-3    | 7             | 1              | 14.29           | 3.30E-02                | 4.00E-02                | 1.60E-01            | J                   | 3.85E-02       | 2.48E-02           | 1.60E-01   | J                      | SS-16                  | 5.36E-02                |                                   |                          |
| Phenanthrene            | 85-01-8    | 7             | 2              | 28.57           | 3.30E-02                | 4.00E-02                | 1.30E-01            | J                   | 6.76E-02       | 3.55E-02           | 2.50E-01   | J                      | SS-16                  | 9.05E-02                |                                   |                          |
| Phenol                  | 108-95-2   | 7             | 2              | 28.57           | 3.30E-02                | 7.90E-02                | 9.60E-02            | J                   | 6.47E-02       | 3.95E-02           | 2.30E-01   | J                      | GEO-175-6              | 7.81E-02                |                                   |                          |
| Pyrene                  | 129-00-0   | 7             | 3              | 42.86           | 3.70E-02                | 4.00E-02                | 1.20E-01            | J                   | 3.07E-01       | 7.64E-02           | 1.00E+00   |                        | SS-17                  | 4.58E-01                |                                   |                          |

**Table 10**  
**Statistical Summary and Selection of COPCs in EU3 Soils (0-16' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent            | 95% UCL<br>mg/kg | Lognormal<br>95% UCL<br>mg/kg | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|------------------------|------------------|-------------------------------|--------------------------------|--|--|-----------------------------------|--------------------------|
| Semivolatiles          |                  |                               |                                |  |  |                                   |                          |
| 2-Methylnaphthalene    | 1.07E-01         | 1.70E-01                      | Unknown                        | 1.70E-01                                 | 8.18E+04                               | no                                | no                       |
| Acenaphthylene         | 1.01E-01         | 2.50E-01                      | Unknown                        | 1.70E-01                                 | 1.23E+05                               | no                                | no                       |
| Antracene              | 1.01E-01         | 2.50E-01                      | Unknown                        | 1.70E-01                                 | 6.13E+05                               | no                                | no                       |
| Benz (a) Pyrene Equiv. | 6.83E-01         | 8.22E+00                      | Unknown                        | 1.12E+00                                 | 7.84E-01                               | YES                               | no                       |
| Benz(ghi)perylene      | 6.33E-01         | 3.79E+01                      | Lognormal                      | 1.20E+00                                 | 6.13E+04                               | no                                | no                       |
| Carbazole              | 6.08E-02         | 8.08E-02                      | Unknown                        | 8.08E-02                                 | 2.86E+02                               | no                                | no                       |
| Di-n-butylphthalate    | 8.12E-02         | 9.38E-02                      | Unknown                        | 9.38E-02                                 | 2.28E+03                               | no                                | no                       |
| Dibenzofuran           | 5.21E-02         | 6.21E-02                      | Unknown                        | 6.21E-02                                 | 8.18E+03                               | no                                | no                       |
| Fluoranthene           | 4.87E-01         | 2.06E+01                      | Lognormal                      | 7.80E-01                                 | 8.17E+04                               | no                                | no                       |
| Naphthalene            | 7.79E-02         | 1.03E-01                      | Unknown                        | 1.03E-01                                 | 8.24E+02                               | no                                | no                       |
| Phenanthrene           | 1.34E-01         | 4.34E-01                      | Unknown                        | 2.50E-01                                 | 6.13E+04                               | no                                | no                       |
| Phenol                 | 1.22E-01         | 3.11E-01                      | Lognormal                      | 2.30E-01                                 | 1.23E+05                               | no                                | no                       |
| Pyrene                 | 6.43E-01         | 5.32E+01                      | Lognormal                      | 1.00E+00                                 | 6.13E+04                               | no                                | no                       |

**Table 11**  
**Statistical Summary and Selection of COPCs in EU5 Soil (θ-16' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                  | CAS Number | Total Samples | Number of Hits | Hit Frequency % | Minimum Detection Limit mg/kg | Maximum Detection Limit mg/kg | Minimum Detected mg/kg | Maximum Detected mg/kg | Logarithmic Mean mg/kg | Mean mg/kg | Maximum Detected mg/kg | Detected Qualifier | Maximum Concentration mg/kg | Location of Maximum Concentration | Standard Deviation mg/kg |
|------------------------------|------------|---------------|----------------|-----------------|-------------------------------|-------------------------------|------------------------|------------------------|------------------------|------------|------------------------|--------------------|-----------------------------|-----------------------------------|--------------------------|
| <b>Semi-volatiles</b>        |            |               |                |                 |                               |                               |                        |                        |                        |            |                        |                    |                             |                                   |                          |
| 2,4-Dimethylphenol           | 105-67-9   | 34            | 5              | 14.71           | 6.70E-02                      | 9.90E+00                      | 7.90E-02               | J                      | 4.99E-01               | 8.34E-02   | 8.90E+00               | J                  | GEO-21/2-3                  | 1.71E+00                          |                          |
| 2-Methylnaphthalene          | 91-57-6    | 34            | 18             | 52.94           | 3.30E-02                      | 4.10E-02                      | 5.10E-02               | J                      | 7.38E+01               | 3.10E-01   | 1.50E-03               | J                  | GEO-21/2-3                  | 2.68E+02                          |                          |
| 2-Methylphenol               | 95-48-7    | 34            | 3              | 8.82            | 3.70E-02                      | 5.00E-00                      | 4.20E-02               | J                      | 4.15E-02               | 7.30E-02   | J                      | J                  | GEO-19/0-1                  | 4.77E-01                          |                          |
| 3- and 4-Methylphenol        | 106-44-5   | 34            | 2              | 5.88            | 7.30E-02                      | 9.90E-00                      | 1.40E-01               | J                      | 3.19E-01               | 7.95E-02   | 2.10E-01               | J                  | GEO-19/0-1                  | 9.35E-01                          |                          |
| Benzo (a) Pyrene Equiv.      | -          | 34            | 22             | 64.71           | NA                            | NA                            | 8.74E-02               | 2.68E+01               | 7.91E-01               | 3.52E+02   | 7.40E+01               | J                  | GEO-21/0-1                  | 7.74E+01                          |                          |
| N-nitrosodiphenylamine       | 86-39-6    | 34            | 1              | 2.94            | 3.70E-02                      | 5.00E-00                      | 2.00E-01               | 1.66E-01               | 3.87E-02               | 2.00E-01   | 4.78E-01               | J                  | GEO-20/0-1                  | 4.78E-01                          |                          |
| Acenaphthene                 | 83-32-9    | 34            | 16             | 47.06           | 3.30E-02                      | 2.00E-01                      | 9.70E-02               | J                      | 5.68E+01               | 3.20E-01   | 1.20E-03               | J                  | GEO-21/2-3                  | 2.11E+02                          |                          |
| Acenaphthylene               | 208-96-8   | 34            | 17             | 50              | 3.30E-02                      | 4.10E-02                      | 4.80E-02               | J                      | 4.42E+00               | 1.81E-01   | 5.00E+01               | J                  | GEO-21/2-3                  | 1.18E+01                          |                          |
| Anthracene                   | 120-12-7   | 34            | 19             | 55.88           | 3.30E-02                      | 4.10E-02                      | 1.30E-01               | J                      | 8.44E+01               | 5.25E-01   | 1.80E+03               | J                  | GEO-21/2-3                  | 3.30E+02                          |                          |
| Benzo(ghi)perylene           | 191-24-2   | 34            | 18             | 52.94           | 3.80E-02                      | 6.70E-02                      | 4.20E-02               | J                      | 6.88E+00               | 2.16E-01   | 9.00E+01               | J                  | GEO-21/0-1                  | 1.97E+01                          |                          |
| Bis[2-(ethylhexyl)]phthalate | 117-81-7   | 34            | 2              | 5.88            | 6.70E-02                      | 9.90E+00                      | 1.40E-01               | J                      | 3.04E-01               | 7.48E-02   | 1.50E-01               | J                  | GEO-32/2-3                  | 9.30E-01                          |                          |
| Carbazole                    | 86-74-8    | 34            | 17             | 50              | 3.30E-02                      | 4.10E-02                      | 2.10E-01               | J                      | 2.91E+01               | 2.76E-01   | 6.20E+02               | J                  | GEO-21/2-3                  | 1.12E+02                          |                          |
| Dibenzofuran                 | 132-64-9   | 34            | 19             | 55.88           | 3.30E-02                      | 4.10E-02                      | 3.90E-02               | J                      | 5.21E+01               | 3.37E-01   | 1.10E-03               | J                  | GEO-21/2-3                  | 1.94E+02                          |                          |
| Fluoranthene                 | 206-44-0   | 34            | 22             | 64.71           | 3.30E-02                      | 4.10E-02                      | 9.80E-02               | J                      | 1.16E+02               | 1.16E+02   | 2.00E+03               | J                  | GEO-21/2-3                  | 3.63E+02                          |                          |
| Fluorene                     | 86-73-7    | 34            | 17             | 50              | 3.30E-02                      | 2.00E-01                      | 1.40E-01               | J                      | 7.91E+01               | 3.86E-01   | 1.50E+03               | J                  | GEO-21/2-3                  | 2.63E+02                          |                          |
| Naphthalene                  | 91-20-3    | 34            | 18             | 52.94           | 3.30E-02                      | 4.10E-02                      | 7.50E-02               | J                      | 1.62E+02               | 3.97E-01   | 3.50E+03               | J                  | GEO-21/2-3                  | 6.16E+02                          |                          |
| Phenanthrene                 | 85-01-8    | 34            | 20             | 58.82           | 3.30E-02                      | 4.10E-02                      | 8.50E-02               | J                      | 1.87E+02               | 8.79E-01   | 4.00E+03               | J                  | GEO-21/2-3                  | 6.98E+02                          |                          |
| Phenol                       | 108-95-2   | 34            | 14             | 41.18           | 3.30E-02                      | 9.90E+00                      | 1.00E-01               | J                      | 3.34E-01               | 1.06E-01   | 3.80E-01               | J                  | GEO-29/0-1                  | 9.20E-01                          |                          |
| Pyrene                       | 129-00-0   | 34            | 22             | 64.71           | 3.80E-02                      | 6.70E-02                      | 9.10E-02               | J                      | 8.35E+01               | 1.04E-00   | 1.30E-03               | J                  | GEO-21/2-3                  | 2.48E+02                          |                          |
| <b>Volatiles</b>             |            |               |                |                 |                               |                               |                        |                        |                        |            |                        |                    |                             |                                   |                          |
| Acetone                      | 67-64-1    | 5             | 5              | 100             | NA                            | NA                            | 9.00E-03               | J                      | 4.40E-02               | 2.95E-02   | 1.00E-01               | J                  | SB-05/10.5-12.5             | 3.79E-02                          |                          |
| Benzene                      | 71-43-2    | 5             | 2              | 40              | 1.00E-03                      | 1.00E-03                      | 5.00E-03               | J                      | 2.70E-03               | 1.34E-03   | 7.00E-03               | J                  | SB-05/10.5-12.5             | 3.09E-03                          |                          |
| Ethylbenzene                 | 100-41-4   | 5             | 3              | 60              | 1.00E-03                      | 1.00E-03                      | 2.40E-02               | 3.82E-02               | 8.02E-03               | 1.20E-01   | SB-05/10.5-12.5        | 4.95E-02           |                             |                                   |                          |
| Styrene                      | 100-42-5   | 5             | 1              | 20              | 1.00E-03                      | 1.00E-03                      | 1.00E-01               | 2.04E-02               | 1.44E-03               | 1.00E-01   | SB-05/10.5-12.5        | 4.45E-02           |                             |                                   |                          |
| Toluene                      | 108-88-3   | 5             | 3              | 60              | 1.00E-03                      | 1.00E-03                      | 1.30E-02               | 3.38E-02               | 5.85E-03               | 1.40E-01   | SB-05/10.5-12.5        | 5.98E-02           |                             |                                   |                          |
| Xylenes (total)              | 1330-20-7  | 5             | 3              | 60              | 1.00E-03                      | 1.00E-03                      | 7.50E-02               | 2.27E-01               | 2.10E-02               | 7.80E-01   | SB-05/10.5-12.5        | 3.30E-01           |                             |                                   |                          |

NA - Not Available

**Table I**  
**Statistical Summary and Selection of COPCs in EU5 Soil (0-16' bgs)**  
**Kerr McGee, Hattiesburg, MS**

| Constituent                | 95% UCL<br>mg/kg | Lognormal<br>95% UCL<br>mg/kg | Distribution<br>99% Confidence | Exposure Point<br>Concentration<br>mg/kg | Tier I Restricted<br>Soil TRG<br>mg/kg | Is the Maximum<br>Detected > TRG? | Is the 95% UCL ><br>TRG? |
|----------------------------|------------------|-------------------------------|--------------------------------|--|--|-----------------------------------|--------------------------|
| <b>Screen volatiles</b>    |                  |                               |                                |  |  |                                   |                          |
| 2,4-Dimethylphenol         | 9.9E-01          | 4.25E-01                      | Unknown                        | 4.25E-01                                 | 4.08E-04                               | no                                | no                       |
| 2-Methylnaphthalene        | 1.52E+02         | 9.46E-03                      | Unknown                        | 1.50E+03                                 | 8.18E-04                               | no                                | no                       |
| 2-Methylphenol             | 3.05E-01         | 1.70E-01                      | Unknown                        | 7.30E-02                                 | 1.02E-05                               | no                                | no                       |
| 3- and 4-Methylphenol      | 5.91E-01         | 3.22E-01                      | Unknown                        | 2.10E-01                                 | 1.02E-04                               | no                                | no                       |
| Benzo (a) Pyrene Equiv.    | 4.93E+01         | 7.78E-02                      | Unknown                        | 3.92E+02                                 | 7.84E-01                               | YES                               | YES - COPC               |
| N-nitrosodiphenylamine     | 3.05E-01         | 1.70E-01                      | Unknown                        | 1.70E-01                                 | 1.17E-03                               | no                                | no                       |
| Acenaphthene               | 1.18E+02         | 7.85E-03                      | Unknown                        | 1.20E+03                                 | 1.23E-05                               | no                                | no                       |
| Acenaphthylene             | 7.86E+00         | 6.72E-01                      | Unknown                        | 5.00E-01                                 | 1.23E-05                               | no                                | no                       |
| Anthracene                 | 1.80E+02         | 1.53E-04                      | Unknown                        | 1.80E-03                                 | 6.13E-05                               | no                                | no                       |
| Benzog(hi)perylene         | 1.26E+01         | 1.21E-02                      | Unknown                        | 9.00E+01                                 | 6.13E-04                               | no                                | no                       |
| Bis(2-ethylhexyl)phthalate | 5.74E-01         | 2.87E-01                      | Unknown                        | 1.50E-01                                 | 4.09E-02                               | no                                | no                       |
| Carbazole                  | 6.17E+01         | 1.01E-03                      | Unknown                        | 6.20E-02                                 | 2.86E-02                               | YES                               | no                       |
| Dibenzofuran               | 1.08E+02         | 5.14E-03                      | Unknown                        | 1.10E+03                                 | 8.18E-03                               | no                                | no                       |
| Fluoranthene               | 2.22E+02         | 1.91E-05                      | Unknown                        | 2.00E+03                                 | 8.17E-04                               | no                                | no                       |
| Fluorene                   | 1.47E+02         | 1.41E-04                      | Unknown                        | 1.50E+03                                 | 8.17E-04                               | no                                | no                       |
| Naphthalene                | 3.42E+02         | 5.87E-04                      | Unknown                        | 3.50E+03                                 | 8.24E-02                               | YES                               | no                       |
| Phenanthrene               | 3.90E+02         | 3.05E-05                      | Unknown                        | 4.00E+03                                 | 6.13E-04                               | no                                | no                       |
| Phenol                     | 6.02E-01         | 4.24E-01                      | Unknown                        | 3.80E-01                                 | 1.23E-05                               | no                                | no                       |
| Pyrene                     | 1.56E+02         | 5.67E-04                      | Unknown                        | 1.30E+03                                 | 6.13E-04                               | no                                | no                       |
| <b>Volatiles</b>           |                  |                               |                                |  |  |                                   |                          |
| Acetone                    | 8.01E-02         | 9.07E-01                      | Normal/Lognormal               | 1.00E-01                                 | 1.04E-05                               | no                                | no                       |
| Benzene                    | 5.65E-03         | 2.77E-01                      | Normal/Lognormal               | 7.00E-03                                 | 1.36E+00                               | no                                | no                       |
| Ethylbenzene               | 8.54E-02         | 1.58E-06                      | Normal/Lognormal               | 1.20E-01                                 | 3.95E+02                               | no                                | no                       |
| Styrene                    | 6.28E-02         | 1.19E-04                      | Unknown                        | 1.90E-01                                 | 3.84E+02                               | no                                | no                       |
| Toluene                    | 9.08E-02         | 1.16E+05                      | Lognormal                      | 1.40E-01                                 | 3.80E+01                               | no                                | no                       |
| Xylene (total)             | 5.41E-01         | 2.82E+13                      | Normal/Lognormal               | 7.80E-01                                 | 3.18E+02                               | no                                | no                       |

**Table 12**  
**Summary of Human Health Exposure Parameters**  
**Kerr McGee, Hattiesburg, MS**

| Receptors:                                |                      | Adolescent<br>Visitor | Maintenance<br>Worker | Construction<br>Worker |   |       |
|---|----------------------|-----------------------|-----------------------|------------------------|---|-------|
| Parameter                                 | Units                |                       |                       |                        |   |       |
| Surface area available for exposure       | cm <sup>2</sup> /day | 3052                  | 1                     | 1820                   | 1 | 5300  |
| Total skin surface area                   | cm <sup>2</sup>      | 12768.3               | 2                     | 20000                  | 2 | 20000 |
| Skin surface area available for exposure  | %                    | 23.9%                 | 2                     | 9.1%                   | 2 | 26.5% |
| Adherence factor                          | mg/cm <sup>2</sup>   | 0.026                 | 2                     | 0.054                  | 2 | 0.104 |
| Dermal absorption factor - benzo(a)pyrene |                      | 0.03                  | 3                     | 0.03                   | 3 | 0.03  |
| Dermal absorption factor - other PAHs     |                      | 0.1                   | 3                     | 0.1                    | 3 | 0.1   |
| Exposure time                             | hours/day            | 1                     | 5                     | NA                     |   | NA    |
| Exposure frequency                        | days/year            | 12                    | 5                     | 150                    | 5 | 80    |
| Exposure duration                         | years                | 10                    | 6                     | 25                     | 6 | 1     |
| Body weight                               | kg                   | 45                    | 6                     | 70                     | 6 | 70    |
| Averaging time - noncarcinogenic          | days                 | 3650                  | 7                     | 9125                   | 7 | 365   |
| Averaging time - carcinogenic             | days                 | 25550                 | 7                     | 25550                  | 7 | 25550 |
| Ingestion rate                            | mg/day               | 100                   | 2                     | NA                     |   | NA    |
| Matrix effect - PAHs                      |                      | 0.29                  | 9                     | NA                     |   | NA    |
| Inhalation rate                           | m <sup>3</sup> /day  | NA                    |                       | NA                     |   | 20    |
| Retention factor - semivolatiles          |                      | NA                    |                       | NA                     |   | 0.75  |

NA - Not Applicable

1 Calculated

2 USEPA 1997, Exposure Factors Handbook

3 USEPA 1995, Region III Technical Guidance Manual: Assessing Dermal Exposure to Soil

4 USEPA 1992, Dermal Exposure Assessment

5 Reasonable Maximum

6 USEPA 1995, Region IV

7 USEPA 1991, HHEM Supplemental Guidances

8 International Commission on Radiological Protection, 1968

9 Magee et al., 1996



**Table 13**  
**Particulate Emission Rate for Vehicular Movement and Excavation**  
**Kerr McGee, Hattiesburg, MS**

| <b>Vehicular Movement</b>   |   |  |
|---|---|--|
| $E = k * (5.9) * (s/12)(S/30) * (W/3)^{0.7}((w/4)^{0.5}) * ((365-p)/365) =$ | 16.49   | lbs/vehicle mile   |
| E = 16.49   | particulate emission rate (lbs/vehicle mile - 30 miles travelled total over 80 - 8 hr days) |  |
| k = 0.5   | particle size multiplier  | US EPA AP-42, 1996   |
| s = 50  | percent silt content  | Site Specific  |
| S = 15  | mean vehicle speed (mi/hr)  | US EPA SEAM, 1988  |
| W = 12.5  | mean vehicle weight (ton)   | US EPA SEAM, 1988  |
| w = 8   | mean number of wheels per vehicle   | US EPA SEAM, 1988  |
| p = 110   | mean number of days with ≥ 0.01 inches of precipitation per year                            | US EPA SEAM, 1988  |
| Emission Rate =   | lbs/sec   | (E lbs/mi) * (30 mi/job) * (job/80 days) * (1 day/8 hrs) * (1 hr/3600 sec) |
|   | 2.15E-04  | lbs/sec  |
|   | 9.74E-02  | g/sec  |
|   | 0.0001  | kg/sec   |
| <b>Excavation</b>   |   |  |
| $E = (1.0 * s^{1.5})/M^{1.4} =$   | 7.90E+00  | lbs/hour   |
| E = 7.90E+00  | particulate emission factor (lbs/hr)  |  |
| s = 50  | percent silt content  | Site Specific  |
| M = 15.1  | percent soil moisture content   | Site Specific  |
| Emission Rate =   | 2.20E-03  | lbs/sec  |
|   | 0.996   | g/sec  |
|   | 0.001   | kg/sec   |



**Table 14**  
**Summary of Windrose Data**  
**Kerr McGee, Hattiesburg, MS**

**GRAPHICAL EXPOSURE MODELING SYSTEM**  
**STAR STATION JACKSON/THOMPSON, MS 1974-1978**

| DIRECTION | FREQUENCY | WINDSPEED | DIRECTION | FREQUENCY | WINDSPEED |
|-----------|-----------|-----------|-----------|-----------|-----------|
| N         | 3.33325   | 0.03      | S         | 0.05336   | 3.08      |
| NNE       | 1.89301   | 0.03      | SSW       | 0.09995   | 3.29      |
| NE        | 3.56791   | 0.07      | SW        | 0.10061   | 3.65      |
| ENE       | 0.12132   | 4.04      | WSW       | 0.14723   | 3.93      |
| E         | 0.04843   | 3.39      | W         | 0.05047   | 3.7       |
| ESE       | 0.04328   | 3.12      | WNW       | 0.04341   | 3.51      |
| SE        | 0.03686   | 3         | NW        | 0.02908   | 3.25      |
| SSE       | 0.05274   | 2.99      | NNW       | 0.0406    | 3.26      |

| STABILITY | FREQUENCY | WINDSPEED | AUXILIARY VARIABLES                |       |
|-----------|-----------|-----------|------------------------------------|-------|
| 1         | 259.2     | 0.13      | Afternoon mixing height (meters)   | 1409  |
| 2         | 0.053     | 0.24      | Nocturnal mixing height (meters)   | 472   |
| 3         | 11.3      | 1         | Ambient air temperature (Kelvin)   | 303.6 |
| 4         | 0.01264   | 2.17      | Precipitation frequency (fraction) | 289.8 |
| 5         | 0.08137   | 2.98      | Precipitation intensity (mm/hour)  | 73.66 |
| 6         | 0.1315    | 3.91      | Grand average windspeed (m/s)      | 4.69  |



**Table 15**  
**Summary of Toxicity Values**  
**Kerr McGee, Hattiesburg, MS**

| Chemical                    | Oral Chronic RD mg/kg-day | Inhalation Chronic RD mg/kg-day | Range of Absorption by G.I. | Dermal Chronic RD mg/kg-day | Oral Subchronic RD mg/kg-day | Inhalation Subchronic RD mg/kg-day | Dermal Subchronic RD mg/kg-day | Oral CSF mg/kg-day | Inhalation CSF mg/kg-day | Source |
|-----------------------------|---------------------------|---------------------------------|-----------------------------|-----------------------------|------------------------------|------------------------------------|--------------------------------|--------------------|--------------------------|--------|
|                             |                           |                                 |                             |                             |                              |                                    |                                |                    |                          |        |
| Sensitizers                 |                           |                                 |                             |                             |                              |                                    |                                |                    |                          |        |
| 2-Methylnaphthalene         | 2.00E-02                  | O                               |                             | 0.5                         | Region IV                    | 1.00E-02                           |                                |                    |                          |        |
| Acenaphthylene              |                           |                                 |                             | 0.5                         | Region IV                    |                                    |                                |                    |                          |        |
| Benz(a)pyrene               |                           |                                 |                             | 0.5                         | Region IV                    |                                    |                                |                    |                          |        |
| Benz(e,h,i)perylene         |                           |                                 |                             | 0.5                         | Region IV                    |                                    |                                |                    |                          |        |
| Bis(2-ethylhexyl) phthalate | 2.00E-02                  | IRIS                            |                             | 1.00E-02                    | 2.00E-02                     | W                                  |                                |                    |                          |        |
| Carbazole                   |                           |                                 |                             | 0.5                         | Region IV                    |                                    |                                |                    |                          |        |
| Dibenzofuran                | 4.00E-03                  | E                               |                             | 0.5                         | Region IV                    | 2.00E-03                           |                                |                    |                          |        |
| Fluoranthene                | 4.00E-02                  | IRIS                            |                             | 0.5                         | Region IV                    | 2.00E-02                           | 4.00E-01                       | H                  |                          |        |
| Fluorene                    | 4.00E-02                  | IRIS                            |                             | 0.5                         | Region IV                    | 2.00E-02                           | 4.00E-01                       | H                  |                          |        |
| Naphthalene                 | 2.00E-02                  | IRIS                            | 9.00E-04                    | IRIS                        | 0.5                          | Region IV                          | 1.00E-02                       |                    |                          |        |
| Phenanthrene                |                           |                                 |                             | 0.5                         | Region IV                    |                                    |                                |                    |                          |        |
| Pyrene                      | 3.00E-02                  | IRIS                            |                             | 0.5                         | Region IV                    | 1.50E-02                           | 3.00E-01                       | H                  |                          |        |
|                             |                           |                                 |                             |                             |                              |                                    | 1.50E-01                       |                    |                          |        |

E - EPAN/CEA Regional Support provisional value from Region III RBC Tables, April 1999

H - Values are published in HEAST, 1997

IRIS - Values are available in IRIS, 1999

O - Values are withdrawn from other EPA documents as presented in the Region III RBC Tables, April 1999  
 Region IV - Region IV default value, 1995

W - Withdrawn from IRIS or HEAST



**Table 16**  
**Summary of Hazard and Risk Calculations**  
**Kerr McGee, Hattiesburg, MS**

| Source/Pathway   | Potentially Exposed Population  | Total Hazard Index   | Total Cancer Risk                | Driving Constituent | Table Referenced |
|--|---|----------------------|----------------------------------|---------------------|------------------|
| Dermal Exposure to Sediment in EU1                                   | Visitor<br>Sub-Total  | NA<br>NA             | 1E-09<br>1E-09                   |                     | 17               |
| Dermal Exposure to Surface Water in EU1                              | Visitor<br>Sub-Total  | 8E-06<br>8E-06       | 1E-07<br>1E-07                   |                     | 18               |
| Dermal Exposure to Sediment in EU4                                   | Visitor   | 7E-03                | 4E-07                            |                     | 19               |
| Oral Exposure to Sediment in EU4                                     | Visitor<br>Sub-Total  | 1E-02<br>2E-02       | NA<br>4E-07                      |                     | 20               |
| Dermal Exposure to Surface Water in EU4                              | Visitor<br>Sub-Total  | 2E-04<br>2E-04       | 2E-07<br>2E-07                   |                     | 21               |
| Dermal Exposure to Surface Soil in EU2                               | Visitor<br>Sub-Total  | NA<br>NA             | 1E-08<br>1E-08                   |                     | 22               |
| Dermal Exposure to Surface Soil in EU3                               | Visitor<br>Sub-Total<br>Visitor Total:  | NA<br>NA<br>2E-02    | 2E-09<br>2E-09<br>7E-07          |                     | 23               |
| Dermal Exposure to Surface Soil in EU2                               | Maintenance Worker<br>Sub-Total   | NA<br>NA             | 3E-07<br>3E-07                   |                     | 24               |
| Dermal Exposure to Surface Soil in EU3                               | Maintenance Worker<br>Sub-Total   | NA<br>NA             | 5E-08<br>5E-08                   |                     | 25               |
| Dermal Exposure to Surface Soil in EUS                               | Maintenance Worker<br>Sub-Total<br>Maintenance Worker Total:                          | NA<br>NA<br>NA       | 2E-05<br>2E-05<br>2E-05          | B(a)P Equiv.        | 26               |
| Dermal Exposure to Soil in EU2<br>Inhalation of Fugitive Dust in EU2 | Construction Worker<br>Construction Worker<br>Sub-Total                               | NA<br>NA<br>NA       | 1E-07<br>4E-08<br>1E-07          |                     | 27<br>28         |
| Dermal Exposure to Soil in EUS<br>Inhalation of Fugitive Dust in EUS | Construction Worker<br>Construction Worker<br>Sub-Total<br>Construction Worker Total: | NA<br>NA<br>NA<br>NA | 2E-06<br>7E-07<br>3E-06<br>3E-06 | B(a)P Equiv.        | 29<br>30         |

B(a)P Equiv. = Benzo(a)pyrene Equivalents



Table 17

Dermal Exposure to EU1 Sediment by an Adolescent Visitor (Aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |             |                        |
|---|------------------------------|-------------|------------------------|
|   | BW*AT                        |             |                        |
| Cs - Concentration in sediment =                            | mg/kg                        | chem. spec. |                        |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day         | 3052        | calculated             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>              | 12768.3     | USEPA 1997, EFH        |
| Fs - Fraction of skin surface area available for exposure = |                              | 23.9%       | USEPA 1997, EFH        |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>           | 0.026       | USEPA 1997, EFH        |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                              | 0.03        | USEPA 1995, Region III |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |                              | 0.1         | USEPA 1995, Region III |
| EF - Exposure frequency =                                   | days/year                    | 12          | reasonable assumption  |
| ED - Exposure duration =                                    | years                        | 10          | USEPA 1995, Region IV  |
| CF - Conversion factor =                                    | kg/mg                        | 1.00E-06    |                        |
| BW - Body weight =  | kg                           | 45          | USEPA 1995, Region IV  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                         | 3650        | USEPA 1991, HHEM       |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                         | 25550       | USEPA 1991, HHEM       |

| Constituent             | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|-------------------------|---------------------------------|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
|                         |                                 |                                |                              |              |   |                                   |             |
| <b>Semivolatiles</b>    |                                 |                                |                              |              |   |                                   |             |
| Benzo (a) Pyrene Equiv. | 5.93E-01                        | 1.03E-09                       | NA                           | NA           | 1.47E-10                                | 7.30E+00                          | 1.08E-09    |
| Acenaphthylene          | 7.80E-02                        | 4.52E-10                       | NA                           | NA           | 6.46E-11                                | NA                                | NA          |
| Benzo(ghi)perylene      | 1.80E-01                        | 1.04E-09                       | NA                           | NA           | 1.49E-10                                | NA                                | NA          |
| Phenanthrene            | 1.70E+00                        | 9.85E-09                       | NA                           | NA           | 1.41E-09                                | NA                                | NA          |

NA - Not Available

Total Cancer Risk = 1.08E-09



Table 18

Dermal Exposure to EU1 Surface Water by an Adolescent Visitor (aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cw*SA*Kp*ABS*ET*EF*ED*CF</u> |           |  |
|---|---------------------------------|-----------|--|
|   | BW*AT                           |           |  |
| Cw - Concentration in surface water =                       | mg/L                            | see below |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup>                 | 3052      | calculated                             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                 | 12768.3   | USEPA 1997, EFH                        |
| Fs - Fraction of skin surface area available for exposure = |                                 | 23.9%     | USEPA 1997, EFH                        |
| Kp - Dermal permeability constant =                         | cm/hr                           | see below |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                                 | 0.03      | USEPA 1995, Region III                 |
| ABS <sub>pbh</sub> - Absorption - PAHs =                    |                                 | 0.1       | USEPA 1995, Region III                 |
| ET - Exposure time =  | hrs/day                         | 1         | USEPA 1992, Dermal Exposure Assessment |
| EF - Exposure frequency =                                   | days/year                       | 12        | reasonable assumption                  |
| ED - Exposure duration =                                    | years                           | 10        | USEPA 1995, Region IV                  |
| CF - Conversion factor =                                    | L/cm <sup>2</sup>               | 1.00E-03  |  |
| BW - Body weight =  | kg                              | 45        | USEPA 1995, Region IV                  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                            | 3650      | USEPA 1991, HHEM                       |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                            | 25550     | USEPA 1991, HHEM                       |

| Constituent            | Concentration<br>in Surface<br>Water |             | Average<br>Daily Intake<br>mg/kg-day | Dermal<br>Chronic RfD<br>mg/kg-day | Hazard Index | Average Lifetime<br>Daily Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) |             | Cancer Risk |
|------------------------|--------------------------------------|-------------|--------------------------------------|------------------------------------|--------------|---|---|-------------|-------------|
|                        | mg/L                                 | Kp<br>cm/hr |                                      |                                    |              |   | 1/(mg/kg-day)                           | Cancer Risk |             |
| <b>Semivolatiles</b>   |                                      |             |                                      |                                    |              |   |   |             |             |
| Benzo (a) Pyrene Equiv | 1.21E-03                             | 1.20E+00    | 9.68E-08                             | NA                                 | NA           | 1.38E-08                                      | 7.30E+00                                | 1.01E-07    |             |
| Pyrene                 | 1.00E-03                             | 5.30E-01    | 1.18E-07                             | 1.50E-02                           | 7.87E-06     | 1.69E-08                                      | NA                                      | NA          |             |

NA - Not Available

Total Hazard Index = 7.87E-06

Total Cancer Risk = 1.01E-07



Table 19

Dermal Exposure to EU4 Sediment by an Adolescent Visitor (Aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                      | <u><math>Cs * SA * AH * ABS * EF * ED * CF</math></u> |                        |  |  |
|---|----------------------|---|------------------------|--|--|
|   |                      | BW * AT   |                        |  |  |
| Cs - Concentration in sediment =                            | mg/kg                | chem. spec.   |                        |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 3052  | calculated             |  |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 12768.3   | USEPA 1997, EFH        |  |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 23.9%   | USEPA 1997, EFH        |  |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.026   | USEPA 1997, EFH        |  |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                      | 0.03  | USEPA 1995, Region III |  |  |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |                      | 0.1   | USEPA 1995, Region III |  |  |
| EF - Exposure frequency =                                   | days/year            | 12  | reasonable assumption  |  |  |
| ED - Exposure duration =                                    | years                | 10  | USEPA 1995, Region IV  |  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06  |                        |  |  |
| BW - Body weight =  | kg                   | 45  | USEPA 1995, Region IV  |  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 3650  | USEPA 1991, HHEM       |  |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550   | USEPA 1991, HHEM       |  |  |

| Constituent             | Concentration in Sediment mg/kg | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day |                                   |             | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|-------------------------|---------------------------------|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|-----------------------------------|-------------|
|                         |                                 |                                |                              |              | Lifetime Daily Intake mg/kg-day         | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |                                   |             |
| <b>Semivolatiles</b>    |                                 |                                |                              |              |   |                                   |             |                                   |             |
| Benzo (a) Pyrene Equiv. | 1.99E+02                        | 3.45E-07                       | NA                           | NA           | 4.93E-08                                | 7.30E+00                          | 3.60E-07    |                                   |             |
| 2-Methylnaphthalene     | 1.50E+03                        | 8.70E-06                       | 1.00E-02                     | 8.70E-04     | 1.24E-06                                | NA                                | NA          |                                   |             |
| Acenaphthylene          | 3.50E+01                        | 2.03E-07                       | NA                           | NA           | 2.90E-08                                | NA                                | NA          |                                   |             |
| Benzo(ghi)perylene      | 3.60E+01                        | 2.09E-07                       | NA                           | NA           | 2.98E-08                                | NA                                | NA          |                                   |             |
| Carbazole               | 5.90E+02                        | 3.42E-06                       | NA                           | NA           | 4.89E-07                                | 2.00E-02                          | 9.77E-09    |                                   |             |
| Dibenzofuran            | 9.40E+02                        | 5.45E-06                       | 2.00E-03                     | 2.72E-03     | 7.78E-07                                | NA                                | NA          |                                   |             |
| Fluoranthene            | 1.60E+03                        | 9.27E-06                       | 2.00E-02                     | 4.64E-04     | 1.32E-06                                | NA                                | NA          |                                   |             |
| Fluorene                | 1.20E+03                        | 6.96E-06                       | 2.00E-02                     | 3.48E-04     | 9.94E-07                                | NA                                | NA          |                                   |             |
| Naphthalene             | 3.00E+03                        | 1.74E-05                       | 1.00E-02                     | 1.74E-03     | 2.48E-06                                | NA                                | NA          |                                   |             |
| Phenanthrene            | 3.20E+03                        | 1.85E-05                       | NA                           | NA           | 2.65E-06                                | NA                                | NA          |                                   |             |
| Pyrene                  | 1.00E+03                        | 5.80E-06                       | 1.50E-02                     | 3.86E-04     | 8.28E-07                                | NA                                | NA          |                                   |             |

NA - Not Available

Total Hazard Index = 6.54E-03

Total Cancer Risk = 3.70E-07



Table 20

Oral Exposure to EU4 Sediment by an Adolescent Visitor (Aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =                                 |                           | $\frac{\text{Cd} * \text{IngR} * \text{EF} * \text{ED} * \text{CF} * \text{ME}}{\text{BW} * \text{AT}}$ |                       |              |  |
|--|---------------------------|---|-----------------------|--------------|--|
| Cd - Concentration in sediment =                     | mg/kg                     | see below   |                       |              |  |
| IngR - Ingestion rate for soil =                     | mg/day                    | 100   | USEPA 1997, EFH       |              |  |
| EF - Exposure frequency =                            | days/year                 | 12  | reasonable assumption |              |  |
| ED - Exposure duration =                             | years                     | 10  | USEPA 1995, Region IV |              |  |
| CF - Conversion factor =                             | kg/mg                     | 1.00E-06  |                       |              |  |
| ME <sub>s</sub> - Matrix effect - PAHs =             |                           | 0.29  | Magee, et al., 1996   |              |  |
| BW - Body weight =                                   | kg                        | 45  | USEPA 1995, Region IV |              |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic = | days                      | 3650  | USEPA 1991, HHEM      |              |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =    | days                      | 25550   | USEPA 1991, HHEM      |              |  |
|  |                           |   |                       |              |  |
| Constituent  | Concentration in Sediment | Average Daily Intake  | Oral Chronic RfD      | Hazard Index | Average Lifetime Daily Intake          |
|  | mg/kg                     | mg/kg-day   | mg/kg-day             |              | Oral Cancer Slope Factor 1/(mg/kg-day) |
| Semivolatiles  |                           |   |                       |              | Cancer Risk                            |
| Benzo (a) Pyrene Equiv.                              | 1.99E+02                  | 4.21E-06  | NA                    | NA           | 6.01E-07                               |
| 2-Methylnaphthalene                                  | 1.50E+03                  | 3.18E-05  | 2.00E-02              | 1.59E-03     | 4.54E-06                               |
| Acenaphthylene                                       | 3.50E+01                  | 7.42E-07  | NA                    | NA           | 1.06E-07                               |
| Benzo(ghi)perylene                                   | 3.60E+01                  | 7.63E-07  | NA                    | NA           | 1.09E-07                               |
| Carbazole  | 5.90E+02                  | 1.25E-05  | NA                    | NA           | 1.79E-06                               |
| Dibenzofuran   | 9.40E+02                  | 1.99E-05  | 4.00E-03              | 4.98E-03     | 2.85E-06                               |
| Fluoranthene   | 1.60E+03                  | 3.39E-05  | 4.00E-02              | 8.47E-04     | 4.84E-06                               |
| Fluorene   | 1.20E+03                  | 2.54E-05  | 4.00E-02              | 6.36E-04     | 3.63E-06                               |
| Naphthalene  | 3.00E+03                  | 6.36E-05  | 2.00E-02              | 3.18E-03     | 9.08E-06                               |
| Phenanthrene   | 3.20E+03                  | 6.78E-05  | NA                    | NA           | 9.69E-06                               |
| Pyrene   | 1.00E+03                  | 2.12E-05  | 3.00E-02              | 7.06E-04     | 3.03E-06                               |

NA - Not Applicable

Total Hazard Index = 1.19E-02



Table 21

Dermal Exposure to EU4 Surface Water by an Adolescent Visitor (aged 7-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cw*SA*Kp*ABS*ET*EF*ED*CF</u> |           |  |  |
|---|---------------------------------|-----------|--|--|
|   | <u>BW*AT</u>                    |           |  |  |
| Cw - Concentration in surface water =                       | mg/L                            | see below |  |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup>                 | 3052      | calculated                             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                 | 12768.3   | USEPA 1997, EFH                        |  |
| Fs - Fraction of skin surface area available for exposure = |                                 | 23.9%     | USEPA 1997, EFH                        |  |
| Kp - Dermal permeability constant =                         | cm/hr                           | see below |  |  |
| ABS <sub>rap</sub> - Absorption - B(a)P =                   |                                 | 0.03      | USEPA 1995, Region III                 |  |
| ABS <sub>par</sub> - Absorption - PAHs =                    |                                 | 0.1       | USEPA 1995, Region III                 |  |
| ET - Exposure time =  | hrs/day                         | 1         | USEPA 1992, Dermal Exposure Assessment |  |
| EF - Exposure frequency =                                   | days/year                       | 12        | reasonable assumption                  |  |
| ED - Exposure duration =                                    | years                           | 10        | USEPS 1995, Region IV                  |  |
| CF - Conversion factor =                                    | L/cm <sup>2</sup>               | 1.00E-03  |  |  |
| BW - Body weight =  | kg                              | 45        | USEPA 1995, Region IV                  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                            | 3650      | USEPA 1991, HHEM                       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                            | 25550     | USEPA 1991, HHEM                       |  |

| Constituent                | Concentration<br>in Surface<br>Water<br>mg/L | Kp<br>cm/hr | Average<br>Daily Intake<br>mg/kg-day | Dermal<br>Chronic RfD<br>mg/kg-day | Hazard<br>Index | Average                               |   |             |
|----------------------------|--|-------------|--------------------------------------|------------------------------------|-----------------|---------------------------------------|---|-------------|
|                            |  |             |                                      |                                    |                 | Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>       |  |             |                                      |                                    |                 |                                       |   |             |
| Benzo (a) Pyrene Equiv     | 2.78E-03                                     | 1.20E+00    | 2.23E-07                             | NA                                 | NA              | 3.18E-08                              | 7.30E+00                                | 2.32E-07    |
| bis(2-ethylhexyl)phthalate | 3.00E-03                                     | 3.30E-02    | 2.21E-08                             | 1.00E-02                           | 2.21E-06        | 3.15E-09                              | 1.40E-02                                | 4.41E-11    |
| Carbazole                  | 1.00E-02                                     | 3.57E-02    | 7.96E-08                             | NA                                 | NA              | 1.14E-08                              | 2.00E-02                                | 2.27E-10    |
| Dibenzofuran               | 1.10E-02                                     | 1.51E-01    | 3.71E-07                             | 2.00E-03                           | 1.85E-04        | 5.29E-08                              | NA                                      | NA          |
| Phenanthrene               | 1.70E-02                                     | 2.30E-01    | 8.72E-07                             | NA                                 | NA              | 1.25E-07                              | NA                                      | NA          |

NA - Not Available

Total Hazard Index = 1.87E-04

Total Cancer Risk = 2.33E-07



Table 22

Dermal Exposure to EU2 Surface Soil by an Adolescent Visitor (aged 10-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |             |                        |
|---|------------------------------|-------------|------------------------|
|   | BW*AT                        |             |                        |
| Cs - Concentration in soil =                                | mg/kg                        | chem. spec. |                        |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day         | 3052        | calculated             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>              | 12768.3     | USEPA 1997, EFH        |
| Fs - Fraction of skin surface area available for exposure = |                              | 23.9%       | USEPA 1997, EFH        |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>           | 0.026       | USEPA 1997, EFH        |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                              | 0.03        | USEPA 1995, Region III |
| EF - Exposure frequency =                                   | days/year                    | 12          | reasonable assumption  |
| ED - Exposure duration =                                    | years                        | 10          | USEPA 1995, Region IV  |
| CF - Conversion factor =                                    | kg/mg                        | 1.00E-06    |                        |
| BW - Body weight =  | kg                           | 45          | USEPA 1995, Region IV  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                         | 3650        | USEPA 1991, HHEM       |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                         | 25550       | USEPA 1991, HHEM       |

| Constituent             | Concentration in<br>Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal<br>Chronic RfD<br>mg/kg-day | Hazard Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day |          |          | Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
|-------------------------|-----------------------------------|--------------------------------------|------------------------------------|--------------|--|----------|----------|--------------------------------------|-------------|
|                         |                                   |                                      |                                    |              |  |          |          |                                      |             |
| <b>Semivolatiles</b>    |                                   |                                      |                                    |              |  |          |          |                                      |             |
| Benzo (a) Pyrene Equiv. | 6.37E+00                          | 1.11E-08                             | NA                                 | NA           | 1.58E-09   | 7.30E+00 | 1.16E-08 |                                      |             |

NA - Not Available

Total Cancer Risk = 1.16E-08



Table 23

Dermal Exposure to EU3 Surface Soil by an Adolescent Visitor (aged 10-16 years)

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |             |                        |
|---|------------------------------|-------------|------------------------|
|   | <u>BW*AT</u>                 |             |                        |
| Cs - Concentration in soil =                                | mg/kg                        | chem. spec. |                        |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day         | 3052        | calculated             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>              | 12768.3     | USEPA 1997, EFH        |
| Fs - Fraction of skin surface area available for exposure = |                              | 23.9%       | USEPA 1997, EFH        |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>           | 0.026       | USEPA 1997, EFH        |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                              | 0.03        | USEPA 1995, Region III |
| EF - Exposure frequency =                                   | days/year                    | 12          | reasonable assumption  |
| ED - Exposure duration =                                    | years                        | 10          | USEPA 1995, Region IV  |
| CF - Conversion factor =                                    | kg/mg                        | 1.00E-06    |                        |
| BW - Body weight =  | kg                           | 45          | USEPA 1995, Region IV  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                         | 3650        | USEPA 1991, HHEM       |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                         | 25550       | USEPA 1991, HHEM       |

| Constituent             | Concentration in Soil mg/kg | Average Daily Intake mg/kg-day | Dermal Chronic RfD mg/kg-day | Hazard Index | Average Lifetime Daily Intake mg/kg-day | Cancer Slope Factor 1/(mg/kg-day) | Cancer Risk |
|-------------------------|-----------------------------|--------------------------------|------------------------------|--------------|---|-----------------------------------|-------------|
|                         |                             |                                |                              |              |   |                                   |             |
| <b>Semivolatiles</b>    |                             |                                |                              |              |   |                                   |             |
| Benzo (a) Pyrene Equiv. | 1.12E+00                    | 1.96E-09                       | NA                           | NA           | 2.79E-10                                | 7.30E+00                          | 2.04E-09    |

NA - Not Available

Total Cancer Risk = 2.04E-09



Table 24

*Dermal Exposure to EU2 Surface Soil by a Maintenance Worker*

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  |                      | <u><math>\frac{Cs*SA*AH*ABS*EF*ED*CF}{BW*AT}</math></u> |                        |  |
|---|----------------------|---|------------------------|--|
| Cs - Concentration in soil =                                | mg/kg                | chem. spec.   |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day | 1820  | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>      | 20000   | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                      | 9.1%  | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>   | 0.054   | USEPA 1997, EFH        |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                      | 0.03  | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year            | 150   | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                | 25  | USEPA 1995, Region IV  |  |
| CF - Conversion factor =                                    | kg/mg                | 1.00E-06  |                        |  |
| BW - Body weight =  | kg                   | 70  | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                 | 9125  | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                 | 25550   | USEPA 1991, HHEM       |  |

| Constituent             | Concentration in<br>Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal<br>Chronic RfD<br>mg/kg-day | Hazard Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day |          |          | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
|-------------------------|-----------------------------------|--------------------------------------|------------------------------------|--------------|--|----------|----------|---|-------------|
|                         |                                   |                                      |                                    |              |  |          |          |   |             |
| <b>Semivolatiles</b>    |                                   |                                      |                                    |              |  |          |          |   |             |
| Benzo (a) Pyrene Equiv. | 6.37E+00                          | 1.10E-07                             | NA                                 | NA           | 3.94E-08   | 7.30E+00 | 2.88E-07 |   |             |

NA - Not Available

Total Cancer Risk = 2.88E-07



**Table 25**  
**Dermal Exposure to EU3 Surface Soil by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |             |                        |
|---|------------------------------|-------------|------------------------|
|   | BW*AT                        |             |                        |
| Cs - Concentration in soil =                                | mg/kg                        | chem. spec. |                        |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day         | 1820        | calculated             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>              | 20000       | USEPA 1997, EFH        |
| Fs - Fraction of skin surface area available for exposure = |                              | 9.1%        | USEPA 1997, EFH        |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>           | 0.054       | USEPA 1997, EFH        |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                              | 0.03        | USEPA 1995, Region III |
| EF - Exposure frequency =                                   | days/year                    | 150         | reasonable assumption  |
| ED - Exposure duration =                                    | years                        | 25          | USEPA 1995, Region IV  |
| CF - Conversion factor =                                    | kg/mg                        | 1.00E-06    |                        |
| BW - Body weight =  | kg                           | 70          | USEPA 1995, Region IV  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                         | 9125        | USEPA 1991, HHEM       |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                         | 25550       | USEPA 1991, HHEM       |

| Constituent             | Concentration<br>in Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal                   |              | Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>1/(mg/kg-day) | Cancer Risk |
|-------------------------|-----------------------------------|--------------------------------------|--------------------------|--------------|---------------------------------------|-------------------------------|-------------|
|                         |                                   |                                      | Chronic RfD<br>mg/kg-day | Hazard Index |                                       |                               |             |
| <b>Semivolatiles</b>    |                                   |                                      |                          |              |                                       |                               |             |
| Benzo (a) Pyrene Equiv. | 1.12E+00                          | 1.95E-08                             | NA                       | NA           | 6.95E-09                              | 7.30E+00                      | 5.08E-08    |

NA - Not Available

Total Cancer Risk = 5.08E-08



**Table 26**  
**Dermal Exposure to EU5 Surface Soil by a Maintenance Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |                |                              |
|---|------------------------------|----------------|------------------------------|
|   | BW*AT                        |                |                              |
| Cs - Concentration in soil =                                | mg/kg                        | chem. spec.    |                              |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day         | 1820           | calculated                   |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>              | 20000          | USEPA 1997, EFH              |
| Fs - Fraction of skin surface area available for exposure = |                              | 9.1%           | USEPA 1997, EFH              |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>           | 0.054          | USEPA 1997, EFH              |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                              | 0.03           | USEPA 1995, Region III       |
| EF - Exposure frequency =                                   | days/year                    | 150            | reasonable assumption        |
| ED - Exposure duration =                                    | years                        | 25             | USEPA 1995, Region IV        |
| CF - Conversion factor =                                    | kg/mg                        | 1.00E-06       |                              |
| BW - Body weight =  | kg                           | 70             | USEPA 1995, Region IV        |
| AT <sub>a</sub> - Averaging time - noncarcinogenic =        | days                         | 9125           | USEPA 1991, HHEM             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                         | 25550          | USEPA 1991, HHEM             |
| Average   |                              |                |                              |
| Concentration   |                              | Lifetime Daily | Cancer Slope                 |
| in Soil   |                              | Intake         | Factor                       |
| Constituent   | mg/kg                        | mg/kg-day      | 1/(mg/kg-day)                |
| Semivolatiles   |                              |                |                              |
| Benzo (a) Pyrene Equiv.                                     | 3.52E+02                     | 6.09E-06       | 2.18E-06                     |
|   |                              |                | 7.30E+00                     |
|   |                              |                | Total Cancer Risk = 1.59E-05 |

NA - Not Available



Table 27

Dermal Exposure to EU2 Soil by a Construction Worker

Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |             |                        |  |
|---|------------------------------|-------------|------------------------|--|
|   | BW*AT                        |             |                        |  |
| Cs - Concentration in soil =                                | mg/kg                        | chem. spec. |                        |  |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day         | 5300        | calculated             |  |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>              | 20000       | USEPA 1997, EFH        |  |
| Fs - Fraction of skin surface area available for exposure = |                              | 26.5%       | USEPA 1997, EFH        |  |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>           | 0.104       | USEPA 1997, EFH        |  |
| ABS <sub>bap</sub> - Absorption - B(a)P =                   |                              | 0.03        | USEPA 1995, Region III |  |
| ABS <sub>pah</sub> - Absorption - PAHs =                    |                              | 0.1         | USEPA 1995, Region III |  |
| EF - Exposure frequency =                                   | days/year                    | 80          | reasonable assumption  |  |
| ED - Exposure duration =                                    | years                        | 1           | reasonable assumption  |  |
| CF - Conversion factor =                                    | kg/mg                        | 1.00E-06    |                        |  |
| BW - Body weight =  | kg                           | 70          | USEPA 1995, Region IV  |  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                         | 365         | USEPA 1991, HHEM       |  |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                         | 25550       | USEPA 1991, HHEM       |  |

| Constituent             | Concentration in<br>Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal                         |                  | Hazard<br>Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
|-------------------------|-----------------------------------|--------------------------------------|--------------------------------|------------------|-----------------|--|---|-------------|
|                         |                                   |                                      | Subchronic<br>RfD<br>mg/kg-day | RfD<br>mg/kg-day |                 |  |   |             |
| <b>Semivolatiles</b>    |                                   |                                      |                                |                  |                 |  |   |             |
| Benzo (a) Pyrene Equiv. | 1.91E+01                          | 9.91E-07                             | NA                             | NA               | NA              | 1.42E-08   | 7.30E+00                                | 1.03E-07    |

NA - Not Available

Total Cancer Risk = 1.03E-07



Table 28

Exposure to Construction Workers from Inhalation of Fugitive Dust in EU2  
Kerr McGee, Hattiesburg, MS

| Intake (mg/kg-day) = $\frac{Ca * InhR * EF * ED * RF}{BW * AT}$   |  |  |  |  |  |  |
|---|--|--|--|--|--|--|
| Ca - Concentration in air =<br>InhR - Inhalation Rate =<br>EF - Exposure Frequency =<br>ED - Exposure Duration =<br>RF <sub>s</sub> - Retention Factor - semivolatiles =<br>AT <sub>b</sub> - Averaging Time noncarcinogenic =<br>AT <sub>c</sub> - Averaging Time carcinogenic =<br>BW - Body Weight = | chem.spec.<br>mg/m <sup>3</sup><br>m <sup>3</sup> /shift<br>shifts/year<br>years | 20<br>80<br>1                                | USEPA 1995, Region IV<br>reasonable assumption<br>reasonable assumption<br>ICRP, 1968<br>USEPA 1991, HHEM<br>USEPA 1991, HHEM<br>USEPA 1995, Region IV | Ei =<br>Ei / (Hb * W * V)<br>Emission Rate of Component (mg/sec) = see below<br>Hb - Downwind Ht (m) = 4.81<br>W - Width (m) = 50<br>V - Wind speed (m/sec) = 4.69<br>Length (downwind distance) (m) = 50<br>r - Roughness Ht. (m) = 0.20<br>z - downwind distance (m) = 50<br>2 = 6.25r[Hb/r * Ln(Hb/r) - 1.58*Hb/r + 1.58] | Ca = Concentration in Air (mg/m <sup>3</sup> ) = Ei / (Hb * W * V) |  |
| E <sub>i</sub> - Emission Rate (mg/sec) = $Cs * (PERv1 + PERe)$   |  |  |  |  |  |  |
| Cs - Concentration in soil =<br>Cs - Concentration in soil =<br>mg/kg   | mg/kg  |  | chem.spec.   |  |  |  |
| Concentration<br>in Soil<br>mg/Kg   | Emision Rate<br>mg/sec   | Concentration<br>in Air<br>mg/m <sup>3</sup> | Average Daily<br>Intake<br>mg/kg-day   | Inhalation<br>Subchronic RID<br>mg/kg-day  | Hazard<br>Index  | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day |
| Chemicals   |  |  |  |  |  |  |
| Semivolatiles   |  |  |  |  |  |  |
| Benzo (a) Pyrene Equiv.   | 1.91E+01   | 2.09E-02                                     | 1.86E-05   | 8.72E-07   | NA   | NA   |
| NA - Not Available  |  |  |  |  |  |  |
|   |  |  |  |  | Total Cancer Risk:   | 3.86E-08   |
|   |  |  |  |  |  | 3.86E-08   |

**Table 29**

**Dermal Exposure to EU5 Soil by a Construction Worker**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u><math>\frac{Cs * SA * AH * ABS * EF * ED * CF}{BW * AT}</math></u> |                                      |  |                 |  |   |             |
|---|---|--------------------------------------|--|-----------------|--|---|-------------|
| Cs - Concentration in soil =                                | mg/kg   | chem. spec.                          |  |                 |  |   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day  | 5300                                 | calculated                               |                 |  |   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>   | 20000                                | USEPA 1997, EFH                          |                 |  |   |             |
| Fs - Fraction of skin surface area available for exposure = |   | 26.5%                                | USEPA 1997, EFH                          |                 |  |   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>  | 0.104                                | USEPA 1997, EFH                          |                 |  |   |             |
| ABS <sub>top</sub> - Absorption - B(a)P =                   |   | 0.03                                 | USEPA 1995, Region III                   |                 |  |   |             |
| EF - Exposure frequency =                                   | days/year   | 80                                   | reasonable assumption                    |                 |  |   |             |
| ED - Exposure duration =                                    | years   | 1                                    | reasonable assumption                    |                 |  |   |             |
| CF - Conversion factor =                                    | kg/mg   | 1.00E-06                             |  |                 |  |   |             |
| BW - Body weight =  | kg  | 70                                   | USEPA 1995, Region IV                    |                 |  |   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days  | 365                                  | USEPA 1991, HHEM                         |                 |  |   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days  | 25550                                | USEPA 1991, HHEM                         |                 |  |   |             |
|   |   |                                      |  |                 |  |   |             |
| Constituent   | Concentration in<br>Soil<br>mg/kg                                     | Average Daily<br>Intake<br>mg/kg-day | Dermal<br>Subchronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |   |                                      |  |                 |  |   |             |
| Benzo (a) Pyrene Equiv.                                     | 3.52E+02  | 1.82E-05                             | NA                                       | NA              | 2.60E-07   | 7.30E+00                                | 1.90E-06    |

NA - Not Available

Total Cancer Risk = 1.90E-06





**Table 30**  
**Exposure to Construction Workers from Inhalation of Fugitive Dust in EUS**  
**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg/day) =                                 | $\frac{Ca * InhR * EF*ED*RF}{BW * AT}$ | Inhalation                             |                                |                                     |               |   |  |
|--|--|--|--------------------------------|-------------------------------------|---------------|---|--|
| Concentration in air =                               | mg/m <sup>3</sup>                      | Concentration in Air                   | Average Daily Intake           | Inhalation Subchronic RfD           | Hazard Index  | Average Lifetime Daily Intake           | Inhalation Cancer Slope Factor               |
| InhR - Inhalation Rate =                             | m <sup>3</sup> /shift                  | mg/sec                                 | mg/kg-day                      | mg/kg-day                           | 1/(mg/kg-day) | mg/kg-day                               | 1/(mg/kg-day)                                |
| EF - Exposure Frequency =                            | shifts/year                            | 80                                     |                                |                                     |               |   |  |
| ED - Exposure Duration =                             | years                                  | 1                                      |                                |                                     |               |   |  |
| RF <sub>s</sub> - Retention Factor - semivolatiles = |  | 0.75                                   |                                |                                     |               |   |  |
| AT <sub>a</sub> - Averaging Time noncarcinogenic =   | days                                   | 365                                    |                                |                                     |               |   |  |
| AT <sub>c</sub> - Averaging Time carcinogenic =      | days                                   | 25550                                  |                                |                                     |               |   |  |
| BW - Body Weight =                                   | kg                                     | 70                                     |                                |                                     |               |   |  |
| E <sub>i</sub> - Emission Rate (mg/sec) =            | $Cs * (PERv + PERe)$                   |  |                                |                                     |               |   |  |
| Cs - Concentration in soil =                         | mg/kg                                  | see below                              |                                |                                     |               |   |  |
|  |  |  |                                |                                     |               |   |  |
| Concentration in Soil                                | Emission Rate mg/kg                    | Concentration in Air mg/m <sup>3</sup> | Average Daily Intake mg/kg-day | Inhalation Subchronic RfD mg/kg-day | Hazard Index  | Average Lifetime Daily Intake mg/kg-day | Inhalation Cancer Slope Factor 1/(mg/kg-day) |
| Chemicals  | mg/sec                                 |  |                                |                                     |               |   |  |
| Semivolatiles  |  |  |                                |                                     |               |   |  |
| Benzo (a) Pyrene Equiv.                              | 3.52E-02                               | 3.85E-01                               | 3.41E-04                       | 1.60E-05                            | NA            | NA                                      | 7.10E-07                                     |
| NA - Not Available                                   |  |  |                                |                                     |               |   |  |
|  |  |  |                                |                                     |               |   |  |
|  |  |  |                                |                                     |               |   | Total Cancer Risk: 7.10E-07                  |

**Table 31**  
**Statistical Summary of Soil Samples in EU5 Post Remediation**  
**Kerr McGee, Hattiesburg, MS**

**Maintenance Worker Scenario:**

Data from sample locations GEO-19/0-1', GEO-21/0-1', and GEO-33/0-1 were eliminated from this statistical analysis.

| Constituent             | Total         |                         |      | Minimum<br>Detection<br>Limit<br>mg/kg | Mean<br>mg/kg | Logarithmic<br>Mean<br>mg/kg | Maximum<br>Detected<br>Qualifer<br>mg/kg | Maximum<br>Detected<br>Qualifer<br>mg/kg | Location of<br>Maximum<br>Concentration |
|-------------------------|---------------|-------------------------|------|--|---------------|------------------------------|--|--|---|
|                         | CAS<br>Number | Number<br>of<br>Samples | Hits |  |               |                              |  |  |   |
| Senivolatiles           |               |                         |      |  |               |                              |  |  |   |
| Benzo (a) Pyrene Equiv. | -             | 6                       | 5    | 83.33                                  | 4.39E-02      | 4.39E-02                     | 5.09E-01                                 | 3.95E+00                                 | 1.43E+00                                |
|                         |               |                         |      |  |               |                              |  |  | 1.29E+01                                |
|                         |               |                         |      |  |               |                              |  |  | GEO-30/0-1'                             |

| Constituent             | Standard  |          |          | 95% Lognormal    |              |                | Exposure Point |       |       |
|-------------------------|-----------|----------|----------|------------------|--------------|----------------|----------------|-------|-------|
|                         | Deviation | UCL      | mg/kg    | 95% UCL          | Distribution | 99% Confidence | Concentration  | mg/kg | mg/kg |
| Senivolatiles           |           |          |          |                  |              |                |                |       |       |
| Benzo (a) Pyrene Equiv. | 4.79E+00  | 7.90E+00 | 1.28E+04 | Normal/Lognormal |              |                | 1.29E+01       |       |       |

**Construction Worker Scenario:**

Data from sample locations GEO-21/0-1' and GEO-21/2-3 were eliminated from this statistical analysis.

| Constituent             | Total         |                         |      | Minimum<br>Detection<br>Limit<br>mg/kg | Mean<br>mg/kg | Logarithmic<br>Mean<br>mg/kg | Maximum<br>Detected<br>Qualifer<br>mg/kg | Maximum<br>Detected<br>Qualifer<br>mg/kg | Location of<br>Maximum<br>Concentration |
|-------------------------|---------------|-------------------------|------|--|---------------|------------------------------|--|--|---|
|                         | CAS<br>Number | Number<br>of<br>Samples | Hits |  |               |                              |  |  |   |
| Senivolatiles           |               |                         |      |  |               |                              |  |  |   |
| Benzo (a) Pyrene Equiv. | -             | 32                      | 20   | 62.5                                   | 4.39E-02      | 7.60E-02                     | 8.74E-02                                 | 8.51E+00                                 | 5.44E-01                                |
|                         |               |                         |      |  |               |                              |  |  | 8.36E+01                                |
|                         |               |                         |      |  |               |                              |  |  | GEO19/0-1'                              |

| Constituent             | Standard  |          |          | 95% Lognormal |              |                | Exposure Point |       |       |
|-------------------------|-----------|----------|----------|---------------|--------------|----------------|----------------|-------|-------|
|                         | Deviation | UCL      | mg/kg    | 95% UCL       | Distribution | 99% Confidence | Concentration  | mg/kg | mg/kg |
| Senivolatiles           |           |          |          |               |              |                |                |       |       |
| Benzo (a) Pyrene Equiv. | 2.07E-01  | 1.47E+01 | 1.29E+02 | Unknown       |              |                | 8.36E+01       |       |       |

**Table 32**

**Dermal Exposure to EU5 Surface Soil by a Maintenance Worker Post Remediation**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u> |             |                        |
|---|------------------------------|-------------|------------------------|
|   | BW*AT                        |             |                        |
| Cs - Concentration in soil =  | mg/kg                        | chem. spec. |                        |
| SA - Surface area available for exposure =                              | cm <sup>2</sup> /day         | 1820        | calculated             |
| SA <sub>t</sub> - Total skin surface area =                             | cm <sup>2</sup>              | 20000       | USEPA 1997, EFH        |
| F <sub>s</sub> - Fraction of skin surface area available for exposure = |                              | 9.1%        | USEPA 1997, EFH        |
| AH - Adherence factor =   | mg/cm <sup>2</sup>           | 0.054       | USEPA 1997, EFH        |
| ABS <sub>bap</sub> - Absorption - B(a)P =                               |                              | 0.03        | USEPA 1995, Region III |
| EF - Exposure frequency =   | days/year                    | 150         | reasonable assumption  |
| ED - Exposure duration =  | years                        | 25          | USEPA 1995, Region IV  |
| CF - Conversion factor =  | kg/mg                        | 1.00E-06    |                        |
| BW - Body weight =  | kg                           | 70          | USEPA 1995, Region IV  |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =                    | days                         | 9125        | USEPA 1991, HHEM       |
| AT <sub>c</sub> - Averaging time - carcinogenic =                       | days                         | 25550       | USEPA 1991, HHEM       |

| Constituent             | Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal<br>Chronic RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily |                                      |             |
|-------------------------|---------------|--------------------------------------|------------------------------------|-----------------|---------------------------|--------------------------------------|-------------|
|                         |               |                                      |                                    |                 | Intake<br>mg/kg-day       | Cancer Slope Factor<br>1/(mg/kg-day) | Cancer Risk |
| <b>Semivolatiles</b>    |               |                                      |                                    |                 |                           |                                      |             |
| Benzo (a) Pyrene Equiv. | 1.29E+01      | 2.24E-07                             | NA                                 | NA              | 8.00E-08                  | 7.30E+00                             | 5.84E-07    |

NA - Not Available

Total Cancer Risk = 5.84E-07

Data from sample locations GEO-19/0-1', GEO-21/0-1', and GEO-33/0-1 were eliminated from this scenario.



**Table 33**

**Dermal Exposure to EU5 Soil by a Construction Worker Post Remediation**

**Kerr McGee, Hattiesburg, MS**

| Intake (mg/kg-day) =  | <u>Cs*SA*AH*ABS*EF*ED*CF</u>      |                                      |  |                 |  |   |             |
|---|-----------------------------------|--------------------------------------|--|-----------------|--|---|-------------|
|   | BW*AT                             |                                      |  |                 |  |   |             |
| Cs - Concentration in soil =                                | mg/kg                             | chem. spec.                          |  |                 |  |   |             |
| SA - Surface area available for exposure =                  | cm <sup>2</sup> /day              | 5300                                 | calculated                               |                 |  |   |             |
| SA <sub>t</sub> - Total skin surface area =                 | cm <sup>2</sup>                   | 20000                                | USEPA 1997, EFH                          |                 |  |   |             |
| Fs - Fraction of skin surface area available for exposure = |                                   | 26.5%                                | USEPA 1997, EFH                          |                 |  |   |             |
| AH - Adherence factor =                                     | mg/cm <sup>2</sup>                | 0.104                                | USEPA 1997, EFH                          |                 |  |   |             |
| ABS <sub>top</sub> - Absorption - B(a)P =                   |                                   | 0.03                                 | USEPA 1995, Region III                   |                 |  |   |             |
| EF - Exposure frequency =                                   | days/year                         | 80                                   | reasonable assumption                    |                 |  |   |             |
| ED - Exposure duration =                                    | years                             | 1                                    | reasonable assumption                    |                 |  |   |             |
| CF - Conversion factor =                                    | kg/mg                             | 1.00E-06                             |  |                 |  |   |             |
| BW - Body weight =  | kg                                | 70                                   | USEPA 1995, Region IV                    |                 |  |   |             |
| AT <sub>n</sub> - Averaging time - noncarcinogenic =        | days                              | 365                                  | USEPA 1991, HHEM                         |                 |  |   |             |
| AT <sub>c</sub> - Averaging time - carcinogenic =           | days                              | 25550                                | USEPA 1991, HHEM                         |                 |  |   |             |
| <hr/>   |                                   |                                      |  |                 |  |   |             |
| Constituent   | Concentration in<br>Soil<br>mg/kg | Average Daily<br>Intake<br>mg/kg-day | Dermal<br>Subchronic<br>RfD<br>mg/kg-day | Hazard<br>Index | Average<br>Lifetime Daily<br>Intake<br>mg/kg-day | Cancer Slope<br>Factor<br>1/(mg/kg-day) | Cancer Risk |
| Semivolatiles   |                                   |                                      |  |                 |  |   |             |
| Benzo (a) Pyrene Equiv.                                     | 8.36E+01                          | 4.33E-06                             | NA                                       | NA              | 6.18E-08   | 7.30E+00                                | 4.51E-07    |

NA - Not Available

Total Cancer Risk = 4.51E-07

Data from sample locations GEO-21/0-1', and GEO-21/2-3 were eliminated from this scenario.





FILE COPY

**PROPOSED WORK PLAN FOR DEVELOPING SITE-SPECIFIC,  
RISK-BASED CLEANUP GOALS FOR THE FORMER  
GULF STATES CREOSOTE SITE IN HATTIESBURG, MISSISSIPPI**

*Privileged and Confidential  
Prepared at the Request of Legal Counsel*

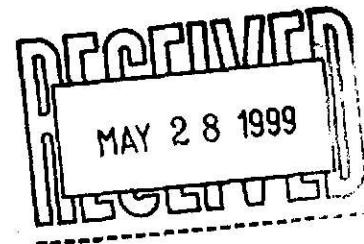
May 25, 1999

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### **Appendix A Qualifications and Experience**

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### **Appendix B Professional Profiles of Key Project Personnel**

## **1.0 Introduction**

### **1.1 Project Background**

In January 1997, Kerr-McGee Chemical Corporation (KMCC) and the Mississippi Department of Environmental Quality (MDEQ) entered into an agreement for the investigation of the Former Gulf States Creosote site in Hattiesburg, Mississippi. KMCC has subsequently conducted extensive investigations to evaluate site conditions and to determine the nature and extent of affected media at the site. All investigative activities have been conducted in accordance with MDEQ-approved work plans.

Work completed at the site in 1997 and 1998 includes the following:

- Characterization of the site-wide stratigraphy to depths of greater than 60 feet below grade;
- Collection and laboratory analysis of over 100 surface and subsurface soil samples;
- Installation of 12 groundwater monitoring wells and 13 temporary well points;
- Collection and laboratory analysis of over 30 groundwater samples;
- Collection of hydrogeologic data (e.g., groundwater flow rate and direction); and
- Collection and laboratory analysis of surface water and sediment samples from two separate drainage pathways.

In a letter dated April 20, 1999, MDEQ concurred that the remedial investigation is complete until such time that additional data needs are identified. The next phase in the remedial evaluation process is to conduct a baseline risk assessment and to develop target cleanup levels for site media, if necessary. This step is provided for in Section 3 of the document *Guidance for the Remediation of Uncontrolled Hazardous Substance Sites in Mississippi*, MDEQ, September 1990.

### **1.2 Risk Assessment Methodologies**

This Work Plan presents the methodologies proposed to be used in conducting a human health risk assessment for the Former Gulf States Creosote site in Hattiesburg, Mississippi. The baseline risk assessment will serve two purposes. First, this assessment will provide a

quantitative analysis of the potential human health risks resulting from residual chemicals found in various media at the site. Second, the risk assessment will be used as a basis for establishing cleanup goals, if necessary.

The five basic phases of a risk assessment described in the US EPA's *Risk Assessment Guidance for Superfund, Volume I, Human Health Evaluation Manual (RAGS)* (US EPA, 1989) are:

- 1) Data evaluation - the process of analyzing site environmental data relevant to potential human health impacts;
- 2) Exposure assessment - the identification of relevant exposure pathways and populations at probable risk, estimation of exposure point concentrations, and estimation of average daily intakes;
- 3) Toxicity assessment - the determination of chemical dose-response relationships and daily intake levels at which no adverse effects or unacceptable cancer risks can be reasonably anticipated to result;
- 4) Risk characterization - a comparison of estimated daily chemical intake levels with acceptable daily intake levels to generate quantitative expressions of hazard and the upper limits of probability of causing cancer (for carcinogens); and
- 5) Uncertainty Analysis - a discussion of the factors throughout the risk assessment process that contribute uncertainty. Depending on the methods and values used, these factors may over- or underestimate risk.

### 1.3 Relevant Guidance

This risk assessment will follow all relevant current United States Protection Agency (US EPA) and Mississippi Department of Environmental Quality (MDEQ) guidance. Key guidance documents that may be referenced to define the risk assessment process, format, and procedures include:

- *Guidance for Remediation of Uncontrolled Hazardous Substance Sites in Mississippi (MDEQ, 1990);*
- *Risk Assessment Guidance for Superfund, Volume 1, Human Health Evaluation Manual/Part A (RAGS/Part A)* (US EPA, 1989);

- *Risk Assessment Guidance for Superfund, Human Health Evaluation Manual/ Part B, Development of Risk-based Preliminary Remediation Goals (RAGS/Part B)* (US EPA, 1991);
- *Human Health Evaluation Manual, Supplemental Guidance: "Standard Default Exposure Factors"* (US EPA, 1991);
- *Exposure Factors Handbook* (US EPA, 1997);
- *Guidelines for Exposure Assessment* (US EPA 1992);
- *Dermal Exposure Assessment: Principles and Applications* (US EPA, 1992);
- *Guidance for Data Usability in Risk Assessment (Part A)* (US EPA, 1992);
- *Supplemental Guidance to RAGS: Calculating the Concentration Term* (US EPA, 1992); and
- *Guiding Principles for Monte Carlo Analysis* (US EPA, 1997).

These documents are not listed in a hierarchical manner; other US EPA guidance documents, MDEQ guidances, or peer-reviewed technical papers may also be referenced in the risk assessment report.

## **2.0 The Human Health Risk Assessment Process**

The methodologies to be employed in each of the five stages of the risk evaluation are detailed in the following sections.

### **2.1 Data Evaluation**

Analytical data to be utilized during the risk assessment will be compiled into a database format and analyzed to generate descriptive statistics that will then be used to select the constituents of potential concern (COPCs) for quantitative evaluation in the risk assessment. Chemicals that are not detected in any sample will be eliminated from the quantitative assessment in accordance with US EPA

guidelines. If a chemical is not detected in an individual sample but is detected in any sample in the media of concern, the chemical will be conservatively assumed present at a concentration equivalent to one-half its sample quantitation limit (SQL). The selection of COPCs may involve constituent frequency of detection and comparison of site concentrations to risk-based screening levels or other appropriate MDEQ-approved criteria. A summary table of detected chemicals with their respective number of samples taken, the number of positive detections, detection frequency, mean concentrations, logarithmic mean concentrations, standard errors, 95% upper confidence limits (UCLs; for normal and lognormal distributions), maximum concentrations detected, and distribution type will be presented. The COPC screening process will also be presented on these tables.

## **2.2 Exposure Assessment**

Subsequent to a thorough review of existing site data and other available site-related information (including the remedial investigation) and the performance of a site reconnaissance visit, a conceptual site model (CSM) will be developed for the Hattiesburg facility. The CSM will help to identify reasonable and realistic human exposure pathways potentially present at the site, currently or in the future. These various exposure scenarios will be employed in the risk assessment to generate conservative estimates of exposure and risks to individuals accessing the site for occupational or recreational purposes. The objective of the exposure assessment is to estimate the type, magnitude, frequency, and duration of exposures to site-related COPCs.

The US EPA guidelines recommend evaluation of only those land-use scenarios that can be reasonably foreseen in the future. Populations who could potentially be exposed to COPCs in affected media at this site will be identified based on the current and potential future use of the site. Those individuals currently accessing the site or potentially accessing the site in the future will be evaluated during the risk assessment process. The specific methodologies for the identification of potential exposure pathways and the elimination of incomplete exposure pathways will be



extensively detailed in the risk assessment report. These methodologies will be consistent with the US EPA *Risk Assessment Guidance for Superfund* (US EPA, 1989), *Guidelines for Exposure Assessment* (US EPA, 1992), *Superfund Exposure Assessment Manual* (US EPA, 1988), *Exposure Factors Handbook* (US EPA, 1997), and other related guidances listed above.

The assumptions and methodologies employed in this analysis will be fully consistent with current US EPA and MDEQ relevant guidelines for exposure assessments. If appropriate, Environmental Standards will use MDEQ- or US EPA-recommended default exposure parameter values. Conservative, but realistic, site-specific assumptions will be used for those exposure parameters where default assumptions do not accurately characterize potential exposures at the site. Appropriate justification of the use of all site-specific exposure assumptions will be included in the risk assessment report.

Chemical intake is expressed as the amount of the agent at the exchange boundaries of an organism (i.e., skin, lungs, gut) that is available for systemic absorption. If the exposure occurs over time, the total exposure is divided by the time period of interest and a body weight (expressed in kilograms) to obtain an average exposure rate (e.g., mg/kg-day). The general equation, as defined by the US EPA, for estimating a time-weighted average intake is:

$$\text{Intake (mg/kg - day)} = \frac{C \times IR \times EF \times ED}{BW \times AT}$$

where:

- C = chemical concentration at the exposure point;
- IR = intake rate (i.e., mg soil/day);
- EF = exposure frequency (days/year);
- ED = exposure duration (years);
- BW = body weight of exposed individual (kg); and
- AT = averaging time (period over which exposure is averaged, usually measured in days).



This basic equation will be modified according to US EPA guidelines to estimate exposure-route-specific intakes, such as dermal absorption or incidental ingestion. Parameter values to be incorporated into this equation for the dermal exposure to soil pathway include skin surface area, adherence factor, and absorption factor. For the soil ingestion exposure pathway, gastrointestinal matrix effect and fraction of soil ingested at the site will be incorporated into the general intake paradigm. To estimate intake via inhalation of volatile organic compound (VOC) vapors and entrained dust, additional exposure parameters will include retention factor and the fraction of PM<sub>15</sub> size particles (particles less than 15 µm in diameter) that are respirable. Oral exposures to groundwater, if appropriate, will use the general intake equation without the incorporation of additional parameters.

If other complete exposure pathways or media of concern are identified, the intake equation will be adjusted to assess those pathways in a manner similar to that above. Any alterations to the general intake equation will be fully documented in the risk assessment report.

The values and assumptions applied for each exposure pathway identified at the site will be presented in a table in the risk assessment report for ease of interpretation. The exposure-point concentration will be determined based on the fit of the chemical-specific data to specified distributions.

The results of the exposure assessment (estimates of daily intake levels) will be combined with chemical-specific toxicity information to characterize potential risks.

### **2.3      Toxicity Assessment**

Toxicity assessment involves the evaluation of available toxicity information for the constituents of concern and characterization of the relationship between exposure concentration and the incidence of adverse health effects. Toxicity values derived from this dose-response relationship will be used



to estimate the potential for the occurrence of adverse effects in individuals exposed to various constituent levels.

Adverse effects can be caused by acute exposure, which is a single or short-term exposure to a toxic substance, or by chronic exposure to constituents on a continuous or repeated basis over an extended period of time. "Acceptable" exposures to non-carcinogens are considered to be exposures without any anticipated adverse effects. Such exposure levels are commonly expressed as Reference Doses (RfDs). An acceptable exposure level is calculated to provide an "adequate margin of safety." Chemical-specific RfD values will be compared to site-specific chemical intake levels for each receptor to determine the potential for the occurrence of non-carcinogenic adverse health effects. This comparison is discussed in more detail in Section 2.4.

The US EPA has derived carcinogenic slope factors (CSFs) for both oral and inhalation pathways, and these will be utilized to quantitatively estimate risks posed to individuals via contact with carcinogenic chemicals. These CSFs will be multiplied by the chemical intake level for each receptor to determine the probability that an individual will experience carcinogenic effects under site-specific exposure assumptions. This calculation is discussed in more detail in Section 2.4. In accordance with US EPA guidance (*Dermal Exposure Assessment, Principles and Applications*, 1992), oral CSFs will be used in the absence of dermal CSFs for the dermal exposure route.

A number of sources of toxicity information (RfDs or CSFs) exist, and these sources vary with regard to the availability and strength of supporting evidence. The protocol detailed below has been established for the determination of toxicity indices; this protocol defines a hierarchy of sources to be consulted and the methodology for the determination of toxicity values and was developed in accordance with current US EPA methodology adopted and/or developed by the National Academy of Sciences. Toxicity values for the chemicals of potential concern at the Site will be obtained with reference to the following hierarchy of sources:



- 1) Toxicity values will be obtained from the *Integrated Risk Information System* (IRIS) database. This database contains the Reference Doses (RfDs) and Cancer Slope Factors (CSFs), which have been verified by the US EPA RfD and Carcinogen Risk Assessment Verification Endeavor (CRAVE) work groups, and is, thus, the agency's preferred source for toxicity values. IRIS supersedes all other information sources.
- 2) For toxicity values that are unavailable on IRIS, the most current source of information is the *Health Effects Assessment Summary Tables* (HEAST), published by the US EPA. HEAST contains interim, as well as verified RfDs and CPFs. Supporting toxicity information for verified values is provided in an extensive reference section of HEAST.
- 3) Toxicity values that cannot be determined in either IRIS or HEAST will be derived from data in toxicological profiles for individual compounds, as compiled by the Agency for Toxic Substances and Disease Registry (ATSDR). These documents provide results from a number of toxicological studies, as well as the methodologies and assumptions used in the studies. Toxicological values for a given compound will be derived from the study summarizing the best available data or the set of data that exhibits either the lowest value for Lowest-Observed-Adverse-Effect-Level (LOAEL) or the highest No-Observed-Adverse-Effect-Level (NOAEL). The LOAEL is the lowest dosage at which some effect is shown. The NOAEL is the dosage at which no observed effect or response is noted. Derivation of the acceptable daily intake will incorporate uncertainty factors for: extrapolation of data from animals to humans, calculation of the human-equivalent dose, and interspecies variability in sensitivity of the toxicant.
- 4) If a toxicological profile from ATSDR is not available, toxicity data will be obtained in a literature search of US EPA sources in the following order:
  - a) Health Assessment Documents,
  - b) Health Effects Assessments,
  - c) Health Advisories, and
  - d) Registry of Toxic Effects of Chemical Substances (RTECS) and Hazardous Substances Data Bank (HSDB).
- 5) If the above sources cannot provide data, Toxline and other related databases and journals will be searched for relevant dose-response studies from which to derive toxicity values; sound principles of toxicology will be used.
- 6) If toxicity data do not exist in any of the above sources, LD<sub>50</sub> data for a given compound will be compiled. The lowest oral LD<sub>50</sub> value for any species will be



- divided by appropriate safety factors, depending upon the anticipated length of exposure.
- 7) For chemicals that lack any toxicity information, the concept of structure-activity relationships will be applied. This concept allows the derivation of an acceptable intake for a chemical by inference and analogy to closely related compounds.

Toxicity indices will be provided for site-related COPCs in order that potential risks can be evaluated quantitatively with the objective of minimizing inherent uncertainty in the process. The toxicity values utilized in the human health risk assessment will be presented in a tabular format.

#### **2.4 Risk Characterization**

The objective of the risk characterization is to combine the results of the exposure and toxicity assessments in a quantitative evaluation of risk. The estimated intakes calculated for each exposure pathway considered and each COPC will be compared to RfDs for noncarcinogenic effects. RfDs have been developed by the US EPA based on the most sensitive noncarcinogenic effects. The chronic RfD for a chemical is an estimate of a lifetime daily exposure level for the human population (including sensitive subpopulations) that is likely to be without an appreciable risk of deleterious effects. A subchronic RfD is similar; however, it is based on less-than-a-lifetime time period (e.g., less than seven years for humans). The potential for non-cancer health effects will be evaluated by comparing an exposure level (intake) over a specified time period with the RfD for a similar exposure period (subchronic or chronic). This ratio of exposure to toxicity is called the hazard quotient.

The non-cancer hazard quotient assumes that there is a threshold level of exposure (i.e., RfD), below which it is unlikely for even the most sensitive populations to experience adverse health effects. If the exposure level exceeds the threshold (i.e., the hazard quotient exceeds a value greater than 1.0), there may be concern for potential non-cancer effects (viz., the greater the value of the hazard quotient above unity, the greater the level of concern for potential health impacts).



To assess the overall potential for non-cancer effects posed by multiple chemicals, a hazard index will be derived by summing the individual hazard quotients for a given receptor. This approach assumes additivity of critical effects of multiple chemicals. This is appropriate only for compounds that induce the same effect by the same mechanism of action. This conservative approach will significantly overestimate the actual potential for adverse health impacts.

In cancer risk assessment, US EPA has required the use of the upper limit, which produces an estimate of risk that has a 95 percent probability of exceeding the actual risk, which may, in fact, be zero. Lifetime average daily intakes for carcinogenic chemicals will be multiplied by the chemical-specific CSF to determine the probability that an individual may experience carcinogenic effects under site-specific conditions. For exposures to multiple carcinogens, the upper limits of cancer risks for each chemical in each exposure pathway will be summed to derive a total cancer risk. The US EPA recognizes that it is not technically appropriate to sum upper confidence limits of the risk to produce a realistic total probability, but requires this approach be used.

Summaries of the combined hazard indices and lifetime cancer risks resulting from exposure to site-related chemicals for the populations of potential concern will be presented in both textual and tabular formats. A description of each exposure pathway and the resulting hazard index and cancer risk will be provided. Hazard index and cancer risk estimates will also be compared to MDEQ or US EPA benchmark values to determine compliance with regulations.

## **2.5      Uncertainty Analysis**

The US EPA has consistently recommended that all risk assessment activities include some degree of uncertainty analysis to provide proper perspective to risk management decision-makers. The uncertainty analysis section of the report will discuss the sources of uncertainty throughout each step of the risk assessment process and how these uncertainties may have over- or underestimated the final hazard and risk calculations.



### **3.0 Development of Site-Specific Cleanup Standards**

In the event that site-specific hazard or risk estimates exceed acceptable MDEQ or US EPA levels, site-specific, health-based cleanup goals will be calculated. The same exposure assumptions used in the baseline risk assessment will be used to calculate site-specific cleanup standards. However, MDEQ or US EPA target hazard and risk benchmarks will also be included in the calculations to ensure adequate protection of human health. The risk assessment process, as described above, will be utilized to determine remediation requirements and to document that achieving these cleanup objectives will afford adequate protection to human health and the environment.

### **4.0 Project Execution**

#### **4.1 Project Team**

The primary risk assessment project team for the Hattiesburg site will consist of Kenneth G. Symms, Ph.D., DABT, Robert J. Fares, and Kathleen A. Koerber, representatives of the Environmental Standards Risk Assessment Department. Environmental Standards background information, risk assessment qualifications, and experience with wood-preserving facilities are provided in Appendix A. Appendix B presents professional profiles for the above-listed individuals. Other Environmental Standards personnel may be called upon to provide expertise (e.g., chemists, geologists) or additional help if needed. Their involvement, however, will be minimal.

#### **4.2 Project Schedule**

A Risk Assessment report will be submitted to MDEQ for its review within approximately 16 weeks subsequent to MDEQ's approval of this Work Plan. A timeline depicting the project schedule is presented in Figure 1.



#### **4.3 Site Visit**

It will be necessary for Environmental Standards risk assessment personnel to visit the Hattiesburg site. This site visit will provide Environmental Standards with a better understanding of the site geography and setting for the development of a site conceptual model and will greatly aid in evaluating site conditions that may affect critical decisions during the risk assessment process. The personnel visiting the site will include Dr. Kenneth G. Symms, Ph.D., DABT, Mr. Robert J. Fares, and Ms. Kathleen A. Koerber. A one-day (eight-hour) visit should provide the time necessary for these personnel to familiarize themselves with the site and surroundings.

#### **4.4 Reporting**

Three copies of the Risk Assessment report will be provided to MDEQ. The report will be bound in a three-ring binder and will include a table of contents, acronym list, bibliography, and relevant tables and figures (as necessary). The Table of Contents for the risk assessment report will be consistent with MDEQ guidance. The example presented below summarizes the salient information that will be included in the Risk Assessment report:

- 1.0 Introduction and General Methodology
- 2.0 Data Evaluation
  - 2.1 COPC Selection Process
- 3.0 Exposure Assessment
  - 3.1 General Intake Equation
  - 3.2 Exposure Parameters
    - 3.2.1 General Exposure Parameters
    - 3.2.2 Route-Specific Exposure Parameters
  - 3.3 Indirect Exposure Paradigms
- 4.0 Toxicity Assessment

**5.0 Risk Characterization**

- 5.1 Non-Carcinogenic Risk Estimates**
- 5.2 Carcinogenic Risk Estimates**

**6.0 Uncertainty Analysis**

If the development of site-specific cleanup standards is necessary, a discussion of the methods utilized will be added as Section 7.0. As recommended in US EPA guidance (Risk Assessment Guidance for Superfund, Part A, 1989), tables will be provided for the following:

- Statistical analysis of site data and COPC selection process for each media evaluated;
- Exposure pathway evaluation;
- Exposure parameter assumptions;
- Toxicity indices;
- Hazard and risk calculation summaries; and
- Detailed daily intake calculations and hazard and risk estimates for each receptor and pathway.

Other tables may be included as the need arises (e.g., detailing local meteorological data or site-specific cleanup standard calculations).



**Figure 1** Former Gulf States Creosoting Site Risk Assessment Schedule

\*Remainder of schedule contingent upon MDEQ approval of the Work Plan in the time provided.