

APPENDIX B

TABLE B - 1
 SAMPLE LOCATIONS AND RATIONALE
 GULF STATE CREOSOTE SITE
 SI - PHASE II

Collection Date	Sample Code	Sample Type	Locations	Rationale	Depth (ft bls)
10/15/91	GS-PW-01	Potable well	Public Well - 1.5 miles east of site - Hattiesburg Water Department	Determine presence or absence of contamination	400 feet
10/16/91	GS-TW-01	Groundwater	Upgradient Temporary Well - Adjacent to Pine Street and Ryan Motors	Background	30 feet
10/16/91	GS-SB-01	Subsurface Soil	Adjacent to Pine Street and Ryan Motors	Background	8 feet
10/16/91	GS-SD-01	Sediment	Collected from Gordons Creek adjacent to Trailer Park	Upstream/Background	NA
10/16/91	GS-SD-02	Sediment	Collected from Gordons Creek approximately 60 feet north/northeast of eastern drainage ditch	Determine presence or absence of contamination	NA
10/17/91	GS-TW-02	Groundwater	Downgradient temporary well located between the two drainage ditches near the Trailer Park	Determine presence or absence of contamination	25 feet
10/17/91	GS-SB-02	Subsurface	Located between the two drainage ditches near the Trailer Park	Determine presence or absence of contamination.	7 feet
	GS-TB-01	Trip Blank	Collected Prior to entry of site	QA/QC	NA

NA - Not Applicable



MISSISSIPPI DEPARTMENT
OF ENVIRONMENTAL QUALITY

CHAIN OF CUSTODY RECORD

OFFICE OF
POLLUTION CONTROL
P. O. Box 10385
Jackson, Mississippi 39289-0385

MSD 985967199				PROJECT LEADER Jim Harbage / Michael Slack		REMARKS * METALS ANALYSIS - SEE ATTACHED PAGE	
PROJECT NAME/LOCATION GULF STATE CREOSOTE / HATTIESBURG, MS				SAMPLER MICHAEL SLACK <i>Michael Slack</i> KEN WHITTEN <i>Ken Whitten</i> JIM HARBAGE <i>Jim Harbage</i> MARK WALTERS <i>Mark E. Walters</i>		DATA TO: MICHAEL SLACK / JIM HARBAGE (MS DEQ)	
ESD SAMPLE TYPES				SAMPLER		ANALYSIS	
1. SURFACE WATER 2. GROUND WATER 3. POTABLE WATER 4. WASTEWATER 5. LEACHATE 11. OTHER TRIP BLANK				6. SOIL/SEDIMENT 7. SLUDGE 8. WASTE 9. AIR 10. FISH		CIRCLE/ADD parameters desired. List no. of containers submitted.	
STATION NO.	SAMPLE TYPE	DATE	TIME	COMP	GRAB	TOTAL CONTAINERS	TAG NO./REMARKS
GS-PW-01	3	10/15/91	2:50 PM	✓		7	2 2 1 1 1 PUBLIC WELL - HALL STREET (WATER DEPT.) SHALLOWEST WELL - WELL #1
GS-TW-01	2	10/16/91	11:50 AM	✓		7	2 2 1 1 1 BACKGROUND TEMPORARY WELL - PINE STREET & RYAN MOTORS
GS-SB-01	6	10/16/91	3:50	✓		4	1 1 1 1 ① BACKGROUND SUBSURFACE SOIL - PINE STREET & RYAN MOTORS
GS-SD-01	6	10/16/91	5:00 PM	✓		4	1 1 1 1 ① BACKGROUND GORDONS CREEK SEDIMENT - TRAILER PARK
GS-SD-02	6	10/16/91	5:20 PM	✓		4	1 1 1 1 ① DOWNGRADIENT GORDONS CREEK SEDIMENT - EAST OF DITCH
GS-TW-02	2	10/17/91	10:05 AM	✓		7	2 2 1 1 1 DOWNGRADIENT TEMPORARY WELL - NEAR TRAILER PARK
GS-SB-02	6	10/17/91	11:30 AM	✓		4	1 1 1 1 ① SUBSURFACE SOIL - ON-SITE BETWEEN THE TWO DITCHES
TRIP BLANK	11					2	2 TRIP BLANK FROM MSU LAB
							UNDER ANALYSIS: ① - DESIGNATES 1 CONTAINER FOR METALS/CYANIDE ANALYSIS 1 - DESIGNATES 1 CONTAINER 2 - DESIGNATES 2 CONTAINERS

RELINQUISHED BY: (PRINT) <i>Ken Whitten</i>	DATE/TIME: 10/18/91	RECEIVED BY: (PRINT) <i>Mark Wooster</i>	RELINQUISHED BY: (PRINT)	DATE/TIME: 10/18/91	RECEIVED BY: (PRINT) <i>D.A. Dollar</i>
(SIGN) <i>Ken Whitten</i>	11:33 AM	(SIGN) <i>Mark Wooster</i>	(SIGN)	6:54 PM	(SIGN) <i>D.A. Dollar</i>
RELINQUISHED BY: (PRINT)	DATE/TIME:	RECEIVED BY: (PRINT)	RELINQUISHED BY: (PRINT)	DATE/TIME:	RECEIVED BY: (PRINT)
(SIGN)		(SIGN)	(SIGN)		(SIGN)

DISTRIBUTION: White and Yellow copies accompany sample shipment to laboratory. Yellow copy retained by laboratory. White copy is returned to samplers; Pink copy retained by samplers.

NOT TO SCALE

SD = SEDIMENT
SB - SUBSURFACE

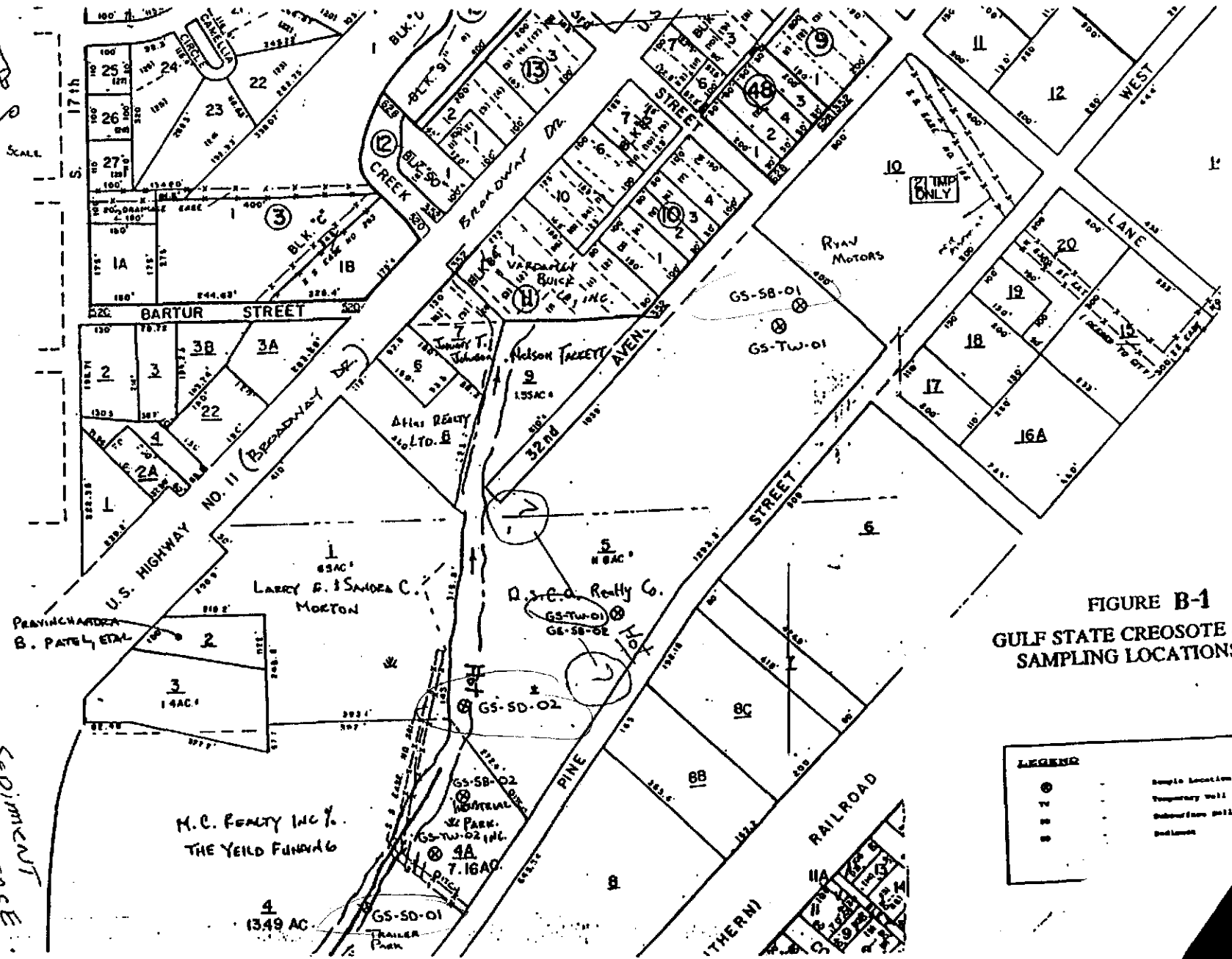
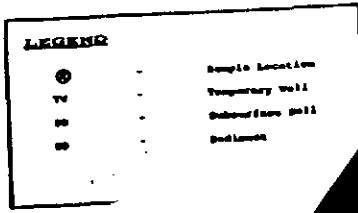
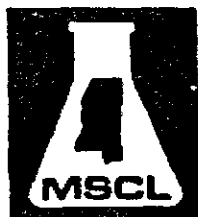


FIGURE B-1
GULF STATE CREOSOTE SITE
SAMPLING LOCATIONS



MISSISSIPPI STATE UNIVERSITY

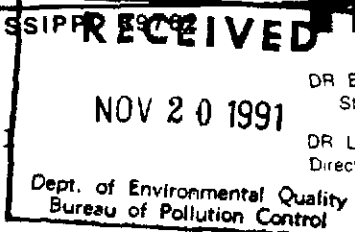


MISSISSIPPI
STATE CHEMICAL LABORATORY

BOX CR - MISSISSIPPI STATE, MISSISSIPPI 39262

TELEPHONE (601) 325-3324

November 18, 1991



DR EARL G. ALLEY
State Chemist
DR LARRY G. LANE
Director, IAS Division

Analysis No. 826,332-335

Analysis of Soil/Sediment

Received on 10-21-91

Address P.O. Box 10385 Jackson, MS 39209

Marked: Gulf State Creosote,
Hattiesburg
from MS Office of Pollution Control
ATTN: Jim Hardage

RESULTS:

<u>MSCL No.</u>	<u>MS DEQ-OPC Identification</u>
826,332	GS-SB-01, Background Subsurface Soil, Pine St. & Ryan Motors
826,333	GS-SD-01, Background Sediment, Gordons Creek Trailer Park
826,334	GS-SD-02, Downgradient Sediment, Gordons Creek East of Ditch
826,335	GS-SB-02, Subsurface Soil, Between two ditches

Results from our gc/mass spec analyses of the above sediment samples for Semivolatile Organic Compounds on the Target Compound List are presented in attached reports.

Analytical Costs

4 ABNs By gc/ms @ \$400 = \$1600

Earl G. Alley

State Chemist

PLEASE GIVE NUMBER WHEN REFERRING TO THIS ANALYSIS

TARGET COMPOUND LIST

SEMIVOLATILES DATA SHEET FOR SOILS

MSCL ANALYSIS NO. 826,332

MARKED Gulf State Creosote

ANALYSIS OF Soil

GS-SB-01

COMPOUNDS	MQL*	Micro g/Kg	COMPOUNDS	MQL*	Micro g/Kg
Phenol	330	ND	2,4-Dinitrophenol	1600	ND
bis(2-Chloroethyl)ether	330	ND	4-Nitrophenol	1600	ND
2-Chlorophenol	330	ND	Dibenzofuran	330	ND
1,3-Dichlorobenzene	330	ND	2,4-Dinitrotoluene	330	ND
1,4-Dichlorobenzene	330	ND	Diethylphthalate	330	ND
Benzyl alcohol	330	ND	4-Chlorophenyl-phenyl ether	330	ND
1,2-Dichlorobenzene	330	ND	Fluorene	330	ND
2-Methylphenol	330	ND	4-Nitroaniline	1600	ND
bis(2-Chloroisopropyl)ether	330	ND	4,6-Dinitro-2-methylphenol	1600	ND
4-Methylphenol	330	ND	N-nitrosodiphenylamine	330	ND
N-Nitroso-di-n-dipropylamine	330	ND	4-Bromophenyl-phenylether	330	ND
Hexachloroethane	330	ND	Hexachlorobenzene	330	ND
Nitrobenzene	330	ND	Pentachlorophenol	1600	ND
Isophorone	330	ND	Phenanthrene	330	ND
2-Nitrophenol	330	ND	Anthracene	330	ND
2,4-Dimethylphenol	330	ND	Di-n-butylphthalate	330	ND
Benzoic acid	1600	ND	Fluoranthene	330	ND
bis(2-Chloroethoxy)methane	330	ND	Pyrene	330	ND
2,4-Dichlorophenol	330	ND	Butylbenzylphthalate	330	ND
1,2,4-Trichlorobenzene	330	ND	3,3'-Dichlorobenzidine	660	ND
Naphthalene	330	ND	Benzo(a)anthracene	330	ND
4-Chloroaniline	330	ND	Chrysene	330	ND
Hexachlorobutadiene	330	ND	bis(2-Ethylhexyl)phthalate	330	ND
4-Chloro-3-methylphenol	330	ND	Di-n-octylphthalate	330	NT
2-Methylnaphthalene	330	ND	Benzo(b)fluoranthene	330	NT
Hexachlorocyclopentadiene	330	ND	Benzo(k)fluoranthene	330	NT
2,4,6-Trichlorophenol	330	ND	Benzo(a)pyrene	330	NT
2,4,5-Trichlorophenol	1600	ND	Indeno(1,2,3-cd)pyrene	330	ND
2-Chloronaphthalene	330	ND	Dibenz(a,h)anthracene	330	ND
2-Nitroaniline	1600	ND	Benzo(g,h,i)perylene	330	ND
Dimethylphthalate	330	ND			
Acenaphthylene	330	ND			
2,6-Dinitrotoluene	330	ND			
3-Nitroaniline	1600	ND			
Acenaphthene	330	ND			

*ND = None Detected

MQL = Minimum Quantifiable Level

SURROGATES	RECOVERY (%)
2-Fluorophenol	100
Phenol-d5	110
Nitrobenzene-d5	88
2-Fluorobiphenyl	104
2,4,6-Tribromophenol	107
p-Terphenyl-d14	119

Multiply MQL's by _____

- No peaks above 40% of internal standard were observed.
- Peaks above 40% of internal standard were not identified.
- Peaks above 40% internal standard not on EPA Appendix IX.
- Additional peaks were observed but not examined.

Earl L. Colley
State Chemist

TARGET COMPOUND LIST

SEMIVOLATILES DATA SHEET FOR SOILS

MSCL ANALYSIS NO. 826,333

MARKED Gulf State Creosote

ANALYSIS OF Soil

GS-SD-01

COMPOUNDS

	MQL*	Micro g/Kg
Phenol	330	ND
bis(2-Chloroethyl)ether	330	ND
2-Chlorophenol	330	ND
1,3-Dichlorobenzene	330	ND
1,4-Dichlorobenzene	330	ND
Benzyl alcohol	330	ND
1,2-Dichlorobenzene	330	ND
2-Methylphene	330	ND
bis(2-Chloroisopropyl)ether	330	ND
4-Methylphenol	330	ND
N-Nitroso-di-n-dipropylamine	330	ND
Hexachloroethane	330	ND
Nitrobenzene	330	ND
Isophorone	330	ND
2-Nitrophenol	330	ND
2,4-Dimethylphenol	330	ND
Benzoic acid	1600	ND
bis(2-Chloroethoxy)methane	330	ND
2,4-Dichlorophenol	330	ND
1,2,4-Trichlorobenzene	330	ND
Naphthalene	330	ND
4-Chloroaniline	330	ND
Hexachlorobutadiene	330	ND
4-Chloro-3-methylphenol	330	ND
2-Methylnaphthalene	330	ND
Hexachlorocyclopentadiene	330	ND
2,4,6-Trichlorophenol	330	ND
2,4,5-Trichlorophenol	1600	ND
2-Chloronaphthalene	330	ND
2-Nitroaniline	1600	ND
Dimethylphthalate	330	ND
Acenaphthylene	330	ND
2,6-Dinitrotoluene	330	ND
3-Nitroaniline	1600	ND
Acenaphthene	330	ND

COMPOUNDS

	MQL*	Micro g/Kg
2,4-Dinitrophenol	1600	ND
4-Nitrophenol	1600	ND
Dibenzofuran	330	ND
2,4-Dinitrotoluene	330	ND
Diethylphthalate	330	ND
4-Chlorophenyl-phenyl ether	330	ND
Fluorene	330	ND
4-Nitroaniline	1600	ND
4,6-Dinitro-2-methylphenol	1600	ND
N-nitrosodiphenylamine	330	ND
4-Bromophenyl-phenylether	330	ND
Hexachlorobenzene	330	ND
Pentachlorophenol	1600	ND
Phenanthrene	330	470
Anthracene	330	ND
Di-n-butylphthalate	330	ND
Fluoranthene	330	700
Pyrene	330	470
Butylbenzylphthalate	330	ND
3,3'-Dichlorobenzidine	660	ND
Benzo(a)anthracene	330	Trace
Chrysene	330	Trace
bis(2-Ethylhexyl)phthalate	330	ND
Di-n-octylphthalate	330	ND
Benzo(b)fluoranthene	330	ND
Benzo(k)fluoranthene	330	ND
Benzo(a)pyrene	330	ND
Indeno(1,2,3-cd)pyrene	330	ND
Dibenz(a,h)anthracene	330	ND
Benzo(g,h,i)perylene	330	ND

*ND = None Detected

MQL = Minimum Quantifiable Level

SURROGATES	RECOVERY (%)
2-Fluorophenol	72
Phenol-d5	92
Nitrobenzene-d5	43
2-Fluorobipheny	78
2,4,6-Tribromophenol	88
p-Terphenyl-d14	92

Multiply MQL's by _____

- _____ No peaks above 40% of internal standard were observed.
- 1 Peaks above 40% of internal standard were not identified.
- 1 Peaks above 40% internal standard not on EPA Appendix IX.
- _____ Additional peaks were observed but not examined.

Earl R. Colley

 State Chemist/

TARGET COMPOUND LIST

SEMIVOLATILES DATA SHEET FOR SOILS

MSCL ANALYSIS NO. 826,335

MARKED Gulf State Creosote

ANALYSIS OF Soil

GS-SB-02

COMPOUNDS

	MQL*	Micro g/Kg
Phenol	330	ND
bis(2-Chloroethyl)ether	330	ND
2-Chlorophenol	330	ND
1,3-Dichlorobenzene	330	ND
1,4-Dichlorobenzene	330	ND
Benzyl alcohol	330	ND
1,2-Dichlorobenzene	330	ND
2-Methylphene	330	ND
bis(2-Chloroisopropyl)ether	330	ND
4-Methylphenol	330	ND
N-Nitroso-di-n-dipropylamine	330	ND
Hexachloroethane	330	ND
Nitrobenzene	330	ND
Isophorone	330	ND
2-Nitrophenol	330	ND
2,4-Dimethylphenol	330	ND
Benzoic acid	1600	ND
bis(2-Chloroethoxy)methane	330	ND
2,4-Dichlorophenol	330	ND
1,2,4-Trichlorobenzene	330	ND
Naphthalene	330	1,900,000
4-Chloroaniline	330	ND
Hexachlorobutadiene	330	ND
4-Chloro-3-methylphenol	330	ND
2-Methylnaphthalene	330	1,400,000
Hexachlorocyclopentadiene	330	ND
2,4,6-Trichlorophenol	330	ND
2,4,5-Trichlorophenol	1600	ND
2-Chloronaphthalene	330	ND
2-Nitroaniline	1600	ND
Dimethylphthalate	330	ND
Acenaphthylene	330	Trace
2,6-Dinitrotoluene	330	ND
3-Nitroaniline	1600	ND
Acenaphthene	330	970,000

COMPOUNDS

	MQL*	Micro g/Kg
2,4-Dinitrophenol	1600	ND
4-Nitrophenol	1600	ND
Dibenzofuran	330	1,000,000
2,4-Dinitrotoluene	330	ND
Diethylphthalate	330	ND
4-Chlorophenyl-phenyl ether	330	ND
Fluorene	330	1,500,000
4-Nitroaniline	1600	ND
4,6-Dinitro-2-methylphenol	1600	ND
N-nitrosodiphenylamine	330	ND
4-Bromophenyl-phenylether	330	ND
Hexachlorobenzene	330	ND
Pentachlorophenol	1600	ND
Phenanthrene *	330	3,500,000
Anthracene *	330	4,200,000
Di-n-butylphthalate	330	ND
Fluoranthene	330	1,600,000
Pyrene	330	770,000
Butylbenzylphthalate	330	ND
3,3'-Dichlorobenzidine	660	ND
Benzo(a)anthracene	330	270,000
Chrysene	330	280,000
bis(2-Ethylhexyl)phthalate	330	ND
Di-n-octylphthalate	330	ND
Benzo(b)fluoranthene	330	113,000
Benzo(k)fluoranthene	330	100,000
Benzo(a)pyrene	330	85,000
Indeno(1,2,3-cd)pyrene	330	ND
Dibenz(a,h)anthracene	330	ND
Benzo(g,h,i)perylene	330	ND

*ND = None Detected

MQL = Minimum Quantifiable Level
*Estimated Value

Multiply MQL's by 250

SURROGATES	RECOVERY (%)
2-Fluorophenol	
Phenol-d5	
Nitrobenzene-d5	
2-Fluorobipheny	
2,4,6-Tribromophenol	
p-Terphenyl-d14	

NA= Had to dilute out of Linear Range

No peaks above 40% of internal standard were observed.

1 Peaks above 40% of internal standard were not identified.

19 Peaks above 40% internal standard, peaks appear to be PAHS not on EPA Appendix IX.

X Additional peaks were observed but not examined.

Carl L. Alley
State Chemist

SEMIVOLATILES DATA SHEET FOR SOILS

MSCL ANALYSIS NO. 826,334

MARKED Gulf State Creosote

ANALYSIS OF Soils

GS-SD-02

COMPOUNDS

	MLQ*	Micro g/Kg
Phenol	330	ND
bis(2-Chloroethyl)ether	330	ND
2-Chlorophenol	330	ND
1,3-Dichlorobenzene	330	ND
1,4-Dichlorobenzene	330	ND
Benzyl alcohol	330	ND
1,2-Dichlorobenzene	330	ND
2-Methylphene	330	ND
bis(2-Chloroisopropyl)ether	330	ND
4-Methylphenol	330	ND
N-Nitroso-di-n-dipropylamine	330	ND
Hexachloroethane	330	ND
Nitrobenzene	330	ND
Isophorone	330	ND
2-Nitrophenol	330	ND
2,4-Dimethylphenol	330	ND
Benzoic acid	1600	ND
bis(2-Chloroethoxy)methane	330	ND
2,4-Dichlorophenol	330	ND
1,2,4-Trichlorobenzene	330	ND
Naphthalene	330	240,000
4-Chloroaniline	330	ND
Hexachlorobutadiene	330	ND
4-Chloro-3-methylphenol	330	ND
2-Methylnaphthalene	330	240,000
Hexachlorocyclopentadiene	330	ND
2,4,6-Trichlorophenol	330	ND
2,4,5-Trichlorophenol	1600	ND
2-Chloronaphthalene	330	ND
2-Nitroaniline	1600	ND
Dimethylphthalate	330	ND
Acenaphthylene	330	Trace
2,6-Dinitrotoluene	330	ND
3-Nitroaniline	1600	ND
Acenaphthene	330	370,000

COMPOUNDS

	MLQ*	Micro g/Kg
2,4-Dinitrophenol	1600	ND
4-Nitrophenol	1600	ND
Dibenzofuran	330	422,000
2,4-Dinitrotoluene	330	ND
Diethylphthalate	330	ND
4-Chlorophenyl-phenyl ether	330	ND
Fluorene	330	554,000
4-Nitroaniline	1600	ND
4,6-Dinitro-2-methylphenol	1600	ND
N-nitrosodiphenylamine	330	ND
4-Bromophenyl-phenylether	330	ND
Hexachlorobenzene	330	ND
Pentachlorophenol	1600	ND
Phenanthrene *	330	18,000,000
Anthracene	330	222,000
Di-n-butylphthalate	330	ND
Fluoranthene	330	772,000
Pyrene	330	490,000
Butylbenzylphthalate	330	ND
3,3'-Dichlorobenzidine	660	ND
Benzo(a)anthracene	330	170,000
Chrysene	330	150,000
bis(2-Ethylhexyl)phthalate	330	NT
Di-n-octylphthalate	330	NT
Benzo(b)fluoranthene	330	32,000
Benzo(k)fluoranthene	330	72,000
Benzo(a)pyrene	330	67,000
Indeno(1,2,3-cd)pyrene	330	Trace
Dibenz(a,h)anthracene	330	ND
Benzo(g,h,i)perylene	330	Trace

*ND = None Detected

MLQ = Minimum Quantifiable Level
*Estimated value

Multiply MLQ's by 100

SURROGATES	RECOVERY (%)
2-Fluorophenol	
Phenol-d5	
Nitrobenzene-d5	
2-Fluorobiphenyl	
2,4,6-Tribromophenol	
p-Terphenyl-d14	

NA= Had to dilute out of Linear Range

No peaks above 40% of internal standard were observed.

1 Peaks above 40% of internal standard were not identified.

19 Peaks above 40% internal standard. compounds appear to be PAHs not on EPA Appendix IX.

X Additional peaks were observed but not examined.

PAHs = Polynuclear Aromatic Hydrocarbons

Carl L. Colley

State Chemist

TARGET COMPOUND LIST
SEMIVOLATILES DATA SHEET FOR WATERS

RECEIVED
JAN 15 1992
Dept. of Environmental Quality
Bureau of Pollution Control

SCL ANALYSIS NO. 826,336

MARKED Gulf State

ANALYSIS OF Water

GS-PW-01

COMPOUNDS

COMPOUNDS

	MLQ*	Micro g/L		MLQ*	Micro g/L
Phenol	10	ND	2,4-Dinitrophenol	50	ND
bis(2-Chloroethyl)ether	10	ND	4-Nitrophenol	50	ND
2-Chlorophenol	10	ND	Dibenzofuran	10	ND
1,3-Dichlorobenzene	10	ND	2,4-Dinitrotoluene	10	ND
1,4-Dichlorobenzene	10	ND	Diethylphthalate	10	ND
Benzyl alcohol	20	ND	4-Chlorophenyl-phenyl ether	10	ND
1,2-Dichlorobenzene	10	ND	Fluorene	10	ND
2-Methylphenol	10	ND	4-Nitroaniline	50	ND
bis(2-Chloroisopropyl)ether	10	ND	4,6-Dinitro-2-methylphenol	50	ND
4-Methylphenol	10	ND	N-nitrosodiphenylamine	10	ND
N-Nitroso-di-n-dipropylamine	10	ND	4-Bromophenyl-phenylether	10	ND
Hexachloroethane	10	ND	Hexachlorobenzene	10	ND
Nitrobenzene	10	ND	Pentachlorophenol	50	ND
Isophorone	10	ND	Phenanthrene	10	ND
2-Nitrophenol	10	ND	Anthracene	10	ND
2,4-Dimethylphenol	10	ND	Di-n-butylphthalate	10	ND
Benzoic acid	20	ND	Fluoranthene	10	ND
bis(2-Chloroethoxy)methane	10	ND	Pyrene	10	ND
2,4-Dichlorophenol	10	ND	Butylbenzylphthalate	10	ND
1,2,4-Trichlorobenzene	10	ND	3,3'-Dichlorobenzidine	20	ND
Naphthalene	10	ND	Benzo(a)anthracene	10	ND
4-Chloroaniline	20	ND	Chrysene	10	ND
Hexachlorobutadiene	10	ND	bis(2-Ethylhexyl)phthalate	10	ND
4-Chloro-3-methylphenol	20	ND	Di-n-octylphthalate	10	ND
2-Methylnaphthalene	10	ND	Benzo(b)fluoranthene	10	ND
Hexachlorocyclopentadiene	10	ND	Benzo(k)fluoranthene	10	ND
2,4,6-Trichlorophenol	10	ND	Benzo(a)pyrene	10	ND
2,4,5-Trichlorophenol	10	ND	Indeno(1,2,3-cd)pyrene	10	ND
2-Chloronaphthalene	10	ND	Dibenz(a,h)anthracene	10	ND
2-Nitroaniline	50	ND	Benzo(g,h,i)perylene	10	ND
Dimethylphthalate	10	ND			
Acenaphthylene	10	ND			
2,6-Dinitrotoluene	10	ND			
3-Nitroaniline	50	ND			
Acenaphthene	10	ND			

*ND = None Detected

MLQ = Minimum Quantifiable Level

SURROGATES	RECOVERY (%)
2-Fluorophenol	53
Phenol-d5	33
Nitrobenzene-d5	99
2-Fluorobiphenyl	98
2,4,6-Tribromophenol	115
p-Terphenyl-d14	104

Multiply MLQ's by _____

____ No peaks above 40% of internal standard were observed.

4 Peaks above 40% of internal standard were not identified.

2 Peaks above 40% internal standard not on EPA Appendix IX. Appear to be substituted chlorinated benzenes at an estimated total concentration of 10 µg/L.

Carl L. Colley
State Chemist