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**U.S. EPA Region III
Office of Analytical Services
and Quality Assurance
Annapolis, Maryland**

ANALYTICAL REPORT

**WILMINGTON DRUM
SUPERFUND REMOVAL/REMEDIAL
Lab Request # REQ99016**

December 15, 1998

AR100005

U.S. EPA Region III
Office of Analytical Services
and Quality Assurance
Annapolis, Maryland

December 15, 1998

ANALYTICAL RESULTS: WILMINGTON DRUM [REQ99016]

Dear William C. Wentworth (3HS33),

Enclosed is our analytical report for the above case. It is organized into several sections: Analytical Request and Sample Descriptions, Organic, Inorganic, and Microbiological Results. All data were peer reviewed.

Analytical Request and Sample Descriptions: (General)

Each laboratory assigned number, station, description, matrix, sample date and locational data is reported. A table summarizes the tests assigned to each sample. A glossary and qualifier code definition is provided.

Inorganic Results:

For requests assigned inorganic tests, results are grouped by service group, e.g., Metals. Sample results are reported; non-detects are provided with the actual quantitation limit. Method description and quality control protocols are described in analyst narratives.

Organic Results:

For requests assigned organic tests, results are grouped by service group, e.g., Volatile Organic Compounds. Only detected analytes are reported. Nominal Quantitation Limit (NQL) tables are provided for each service group. Specific information for the calculation of Actual Quantitation Limits (AQL) achieved for a given sample is included. Quality control values are provided in summary tables with acceptance criteria. Method description and quality control protocols are described in analyst narratives.

Microbiological Results:

For requests assigned microbiological tests, sample results and quality control values are incorporated into a single table. Method description and quality control protocols are described in analyst narratives.

If you have any questions we may be reached at 410-573-2600.

Approval for Release:

Susan Warner

Site Contact(s): Kevin Burns (WESTON)
William C. Wentworth (3HS33)

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Facility: WILMINGTON DRUM

Program: SUPERFUND REMOVAL/REMEDIAL

Lab Request #: REQ99016

SAMPLE DESCRIPTIONS

Sample #	Station Description	Matrix	Type	End Collection	
				Date	Time
98102101	STA SD-01, SD-01	Soil	GRAB	10/19/98	15:40
98102102	STA SD-02, SD-02	Soil	GRAB	10/19/98	11:00
98102103	STA SD-03, SD-03	Soil	GRAB	10/19/98	11:30
98102104	STA SD-06, SD-06	Soil	GRAB	10/19/98	13:00
98102105	STA SD-05, SD-05	Soil	GRAB	10/19/98	10:00
98102106	STA SD-04, SD-04	Soil	GRAB	10/19/98	12:30
98102107	STA TB-1, TB-1	Aqueous Blank	GRAB	10/19/98	08:00
98102108	STA PB-1, PB-1	Aqueous Blank	GRAB	10/19/98	12:00
98102109	STA SW-04, SW-04	Water - Type Unspecified	GRAB	10/19/98	12:30
98102110	STA SW-06, SW-06	Water - Type Unspecified	GRAB	10/19/98	13:00
98102111	STA SW-01, SW-01	Water - Type Unspecified	GRAB	10/19/98	15:40
98102112	STA SW-02, SW-02	Water - Type Unspecified	GRAB	10/19/98	11:00
98102113	STA SW-03, SW-03	Water - Type Unspecified	GRAB	10/19/98	11:30
98102114	STA SW-05, SW-05	Water - Type Unspecified	GRAB	10/19/98	10:00

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TESTS REQUESTED

Inorganic Test Assigned	Sample No. 981021-													
	01	02	03	04	05	06	07	08	09	10	11	12	13	14
Metals Analysis	X	X	X	X	X	X		X	X	X	X	X	X	X
Percent Dry Weight	X	X	X	X	X	X								
Total Cyanide	X	X	X	X	X	X		X	X	X	X	X	X	X

Organic Test Assigned	Sample No. 981021-													
	01	02	03	04	05	06	07	08	09	10	11	12	13	14
PCBs and Pesticides by Gas Chromatography	X	X	X	X	X	X		X	X	X	X	X	X	X
Semivolatile Organics by GC/MS	X	X	X	X	X	X		X	X	X	X	X	X	X
Volatile Organics by Purge & Trap GC/MS	X	X	X	X	X	X		X	X	X	X	X	X	X

QUALIFIER CODE AND GLOSSARY DEFINITIONS

Qualifier Codes:

<	Sample value is below the quantitation limit. Quantitation limit reported.
<=	Reported value is estimated. Sample was analyzed in duplicate, one value is equal to or above the quantitation limit and one below. Average of quantitation limit and detected value reported.
>	Sample value is above the quantitation range.
A	Quality control value is outside acceptance limits.
B	Not detected substantially above (10 times) the level reported in the laboratory or field blanks (includes field, trip, rinsate, and equipment blanks).
C	See report narrative for analyst's observations concerning this result.
D	Sample and duplicate values are below the quantitation limit. Quantitation limit reported.
E	Value exceeds a theoretically greater value (e.g. dissolved > total, orthophosphate > total phosphorus). However, the difference is within the expected precision of the analytical techniques and is not statistically significant.
I	An interference exists which masks true response. See report narrative for explanation.
J	Analyte present. Reported value is estimated, concentration is outside the range for accurate quantitation.
K	Analyte present. Reported value may be biased high. Actual value is expected to be lower.
L	Analyte present. Reported value may be biased low. Actual value is expected to be higher.
N	Presumptive evidence indicates the presence of the compound. Special methods and/or method modifications may be needed to confirm its presence or absence in future sampling efforts.
NA	Analysis was not requested.
Q	No analytical results. See report for explanation.
R	Unreliable results. Analyte may or may not be present in the sample. Supporting data is necessary to confirm results.
T	Tentatively Identified Compound. Identified as a result of a library search using the EPA/NIH Mass Spectral Library. Authentic standards were not available to properly identify and quantitate the compound. The reported concentration is an estimate.
TD	Spike recovery too dilute for accurate quantitation.
UI	Not detected. Quantitation limit is estimated.
UL	Not detected. Quantitation limit is probably higher.

Glossary:

()	Numbers in parentheses are analytical spike recoveries (e.g. post-digestion spikes).
[]	Numbers in brackets are matrix spike recoveries (e.g. pre-digestion spikes).
MS/MSD	Matrix spike/matrix spike duplicate; a known increment of target analyte made to a sample before preparation or analyses.
MSA	Method of Standard Additions
RPD	Relative Percent Difference; the results for duplicate analyses are presented as the mean and the relative percent difference.

$$RPD = \frac{|\text{Replicate1} - \text{Replicate2}|}{(\text{Replicate1} + \text{Replicate2})/2} \times 100$$

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(07/01/97)

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INORGANIC ANALYTICAL REPORT

WILMINGTON DRUM
SUPERFUND REMOVAL/REMEDIAL
Lab Request # REQ99016

AR100010

U.S. EPA Region III
Office of Analytical Services
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Annapolis, Maryland

Facility: WILMINGTON DRUM
Program: SUPERFUND REMOVAL/REMEDIAL

Lab Request #: REQ99016

INORGANIC ANALYTICAL SAMPLE RESULTS

Sample Number:	98102101	98102102	98102103	98102104
Station ID :	SD-01	SD-02	SD-03	SD-06
ANALYTES	SAMPLE	SAMPLE	SAMPLE	SAMPLE
Metals Analysis				
	Units: ug/g	ug/g	ug/g	ug/g
Aluminum	3380	23800	12980	23400
Antimony	<0.3	<0.3	3.2	0.8
Arsenic	0.6	57.7	25.7	11.5
Barium	32.3	214	342	346
Beryllium	0.2	1.4	<1.2	1.4
Cadmium	<0.5	<2.5	6.1	<2.5
Calcium	2680	3900	4330	3630
Chromium	9.2	82.0	145	103
Cobalt	3.5	22.0	22.1	22.2
Copper	14.3	33.2	9270	90.5
Iron	7400	61600	52800	42300
Lead	20.5	91.1	12500	1970
Magnesium	1280	5980	3570	6120
Manganese	288	1930	1020	1590
Mercury	<0.1	0.3	0.6	0.4
Nickel	6.9	31.6	100	44.0
Potassium	503	2610	1290	2650
Selenium	<0.5	0.6	1.4	0.7
Silver	<1.0	<5.0	<5.0	<5.0
Sodium	<100	691	599	779
Thallium	<0.3	<0.3	<0.3	<0.3
Vanadium	11.3	61.3	202	59.4
Zinc	73.0	176	869	675
Percent Dry Weight	Units: %	%	%	%
Percent Dry Weight (105C)	71.4	56.2	61.2	56.9
Percent Dry Weight (60C)	73.9	59.7	64.0	57.1

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INORGANIC ANALYTICAL SAMPLE RESULTS

Sample Number:	98102101	98102102	98102103	98102104
Station ID :	SD-01	SD-02	SD-03	SD-06
ANALYTES (continued)	SAMPLE	SAMPLE	SAMPLE	SAMPLE

Total Cyanide	Units: mg/kg	% Rec	RPD	mg/kg	% Rec	RPD	mg/kg	% Rec	RPD
Cyanide	<1.0			<1.0	(98)		<1.0		

Sample Number:	98102105	98102106	98102108	98102109
Station ID :	SD-05	SD-04	PD-1	SW-04
ANALYTES	SAMPLE	SAMPLE	FIELD BLANK	SAMPLE

Metals Analysis	Units: ug/g	% Rec	RPD	ug/g	% Rec	RPD	ug/L	% Rec	RPD
Aluminum	17100			21300		6	721	(89)	16
Antimony	0.6	(94)		0.7	(88)	26	<60		D
Arsenic	20.7	(89)		15.6	(92)	4	<1	(108)	18
Barium	161			192	(106)	1	<200	(93)	D
Beryllium	<1.2			<1.2			<5	(93)	D
Cadmium	<2.5			<2.5			<5	(94)	D
Calcium	5120			3560	(104)	2	38300		11
Chromium	85.9			107	(101)	2	<10	(96)	D
Cobalt	16.1			21.4	(103)	7	<50	(94)	D
Copper	123			272		5	<25	(92)	D
Iron	47200			71400		3	<100	(87)	13
Lead	214			572	(113)	6	<1	(95)	0
Magnesium	4760			5740	(94)	2	<500		3
Manganese	1600			1520		5	<15	(94)	12
Mercury	0.3			0.4	(99)	2	<0.2	(105)	D
Nickel	33.8			56.0	(106)	5	<40	(95)	D
Potassium	1780			2450	(95)	0	<1000	(93)	7
Selenium	0.7	(MSA)		0.7	(114)	12	<2	(90)	D
Silver	<5.0			<5.0	(112)	D	<10	(90)	D
Sodium	558			714	(100)	0	<1000		1
Thallium	<0.3	(108)		<0.3	(99)	D	<3	(94)	D

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INORGANIC ANALYTICAL SAMPLE RESULTS

Sample Number:

Station ID :

ANALYTES (continued)

98102105
SD-05
SAMPLE

98102106
SD-04
SAMPLE

98102109
PB-1
FIELD BLANK

98102109
SW-04
SAMPLE

Metal Analysis

Vanadium

Zinc

Percent Dry Weight

Percent Dry Weight (105C)

Percent Dry Weight (60C)

Total Cyanide

Cyanide

Units: ug/g	% Rec	RPD	ug/g	% Rec	RPD	ug/L	% Rec	RPD	ug/L	% Rec	RPD
54.1			56.8	[99]	3	<50			<50	[93]	D
402			2380		6	<20			<20	[99]	D
Units: %	% Rec	RPD	%	% Rec	RPD						
69.4			55.3		1						
67.5			57.9		3						
Units: mg/kg	% Rec	RPD	mg/kg	% Rec	RPD	mg/L	% Rec	RPD	mg/L	% Rec	RPD
<1.0			<1.0	[107]	D	<0.02			<0.02	[94]	D

Sample Number:

Station ID :

ANALYTES

98102110
SW-06
SAMPLE

98102111
SW-01
SAMPLE

98102112
SW-02
SAMPLE

98102113
SW-03
SAMPLE

Metal Analysis

Aluminum

Antimony

Arsenic

Barium

Beryllium

Cadmium

Calcium

Chromium

Cobalt

Copper

Iron

Lead

Magnesium

Units: ug/L	% Rec	RPD	ug/L	% Rec	RPD	ug/L	% Rec	RPD	ug/L	% Rec	RPD
1990			<200			446			537		
<60			<60	(102)		<60			<60		
2	(112)		<1	(112)		1	(115)		1	(112)	
<200			<200			<200			<200		
<5			<5			<5			<5		
<5			<5			<5			<5		
37800			27500	(102)		34400			36900		
11			<10			<10			<10		
<50			<50			<50			<50		
<25			<25			<25			<25		
3010			215			784			899		
16	(101)		<1	(110)		2	(106)		2	(105)	
48900			10300	(103)		36400			46500		

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INORGANIC ANALYTICAL SAMPLE RESULTS

Sample Number:

Station ID :

ANALYTES (continued)

Metals Analysis

Manganese

Mercury

Nickel

Potassium

Selenium

Silver

Sodium

Thallium

Vanadium

Zinc

98102110

SW-06

SAMPLE

Units: ug/L

156

<0.2

<40

18600

<2

<10

359000

<3

<50

43

% Rec

(97)

(101)

(103)

(102)

(107)

(102)

(107)

(102)

(107)

(102)

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

98102111

SW-01

SAMPLE

ug/L

28

<0.2

<40

4070

<2

<10

15000

<3

<50

<20

% Rec

(103)

(102)

(107)

(102)

(107)

(102)

(107)

(102)

(107)

(102)

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

98102112

SW-02

SAMPLE

ug/L

89

<0.2

<40

14300

<2

<10

270000

<3

<50

<20

% Rec

(99)

(102)

(107)

(102)

(107)

(102)

(107)

(102)

(107)

(102)

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

98102113

SW-03

SAMPLE

ug/L

87

<0.2

<40

17700

<2

<10

344000

<3

<50

<20

% Rec

(98)

(101)

(101)

(101)

(101)

(101)

(101)

(101)

(101)

(101)

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

RPD

Sample Number:

Station ID :

ANALYTES

Metals Analysis

Aluminum

Antimony

Arsenic

Barium

Beryllium

Cadmium

Calcium

Chromium

Cobalt

98102114

SW-05

SAMPLE

Units: ug/L

1550

<50

1

<200

<5

<5

32400

<10

<50

% Rec

(115)

RPD

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INORGANIC ANALYTICAL SAMPLE RESULTS

Sample Number: 98102114

Station ID: SW-05

ANALYTES (continued)
SAMPLE

Metal Analysis	Units: ug/L	% Rec	RPD
Copper	<25		
Iron	2710		
Lead	7	(106)	
Magnesium	25500		
Manganese	145		
Mercury	<0.2		
Nickel	<40		
Potassium	14000		
Selenium	<2	(99)	
Silver	<10		
Sodium	259000		
Thallium	<3	(102)	
Vanadium	<50		
Zinc	36		
Total Cyanide	units: mg/L	% Rec	RPD
Cyanide	<0.02	(100)	

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INORGANIC WET/DRY CONVERSION

Sample Number:		98102101		98102102		98102103		98102104		98102105	
ANALYTES		WET	DRY	WET	DRY	WET	DRY	WET	DRY	WET	DRY
		SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE
Percent Dry Weight		Units: %		%		%		%		%	
Percent Dry Weight (105C)		71.4		56.2		61.2		56.9		69.4	
Percent Dry Weight (60C)		73.9		59.7		64.0		57.1		67.5	
Metals Analysis		Units: ug/g		ug/g		ug/g		ug/g		ug/g	
Aluminum	2500	3360	3360	14200	23800	8260	12900	13400	23400	11500	17100
Antimony	< 0.2	< 0.3	< 0.3	< 0.2	< 0.3	2.0	3.2	0.5	0.8	0.4	0.6
Arsenic	0.4	0.6	57.7	16.4	25.7	16.4	25.7	6.6	11.5	14.0	20.7
Barium	23.9	32.3	214	219.9	342	219.9	342	197.6	346	108.7	161
Beryllium	0.1	0.2	1.4	0.9	1.4	< 0.8	< 1.2	0.8	1.4	< 0.8	< 1.2
Cadmium	< 0.4	< 0.5	< 2.5	3.9	6.1	< 1.4	< 2.5	< 1.4	< 2.5	< 1.7	< 2.5
Calcium	1980	2680	3900	2770	4330	2190	3830	56.6	103	3460	5120
Chromium	6.8	9.2	62.0	92.6	145	12.7	22.2	12.7	22.2	58.0	85.9
Cobalt	2.6	3.5	22.0	14.1	22.1	12.7	22.2	12.7	22.2	10.9	16.1
Copper	10.6	14.3	33.2	530	9270	51.7	90.5	51.7	90.5	83.0	123
Iron	5470	7400	61600	33800	52800	24200	42300	24200	42300	31900	47200
Lead	15.1	20.5	91.1	8000	12500	1120	1970	1120	1970	144.5	214
Magnesium	945.9	1280	5980	2280	3570	3500	6130	3500	6130	3210	4760
Manganese	212.8	288	1930	652.8	1020	907.9	1590	907.9	1590	1080	1600
Mercury	< 0.1	< 0.1	0.3	0.5	0.8	0.2	0.4	0.2	0.4	0.2	0.3
Nickel	5.1	6.9	31.6	64.0	100	25.1	44.0	25.1	44.0	22.8	33.6
Potassium	371.7	503	2610	825.6	1290	1510	2650	1510	2650	1200	1780
Selenium	< 0.4	< 0.5	0.6	0.9	1.4	0.4	0.7	0.4	0.7	0.5	0.7
Silver	< 0.7	< 1.0	< 5.0	< 3.2	< 5.0	< 2.9	< 5.0	< 2.9	< 5.0	< 3.4	< 5.0
Sodium	< 73.9	< 100	691	382.7	598	444.8	779	444.8	779	444.2	658
Thallium	< 0.2	< 0.3	< 0.3	< 0.2	< 0.3	< 0.2	< 0.3	< 0.2	< 0.3	< 0.2	< 0.3
Vanadium	8.4	11.3	61.3	129.3	202	33.9	59.4	33.9	59.4	36.5	54.1
Zinc	53.9	73.0	176	556.2	969	395.4	675	395.4	675	271.4	402

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Facility: WILMINGTON DRUM
Program: SUPERFUND REMOVAL/REMEDIAL

Lab Request #: REQ99016

INORGANIC WET/DRY CONVERSION

Sample Number:

98102101	98102102	98102103	98102104	98102105
WET	WET	WET	WET	WET
DRY	DRY	DRY	DRY	DRY
SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE

ANALYTES (continued)

Total Cyanide
Cyanide

Units: mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
<1.0	<1.4	<1.0	<1.0	<1.4

Sample Number:

98102106
WET
DRY
SAMPLE

ANALYTES

Percent Dry Weight

Units: %
Percent Dry Weight (105C)
Percent Dry Weight (60C)

Metals Analysis

Aluminum	21300	21300
Antimony	0.4	0.7
Arsenic	9.1	15.8
Barium	111.2	192
Beryllium	<= 0.7	<= 1.2
Cadmium	<= 1.4	<= 2.5
Calcium	2060	3560
Chromium	62.0	107
Cobalt	12.4	21.4
Copper	157.5	272
Iron	41300	71400
Lead	331.2	572
Magnesium	3320	5740
Manganese	880.1	1520
Mercury	0.2	0.4
Nickel	32.4	56.0
Potassium	1420	2450
Selenium	0.4	0.7
Silver	< 2.9	< 5.0

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INORGANIC WET/DRY CONVERSION

Sample Number:

98102106

WET	DRY
SAMPLE	SAMPLE

ANALYTES (continued)

Metal Analysis	Units: ug/g	ug/g
Sodium	413.4	714
Thallium	< 0.2	< 0.3
Vanadium	32.9	56.6
Zinc	1380	2380
Total Cyanide	Units: mg/kg	mg/kg
Cyanide	< 1.0	< 1.8

The wet/dry weight conversion table is presented for the convenience of the data user. The attached analyst narratives state whether a test was determined on a wet or dry weight basis. Any difference in significant figures between the two columns is an artifact of the reporting function and does not imply that the calculated value has greater precision than the measured value.

AR100018

Percent Dry Weight Determinations

Analyst:

Steven A. Bartimo
LMES Chemist

Method:

The soil samples from the WILMINGTON DRUM site (REQ99016) were analyzed for Percent Dry Weight as required by EPA analytical methods. The samples were dried at 105°C following the procedure outlined in the EPA Region III OASQA laboratory SOP #R3QA056.0.

These results are to be used to convert analyte concentrations to a dry weight basis for organic and non-metal analyses. Normally, analytical values are reported on a wet weight basis for organic and non-metals. All metals reported use a 60°C drying temperature for the percent dry weight determinations, as required by the methodology. The 60°C percent dry weight values are reported with the metals results, if applicable.

Metals Determinations**Analysts:**

R. Fisher M. Mecanic
LMES Chemist LMES Chemist

Methods:

Samples from WILMINGTON DRUM site (REQ99016) were prepared for analysis by acid digestion and analyzed by graphite furnace atomic absorption spectroscopy (GFAAS) and inductively coupled plasma-atomic emission spectrometry (ICP-AES). The results are presented in the attached table. The following are the digestion and analytical techniques and methods employed:

Digestion Methods

Method from CLP SOW 9/91 revision, p. D-5, A.1. for GFAAS (excluding antimony)

Method from CLP SOW 9/91 revision, p. D-5, A.2. for ICP-AES

EPA Method 3050A, excluding HCl for GFAAS (excluding antimony) (solid samples) (2)

EPA Method 3050A, for ICP-AES and antimony by GFAAS (solid samples) (2)

Analytical Methods

EPA Method 204.2 and Internal SOP R3-QA132, antimony by GFAAS (1)

EPA Method 206.2 and Internal SOP R3-QA132, arsenic by GFAAS (1)

EPA Method 239.2 and Internal SOP R3-QA132, lead by GFAAS (1)

EPA Method 270.2 and Internal SOP R3-QA132, selenium by GFAAS (1)

EPA Method 279.2 and Internal SOP R3-QA132, thallium by GFAAS (1)

EPA Method 200.7 and Internal SOP R3-QA132, remaining elements by ICP-AES (1)

(1) EPA Manual of Methods for Chemical Analysis of Water and Wastes, 1979/83.

(2) SW-846, Test Methods for Evaluating Solid Waste Physical/Chemical Methods, Volume IA, Final Update I, July 1992.

Results for solid samples are reported in ug/g (ppm) DRY weight at 60 degrees centigrade. This, Percent Dry Weight test, pertains only to metals results. The drying temperature of 60 degrees centigrade is selected to retain volatile elements. The % Dry Weight (60 °C) is reported to allow for conversion to wet weight.

Note: The quantitation limits for silver, beryllium and cadmium in the following samples are elevated due to dilutions for matrix interferences:

Silver: 981021-02 through 981021-06
Beryllium: 981021-03, 981021-05, 981021-06
Cadmium: 981021-02, 981021-04, 981021-05, 981021-06

Mercury Determinations

Analysts:

Odis Cain
LMES Chemist

Steve Bartimo
LMES Chemist

Method:

Samples from WILMINGTON DRUM site (REQ99016) were analyzed for total mercury using EPA Methods 245.1¹ and 245.5¹.

Results for solid samples are reported in ug/g Dry weight at 60 degrees centigrade. This Percent Dry Weight test pertains only to metals results. The drying temperature of 60 degrees centigrade is selected to retain volatile elements. The Percent Dry Weight (60 degrees C) is reported to allow for conversion to wet weight.

¹Methods for Chemical Analysis of Water and Wastes, EPA 600/4-79-020.

Total Cyanide Determinations

Analyst:

Odis W. Cain
LMES CHEMIST

Method:

Samples from WILMINGTON DRUM site (REQ99016) were analyzed for Total Cyanide determination using EPA Method 335.4.

Soil results are reported on a wet weight basis.

U.S. EPA Region III
Office of Analytical Services
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Annapolis, Maryland

ORGANIC ANALYTICAL REPORT

WILMINGTON DRUM
SUPERFUND REMOVAL/REMEDIAL
Lab Request # REQ99016

AR100024

U.S. EPA Region III
Office of Analytical Services
and Quality Assurance
Annapolis, Maryland

Facility: WILMINGTON DRUM
Program: SUPERFUND REMOVAL/REMEDIAL

Lab Request #: REQ99016

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ORGANIC ANALYTICAL SAMPLE RESULTS

Sample Number:	98102101	98102102	98102103	98102104	98102105	98102106	98102107	98102108
Station ID :	SD-01	SD-02	SD-03	SD-06	SD-05	SD-04	TR-1	FB-1
ANALYTES	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	TRIP BLANK	FIELD BLANK
Percent Dry Weight	71.4	56.2	61.2	56.9	69.4	55.3		
Percent Dry Weight (105C)								
NQL Factor:	1	1	1	1	1	1	1	1
PCBs and Pesticides by Gas Chromatography Units:	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	ug/L
Alpha Chlordane			0.014 R,I					
Aroclor 1260			0.25	0.090	0.72	0.081		
4,4'-DDD				0.0088 I				
4,4'-DDE			0.016 R,I	0.0074 R,I				
4,4'-DDT			0.050 R,I	0.0095 R,I	0.030 R,I	0.0079 R,I		
Endosulfon-II					0.024 I			
Endrin			0.011 R,I	0.0097 I				
Gamma Chlordane								
NQL Factor:	1	1	1	1	1	1	1	1
Semivolatile Organics by GC/MS	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	ug/L
Anthracene			0.06 J	0.04 J	0.07 J			
Benzo(a)Anthracene	0.2 J	0.3 J	0.3 J	0.2 J	0.3 J	0.1 J		
Benzo(a)Pyrene	0.1 J	0.3 J	0.3 J	0.2 J	0.2 J	0.1 J		
Benzo(b)Fluoranthene	0.1 J	0.3 J	0.3 J	0.2 J	0.2 J	0.1 J		
Benzo(g,h,i)Perylene	0.08 J	0.2 J	0.2 J	0.1 J	0.2 J	0.08 J		
Benzo(k)Fluoranthene	0.2 J	0.2 J	0.2 J	0.2 J	0.2 J	0.2 J		
Bis(2-Ethylhexyl)phthalate			0.07 B	0.04 B	0.2 B	0.38		
Carbazole	UU	UU	0.05 J	UU	UU	UU	UU	UU
4-Chloroaniline	UU	UU	UU	UU	UU	UU	UU	UU
Chrysene	0.03 J	0.2 J	0.35	0.2 J	0.3 J	0.1 J		
Di-n-Butylphthalate								
Dibenzo(a,h)Anthracene			0.04 J					
4,6-Dinitro-2-Methylphenol		0.05 J	0.09 J	0.06 J	0.1 J	0.04 J		
2,4-Dinitrophenol	UU	UU	UU	UU	UU	UU		
Fluoranthene	0.04 J	0.2 J	0.49	0.3 J	0.47	0.2 J		
Indeno(1,2,3-cd)Pyrene		0.1 J	0.2 J	0.1 J	0.2 J	0.09 J		
2-Methylnaphthalene			0.04 J					
Naphthalene			0.06 J	0.08 J	0.08 J			

ORGANIC ANALYTICAL SAMPLE RESULTS

Sample Number:	98102101	98102102	98102103	98102104	98102105	98102106	98102107	98102108
Station ID :	SD-01	SD-02	SD-03	SD-06	SD-05	SD-04	TS-1	FB-1
ANALYTES (continued)	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	TRIP BLANK	FIELD BLANK
NOL Factor:	1	1	1	1	1	1	1	1
Semivolatile Organics by GC/MS	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/L	ug/L
3-Nitroaniline	UU	UU	UU	UU	UU	UU	UU	UU
Phenanthrene	0.03 J	0.09 J	0.3 J	0.2 J	0.3 J	0.07 J	0.07 J	0.07 J
Pyrene	0.03 J	0.2 J	0.51	0.3 J	0.52	0.2 J	0.2 J	0.2 J
NOL Factor:	1	1	1	1	1	1	1	1
Volatile Organics by Purge & Trap GC/MS	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/kg	ug/L	ug/L
Bromobenzene	UU	UU	UU	UU	UU	UU	UU	UU
2-Chloroethylvinyl Ether	UU	UU	UU	UU	UU	UU	UU	UU
2-Chlorotoluene	UU	UU	UU	UU	UU	UU	UU	UU
4-Chlorotoluene	UU	UU	UU	UU	UU	UU	UU	UU
1,2-Dibromo-3-Chloropropane	UU	UU	UU	UU	UU	UU	UU	UU
2,2-Dichloropropane	UU	UU	UU	UU	UU	UU	UU	UU
Naphthalene	UU	UU	UU	UU	UU	UU	UU	UU
N-Propylbenzene	UU	UU	UU	UU	UU	UU	UU	UU
1,2,3-Trichlorobenzene	UU	UU	UU	UU	UU	UU	UU	UU
1,2,3-Trichloropropane	UU	UU	UU	UU	UU	UU	UU	UU
1,2,4-Trimethylbenzene	UU	UU	UU	UU	UU	UU	UU	UU
Vinyl Acetate	UU	UU	UU	UU	UU	UU	UU	UU

Lab Request #: RQ99016

Facility: WILMINGTON DRUM
Program: SUPERFUND REMOVAL/REMEDIAL

ORGANIC ANALYTICAL SAMPLE RESULTS

Sample Number:	981021109	981021110	981021111	981021112	981021113	981021114
Station ID :	SW-04	SW-06	SW-01	SW-02	SW-03	SW-05
ANALYTES	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE
NOL Factor:	1	1	1	1	1	1
Semivolatile Organics by GC/MS	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Carbazole	UU	UU	UU	UU	UU	UU
4-Chloroaniline	UU	UU	UU	UU	UU	UU
4,6-Dinitro-2-Methylphenol	UU	UU	UU	UU	UU	UU
2,4-Dinitrophenol	UU	UU	UU	UU	UU	UU
3-Nitroaniline	UU	UU	UU	UU	UU	UU
NOL Factor:	1	1	1	1	1	1
Volatile Organics by Purge & Trap GC/MS	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Bromobenzene	UU	UU	UU	UU	UU	UU
2-Chloroethylvinyl Ether	Q	Q	Q	Q	Q	Q
2-Chlorotoluene	UU	UU	UU	UU	UU	UU
4-Chlorotoluene	UU	UU	UU	UU	UU	UU
Naphthalene	UU	UU	UU	UU	UU	UU
N-Propylbenzene	UU	UU	UU	UU	UU	UU
1,2,3-Trichloropropane	UU	UU	UU	UU	UU	UU
1,2,4-Trimethylbenzene	UU	UU	UU	UU	UU	UU
Vinyl Acetate	UU	UU	UU	UU	UU	UU

U.S. EPA Region III
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Annapolis, Maryland

Facility: WILMINGTON DRUM
Program: SUPERFUND REMOVAL/REMEDIAL

Lab Request #: REQ99016

ORGANIC QUALITY CONTROL (SURROGATE RECOVERIES)

Matrix: SOLIDS

Surrogate	98102101	98102102	98102103	98102104	98102105	98102105
	SD-01	SD-02	SD-03	SD-06	SD-05	SD-04
Limits	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE

PCBs and Pesticides by Gas Chromatography

Range	%	%	%	%	%	%
(60-150)	112	109	160 A	114	120	109
(60-150)	106	104	101	103	107	99

Semivolatile Organics by GC/MS

Range	%	%	%	%	%	%
(30-115)	72	77	80	80	81	61
(25-121)	68	59	69	66	66	48
(23-120)	66	54	65	63	64	47
(24-113)	69	69	75	72	72	52
(18-137)	93	90	94	94	98	73
(19-122)	47	44	48	47	55	27

Volatile Organics by Purge & Trap GC/MS

Range	%	%	%	%	%	%
(59-113)	82	109	80	97	87	80
(70-121)	104	105	109	102	100	108
(84-138)	98	97	105	115	101	103

Matrix: WATER

Surrogate	98102107	98102108	98102109	98102110	98102111	98102112	98102113	98102114
	TR-1	TR-1	SW-04	SW-06	SW-01	SW-02	SW-03	SW-05
Limits	TRP BLANK	TRP BLANK	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE

PCBs and Pesticides by Gas Chromatography

Range	%	%	%	%	%	%	%	%
(60-150)	89	49 A	49 A	56 A	83	51 A	54 A	52 A
(60-150)	101	90	90	92	104	94	96	98

Semivolatile Organics by GC/MS

Range	%	%	%	%	%	%	%	%
(43-116)	69	66	66	67	64	73	66	66
(21-110)	73	72	72	70	63	74	56	70
(35-114)	75	69	69	68	68	79	63	71
(10-110)	63	64	64	63	66	75	63	71
(13-141)	60	60	60	60	60	60	60	60

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SECRET

ORGANIC QUALITY CONTROL (SURROGATE RECOVERIES)

SURROGATES (continued)		99102107	99102108	99102109	99102110	99102111	99102112	99102113	99102114
Limits		TD-1	PD-1	SW-04	SW-06	SW-01	SW-02	SW-03	SW-05
		TRP BLANK FLD BLANK	TRP BLANK FLD BLANK	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE	SAMPLE
Semivolatile Organics by GC/MS	Range	0	0	0	0	0	0	0	0
	(10-123)	62	74	74	86	86	86	86	86
Volatile Organics by purge & trap GC/MS	Range	0	0	0	0	0	0	0	0
	(66-115)	103	99	109	101	99	100	105	103
	(76-114)	92	110	105	105	117 A	110	106	113
	(88-110)	101	91	95 A	92	90	93	92	90
	d8-Toluene								

Facility: WILMINGTON DRUM
Program: SUPERFUND REMOVAL/REMEDIAL

Lab Request #: REQ99016

ORGANIC QUALITY CONTROL (MATRIX SPIKE RECOVERIES)

Matrix: SOLIDS

98102106

SD-04

ANALYTES

Pest and Pesticides by Gas Chromatography

ANALYTES	Spike Recovery		Recovery		RPD	
	MS	MSD	Limit	Range	RPD	Limit
Aldrin	90	99	(34-132)		2	43
4,4' DDT	88	155 A	(23-134)		55 A	50
Dieldrin	98	97	(31-134)		1	38
Endrin	111	110	(42-139)		1	45
Gamma BHC (Lindane)	97	94	(46-127)		3	50
Heptachlor	121	115	(35-130)		5	31

Semivolatile Organics by GC/MS

ANALYTES	Spike Recovery		Recovery		RPD	
	%	%	Limit	Range	RPD	Limit
Acenaphthene	81	81	(31-137)		0	19
4-Chloro-3-Methylphenol	84	83	(26-103)		2	33
2-Chlorophenol	60	64	(25-102)		6	50
Di-n-Butylphthalate	94	94	(11-117)		1	40
1,4-Dichlorobenzene	43	52	(28-104)		18	27
2,4-Dinitrotoluene	72	73	(28-89)		1	47
4-Nitrophenol	79	79	(11-114)		0	50
N-Nitroso-di-n-Propylamine	54	57	(41-126)		6	38
Pentachlorophenol	15 A	21	(17-109)		35	47
Phenol	65	67	(26-90)		3	35
Pyrene	93	91	(35-142)		3	36
1,2,4-Trichlorobenzene	52	58	(38-107)		11	23

Volatile Organics by purge & Trap GC/MS

ANALYTES	Spike Recovery		Recovery		RPD	
	%	%	Limit	Range	RPD	Limit
Benzene	98	103	(66-142)		5	21
Chlorobenzene	98	109	(60-133)		10	21
1,1-Dichloroethene	75	80	(59-172)		6	22
Toluene	94	111	(59-139)		16	21
Trichloroethene	99	92	(62-137)		4	24

AP-000030

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ORGANIC QUALITY CONTROL (MATRIX SPIKE RECOVERIES)

Matrix: WATER

98102109
SM-04

ANALYTES

PCBs and Pesticides by Gas Chromatography

	Spike Recovery		Recovery	RPD	
	MS	MSD	Limit	RPD	Limit
Aldrin	59	92	(40-120)	44 A	22
4,4'-DDT	46	75	(38-127)	48 A	27
Dieldrin	70	99	(52-126)	35 A	18
Endrin	74	106	(56-121)	35 A	21
Gamma BHC (Lindane)	71	100	(56-123)	34 A	15
Heptachlor	61	91	(40-131)	39 A	20

Semivolatile Organics by GC/MS

	%	Range	RPD	Limit
Acenaphthene	78	(46-118)	10	31
4-Chloro-3-Methylphenol	86	(23- 97)	3	42
2-Chlorophenol	55	(27-123)	16	40
Di-n-Butylphthalate	95	(11-117)	9	40
1,4-Dichlorobenzene	55	(36- 97)	2	28
2,4-Dinitrotoluene	74	(24- 96)	12	38
4-Nitrophenol	99 A	(10- 80)	5	50
N-Nitroso-di-n-Propylamine	56	(41-116)	18	38
Pentachlorophenol	73	(9-103)	3	50
Phenol	54	(12-110)	14	42
Pyrene	98	(26-127)	6	31
1,2,4 Trichlorobenzene	53	(39- 98)	6	28

Volatile Organics by Purge & Trap GC/MS

	%	Range	RPD	Limit
Benzene	93	(76-127)	5	11
Chlorobenzene	95	(75-130)	2	13
1,1-Dichloroethene	79	(61-145)	5	14
Toluene	78	(76-125)	0	13
Trichloroethene	93	(71-120)	11	14

ARI00031

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LABORATORY REAGENT BLANK RESULTS

Method: PCBs and Pesticides by Gas Chromatography

Run: OC981022 1 LRB

SURROGATES

Decachlorobiphenyl
Tetrachloro-M-Xylene

NOL Factor	% Rec
1	97
1	91

Run: OC981026 1 LRB

SURROGATES

Decachlorobiphenyl
Tetrachloro-M-Xylene

NOL Factor	% Rec
1	104
1	100

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Facility: WILMINGTON DRUM
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LABORATORY REAGENT BLANK RESULTS

Method: Semivolatile Organics by GC/MS

Run: 0198B231 18 LRB

SURROGATES

ANALYTES	NOL Factor	% Rec
2-Fluoro-1,1'-Biphenyl	1	68
2-Fluorophenol	1	78
d5-Nitrobenzene	1	69
d5-Phenol	1	65
d14-Terphenyl	1	96
2,4,6-Tribromophenol	1	50

ANALYTES

ANALYTES	NOL Factor	mg/kg
Bis(2-Ethylhexyl)Phthalate	1	0.03 U
Carbazole	1	UU
4-Chloroaniline	1	UU
2,4-Dinitrophenol	1	UU
3-Nitroaniline	1	UU

AD-1000033

Run: 0198B231 19 LRB

SURROGATES

ANALYTES	NOL Factor	% Rec
2-Fluoro-1,1'-Biphenyl	1	78
2-Fluorophenol	1	80
d5-Nitrobenzene	1	80
d5-Phenol	1	70
d14-Terphenyl	1	100
2,4,6-Tribromophenol	1	70

ANALYTES

ANALYTES	NOL Factor	ug/L
Carbazole	1	UU
4-Chloroaniline	1	UU
2,4-Dinitrophenol	1	UU
3-Nitroaniline	1	UU

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LABORATORY REAGENT BLANK RESULTS

Method: Volatile Organics by Purge & Trap GC/MS

Run: 08981123 17 LRB

SURROGATES

Bromofluorobenzene	NOL Factor	% Rec
d4-1,2-Dichloroethane	1	81
d8-Toluene	1	102
	1	97

ANALYTES

Benzene	NOL Factor	ug/kg
Chlorobenzene	1	2 J
2-Chloroethylvinyl Ether	1	2 J
1,2-Dibromo-3-Chloropropane	1	UU
Methylene Chloride	1	UU
Naphthalene	1	1 J
Toluene	1	UU
1,2,3-Trichlorobenzene	1	2 J
Trichloroethene	1	UU
	1	2 J

AR 100034

Run: 08981123 18 LRB

SURROGATES

Bromofluorobenzene	NOL Factor	% Rec
d4-1,2-Dichloroethane	1	103
d8-Toluene	1	104
	1	86

ANALYTES

Bromobenzene	NOL Factor	ug/L
2-Chloroethylvinyl Ether	1	UU
2-Chlorotoluene	1	Q
4-Chlorotoluene	1	UU
Methylene Chloride	1	UU
Naphthalene	1	2 J
N-Propylbenzene	1	UU
	1	UU

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LABORATORY REAGENT BLANK RESULTS

1,2,4-Trimethylbenzene

1 UJ

Run: CH981123 19 LRB

SURROGATES

Bromofluorobenzene

NQL Factor 1 Rec 143 A

d4-1,2-Dichloroethane

111

d8-Toluene

80 A

ANALYTES

Bromobenzene

NQL Factor 1 Rec UJ

2-Chloroethylvinyl Ether

Q

2-Chlorotoluene

UJ

4-Chlorotoluene

UJ

Dibromochloromethane

UJ

2,2-Dichloropropane

UJ

N-Propylbenzene

UJ

1,2,3-Trichloropropane

UJ

Run: CH981123 20 LRB

SURROGATES

Bromofluorobenzene

NQL Factor 1 Rec 83

d4-1,2-Dichloroethane

102

d8-Toluene

101

ANALYTES

2-Chloroethylvinyl Ether

NQL Factor 1 Rec UJ

1,2-Dibromo-3-Chloropropane

UJ

2,2-Dichloropropane

UJ

Naphthalene

UJ

1,2,3-Trichlorobenzene

UJ

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LABORATORY REAGENT BLANK RESULTS

Method: Volatile Organics by Purge & Trap GC/MS

Run: 0H981123 21 LRB

SURROGATES

	NQL Factor	% Rec
Bromofluorobenzene	1	104
d4-1,2-Dichloroethane	1	89
d8-Toluene	1	102

ANALYTES

	NQL Factor	ug/L
Methylene Chloride	1	1 J
Vinyl Acetate	1	UV

AR100036

Office of Analytical Services and Quality Assurance
Volatile Organics Analysis
Nominal Quantitation Limits (NQL)

Units: Water = ug/L NPTC = Non-Priority Pollutant Target Compound

Actual Quantitation Limit = (NQL Factor) X NQL

CAS	ANALYTE	NQL
75-71-8	Dichlorodifluoromethane	5
74-87-3	Chloromethane	5
75-01-4	Vinyl Chloride	5
74-83-9	Bromomethane	5
75-00-3	Chloroethane	5
75-69-4	Trichlorofluoromethane	5
75-35-4	1,1-Dichloroethylene	5
75-15-0	Carbon Disulfide	NPTC 5
67-64-1	Acetone	NPTC 5
75-09-2	Methylene Chloride	5
156-60-5	trans-1,2-Dichloroethane	5
75-34-3	1,1-Dichloroethane	5
108-05-4	Vinyl Acetate	NPTC 5
590-20-7	2,2-Dichloropropane	5
156-59-4	cis-1,2-Dichloroethane	NPTC 5
78-93-3	2-Butanone	NPTC 5
74-97-5	Bromochloromethane	NPTC 5
65-66-3	Chloroform	5
71-55-6	1,1,1-Trichloroethane	5
56-23-5	Carbon Tetrachloride	5
63-58-6	1,1-Dichloropropene	5
71-43-2	Benzene	5
107-06-2	1,2-Dichloroethane	5
79-01-6	Trichloroethylene	5
78-87-5	1,2-Dichloropropane	5
74-95-3	Dibromomethane	NPTC 5
75-27-4	Bromodichloromethane	5
110-75-8	2-Chloroethylvinyl ether	5
10061-01-6	trans-1,3-Dichloropropene	NPTC 5
108-10-1	4-Methyl-2-pentanone	5
108-83-3	Toluene	5
10061-01-5	cis-1,3-Dichloropropene	5
79-00-5	1,1,2-Trichloroethane	5
127-18-4	Tetrachloroethylene	5

CAS	ANALYTE	NPTC	NQL
142-28-9	1,3-Dichloropropane	NPTC	5
591-78-6	2-Hexanone	NPTC	5
124-48-1	Dibromochloromethane		5
106-93-4	1,2-Dibromoethane (EDB)	NPTC	5
108-90-7	Chlorobenzene		5
620-20-6	1,1,1,2-Tetrachloroethane	NPTC	5
100-41-4	Ethylbenzene		5
108-38-3	m-Xylene (m & p isomers)	NPTC	5
106-42-3	p-Xylene (together)	NPTC	5
95-47-6	o-Xylene	NPTC	10
100-42-5	Styrene	NPTC	10
75-25-2	Bromoform		5
98-82-8	Isopropylbenzene	NPTC	5
108-86-1	Bromobenzene	NPTC	5
79-34-5	1,1,2,2-Tetrachloroethane		5
96-18-4	1,2,3-Trichloropropane	NPTC	5
103-65-1	n-Propylbenzene	NPTC	5
95-49-8	2-Chlorotoluene	NPTC	5
106-43-4	4-Chlorotoluene	NPTC	5
108-67-8	1,3,5-Trimethylbenzene	NPTC	5
98-06-6	tert-Butylbenzene	NPTC	5
93-63-6	1,2,4-Trimethylbenzene	NPTC	5
135-98-8	sec-Butylbenzene	NPTC	5
541-73-1	1,3-Dichlorobenzene		5
106-46-7	1,4-Dichlorobenzene		5
99-87-6	p-Isopropyltoluene	NPTC	5
95-50-1	1,2-Dichlorobenzene		5
104-51-8	n-Butylbenzene	NPTC	5
96-12-8	1,2-Dibromo-3-chloropropane		5
120-82-1	1,2,4-Trichlorobenzene		5
91-20-3	Naphthalene		5
87-68-3	Hexachlorobutadiene		5
87-61-6	1,2,3-Trichlorobenzene	NPTC	5

The "Nominal Quantitation Limit" factor is an overall correction factor applied to the method's NQLs for analytical adjustments made during the analysis (i.e., for analyses of less than the ideal 5 mls for aqueous samples, and for sample dilutions prior to analysis).

For example, if the NQL factor for an OASQA water sample is 2, the estimated Actual Quantitation Limit for vinyl chloride would be 10 ug/L (i.e., 2 x 5 ug/L).

Rev. October 16, 1996

AR100037

Office of Analytical Services and Quality Assurance
Volatile Organics Analysis
Nominal Quantitation Limits (NQL)

Units: Solids = ug/kg (wet) NPTC = Non-Priority Pollutant Target Compound

Actual Quantitation Limit = (NQL Factor) X NQL

CAS	ANALYTE	NQL
75-71-8	Dichlorodifluoromethane	5
74-87-3	Chloromethane	5
75-01-4	Vinyl Chloride	5
74-83-9	Bromomethane	5
75-00-3	Chloroethane	5
75-69-4	Trichlorofluoromethane	5
75-35-4	1,1-Dichloroethylene	5
75-15-0	Carbon Disulfide	NPTC 5
67-64-1	Acetone	NPTC 5
75-09-2	Methylene Chloride	5
156-60-5	trans-1,2-Dichloroethene	5
75-34-3	1,1-Dichloroethane	5
108-05-4	Vinyl Acetate	NPTC 5
590-20-7	2,2-Dichloropropane	5
156-59-4	cis-1,2-Dichloroethene	NPTC 5
78-93-3	2-Butanone	NPTC 5
74-97-5	Bromochloromethane	NPTC 5
65-65-3	Chloroform	5
71-55-6	1,1,1-Trichloroethane	5
56-23-5	Carbon Tetrachloride	5
563-58-6	1,1-Dichloropropene	5
71-43-2	Benzene	5
107-06-2	1,2-Dichloroethane	5
79-01-6	Trichloroethylene	5
78-87-5	1,2-Dichloropropane	5
74-95-3	Dibromomethane	NPTC 5
75-27-4	Bromodichloromethane	5
110-75-8	2-Chloroethylvinyl ether	5
10061-01-6	trans-1,3-Dichloropropene	NPTC 5
108-10-1	4-Methyl-2-pentanone	5
108-88-3	Toluene	5
10061-01-5	cis-1,3-Dichloropropene	5
79-00-5	1,1,2-Trichloroethane	5
127-18-4	Tetrachloroethylene	5

CAS	ANALYTE	NQL
142-28-9	1,3-Dichloropropane	NPTC 5
591-78-6	2-Hexanone	NPTC 5
124-48-1	Dibromochloromethane	5
106-93-4	1,2-Dibromoethane (EDB)	NPTC 5
108-90-7	Chlorobenzene	5
630-20-6	1,1,1,2-Tetrachloroethane	NPTC 5
100-41-4	Ethylbenzene	5
108-38-3	m-Xylene (m & p isomers)	NPTC 5
106-42-3	p-Xylene (together)	NPTC 5
95-47-6	o-Xylene	NPTC 10
100-42-5	Styrene	NPTC 10
75-25-2	Bromoform	5
98-82-8	Isopropylbenzene	NPTC 5
108-86-1	Bromobenzene	NPTC 5
79-34-5	1,1,2,2-Tetrachloroethane	5
96-18-4	1,2,3-Trichloropropane	NPTC 5
103-65-1	n-Propylbenzene	NPTC 5
95-49-8	2-Chlorotoluene	NPTC 5
106-43-4	4-Chlorotoluene	NPTC 5
108-67-8	1,3,5-Trimethylbenzene	NPTC 5
98-06-6	tert-Butylbenzene	NPTC 5
93-63-6	1,2,4-Trimethylbenzene	NPTC 5
135-98-8	sec-Butylbenzene	NPTC 5
541-73-1	1,3-Dichlorobenzene	5
106-46-7	1,4-Dichlorobenzene	5
99-87-6	p-Isopropyltoluene	NPTC 5
95-50-1	1,2-Dichlorobenzene	5
104-51-8	n-Butylbenzene	NPTC 5
96-12-8	1,2-Dibromo-3-chloropropane	5
120-82-1	1,2,4-Trichlorobenzene	5
91-20-3	Naphthalene	5
87-68-3	Hexachlorobutadiene	5
87-61-6	1,2,3-Trichlorobenzene	NPTC 5

The "Nominal Quantitation Limit" factor is an overall correction factor applied to the method's NQLs for analytical adjustments made during the analysis (i.e., for analyses of more or less than the ideal 5 grams for soil samples, for sample dilutions prior to analysis).

For example, if the NQL factor for an OASQA soil sample is 2, the estimated Actual Quantitation Limit for vinyl chloride would be 10 ug/kg (i.e., 2 x 5 ug/kg).

Office of Analytical Services and Quality Assurance
Pesticide and PCB Analysis
Nominal Quantitation Limits (NQL)

Units: Water = ug/L NPIC = Non-Priority Pollutant Target Compound

Actual Quantitation Limit = (NQL Factor) X NQL

CAS	Pesticide	NQL
319-84-6	Alpha-BHC	0.05
319-85-7	Beta-BHC	0.05
319-86-8	Delta-BHC	0.05
58-89-8	Gamma-BHC	0.05
76-44-8	Heptachlor	0.05
309-00-2	Aldrin	0.05
1024-57-3	Heptachlor Epoxide	0.05
959-98-8	Endosulfan I	0.05
60-57-1	Dieldrin	0.10
72-55-9	4,4'-DDE	0.10
72-20-8	Endrin	0.10
33213-65-9	Endosulfan II	0.10
72-54-8	4,4'-DDD	0.10
1031-07-8	Endosulfan Sulfate	0.10
50-29-3	4,4'-DDT	0.10
7421-93-4	Endrin Aldehyde	0.10
53494-70-5	Endrin Ketone	NPIC 0.10
72-43-5	Methoxychlor	NPIC 0.50
5103-71-9	Alpha-Chlordane	0.05
5103-74-2	Gamma-Chlordane	0.05
57-74-9	Chlordane	1.0
8001-35-2	Toxaphene	5.0

CAS	PCB	NQL
12674-11-2	Atoclor-1016	1.0
1104-28-2	Atoclor-1221	1.0
11141-16-5	Atoclor-1232	1.0
53469-21-9	Atoclor-1242	1.0
12672-29-6	Atoclor-1248	1.0
11097-69-1	Atoclor-1254	1.0
11096-82-5	Atoclor-1260	1.0

The "Nominal Quantitation Limit" listed for each target compound is based on the Superfund CLP Protocol. The Actual Quantitation Limits are related to the NQLs by the NQL Factor. This NQL factor reflects procedural steps, e.g., extract dilution, which influence quantitation limits.

Office of Analytical Services and Quality Assurance
Pesticide and PCB Analysis
Nominal Quantitation Limits (NQL)

Units: Solids = mg/kg NPTC = Non-Priority Pollutant Target Compound

Actual Quantitation Limit = (NQL Factor) X NQL

CAS	Pesticide	NQL
319-84-6	Alpha-BHC	0.002
319-85-7	Beta-BHC	0.002
319-86-8	Delta-BHC	0.002
58-89-8	Gamma-BHC	0.002
76-44-8	Heptachlor	0.002
309-00-2	Aldrin	0.002
1024-57-3	Heptachlor Epoxide	0.002
959-98-8	Endosulfan I	0.002
60-57-1	Dieldrin	0.003
72-55-9	4,4'-DDE	0.003
72-20-8	Endrin	0.003
33213-65-9	Endosulfan II	0.003
72-54-8	4,4'-DDD	0.003
1031-07-8	Endosulfan Sulfate	0.003
50-29-3	4,4'-DDT	0.003
7421-93-4	Endrin Aldehyde	0.003
53494-70-5	Endrin Ketone	NPTC 0.003
72-43-5	Methoxychlor	NPTC 0.017
5103-71-9	Alpha-Chlordane	0.002
5103-74-2	Gamma-Chlordane	0.002
57-74-9	Chlordane	0.033
8001-35-2	Toxaphene	0.167

CAS	PCB	NQL
12674-11-2	Aroclor-1016	0.033
1104-28-2	Aroclor-1221	0.033
11141-16-5	Aroclor-1232	0.033
53469-21-9	Aroclor-1242	0.033
12672-29-6	Aroclor-1248	0.033
11097-69-1	Aroclor-1254	0.033
11096-82-5	Aroclor-1260	0.033

The "Nominal Quantitation Limit" listed for each target compound is based on the Superfund CLP Protocol. The Actual Quantitation Limits are related to the NQLs by the NQL Factor. This NQL factor reflects procedural steps, e.g., extract dilution, which influence quantitation limits.

Office of Analytical Services and Quality Assurance
Extractable Organics Analysis
Nominal Quantitation Limits (NQL)

Units: Water = ug/L NPTC = Non-Priority Pollutant Target Compound

Actual Quantitation Limit = (NQL Factor) X NQL

CAS	ANALYTE	NQL
62-75-9	N-Nitrosodimethylamine	10
108-95-2	Phenol	10
62-53-3	Aniline	NPTC 10
111-44-4	Bis(2-Chloroethyl) Ether	10
95-57-8	2-Chlorophenol	10
541-73-1	1,3-Dichlorobenzene	10
106-46-7	1,4-Dichlorobenzene	10
100-51-6	Benzyl Alcohol	NPTC 10
95-50-1	1,2-Dichlorobenzene	10
95-48-7	2-Methylphenol	NPTC 10
108-60-1	Bis(2-chloroisopropyl) Ether	10
106-44-5	4-Methylphenol	NPTC 10
621-64-7	N-Nitroso-di-n-Propylamine	10
67-72-1	Hexachloroethane	10
98-95-3	Nitrobenzene	10
78-59-1	Isophorone	10
88-75-5	2-Nitrophenol	10
105-67-9	2,4-Dimethylphenol	10
65-85-0	Benzoic Acid	NPTC 50
111-91-1	Bis(2-Chloroethoxy) Methane	10
10-83-2	2,4-Dichlorophenol	10
120-82-1	1,2,4-Trichlorobenzene	10
91-20-3	Naphthalene	10
106-47-8	4-Chloroaniline	NPTC 10
87-68-3	Hexachlorobutadiene	10
59-50-7	4-Chloro-3-Methylphenol	10
91-57-6	2-Methylnaphthalene	NPTC 10
77-47-4	Hexachlorocyclopentadiene	10
88-06-2	2,4,6-Trichlorophenol	10
88-95-4	3,4,5-Trichlorophenol	NPTC 50
91-58-7	2-Chloronaphthalene	10
88-74-4	2-Nitroaniline	NPTC 50
131-11-3	Dimethylphthalate	10
208-96-8	Acenaphthylene	10

CAS	ANALYTE	NQL
99-09-2	3-Nitroaniline	NPTC 50
83-32-9	Acenaphthene	10
51-28-5	2,4-Dinitrophenol	50
100-02-7	4-Nitrophenol	50
132-64-9	Dibenzofuran	NPTC 10
606-20-2	2,6-Dinitrotoluene	10
121-14-2	2,6-Dinitrotoluene	10
84-66-2	Diethylphthalate	10
7005-72-3	4-Chlorophenylphenylether	10
86-73-7	Fluorene	10
100-01-6	4-Nitroaniline	NPTC 50
86-30-6	N-Nitrosodiphenylamine (1)	10
534-52-1	4,6-Dinitro-2-Methylphenol	50
101-55-3	4-Bromophenylphenylether	10
118-74-1	Hexachlorobenzene	10
87-86-5	Pentachlorophenol	50
85-01-8	Phenanthrene	10
120-12-7	Anthracene	10
86-74-8	Carbazole	NPTC 10
84-74-2	Di-n-Butylphthalate	10
206-44-0	Fluoranthene	10
92-87-5	Benzidine	50
129-00-0	Pyrene	10
85-68-7	Butylbenzylphthalate	10
91-94-1	3,3'-Dichlorobenzidine	20
56-55-3	Benzo(a)Anthracene	10
117-81-7	Bis(2-Ethylhexyl) Phthalate	10
218-01-9	Chrysene	10
117-84-0	Di-n-Octylphthalate	10
205-99-2	Benzo(b) Fluoranthene	10
207-08-9	Benzo(x) Fluoranthene	10
50-32-8	Benzo(a) Pyrene	10
193-39-5	Indeno(1,2,3-cd) Pyrene	10
53-70-3	Dibenzo(a,h) Anthracene	10
191-24-2	Benzo(g,h,i) Perylene	10

(1) N-nitrosodiphenylamine decomposes in the gas chromatographic inlet and cannot be separated from diphenylamine.

The "Nominal Quantitation Limit" factor is an overall correction factor applied to the method's NQLs for analytical adjustments made during the analysis (i.e., for extractions of more or less than the ideal 1000 ml for aqueous samples, for sample extracts not concentrated to 100 ml due to excessive foaming/darkness of the extract, and for sample extract dilutions prior to analyses).

Office of Analytical Services and Quality Assurance
Extractable Organics Analysis
Nominal Quantitation Limits (NQL)

Units: Solids = mg/kg (wet) NPTC = Non-Priority Pollutant Target Compound

Actual Quantitation Limit = (NQL Factor) X NQL

CAS	ANALYTE	NQL
62-75-9	N-Nitrosodimethylamine	0.33
108-95-2	Phenol	0.33
62-53-3	Aniline	NPTC 0.33
111-44-4	bis(2-Chloroethyl) Ether	0.33
95-57-8	2-Chlorophenol	0.33
541-73-1	1,3-Dichlorobenzene	0.33
106-46-7	1,4-Dichlorobenzene	0.33
100-51-6	Benzyl Alcohol	NPTC 0.33
95-50-1	1,2-Dichlorobenzene	0.33
95-48-7	2-Methylphenol	NPTC 0.33
108-60-1	bis(2-chloroisopropyl) Ether	0.33
106-44-5	4-Methylphenol	NPTC 0.33
621-64-7	N-Nitroso-di-n-Propylamine	0.33
67-72-1	Hexachloroethane	0.33
98-95-3	Nitrobenzene	0.33
78-59-1	Isophorone	0.33
88-75-5	2-Nitrophenol	0.33
105-67-9	2,4-Dimethylphenol	0.33
65-85-0	Benzoic Acid	NPTC 1.67
111-91-1	bis(2-Chloroethoxy)Methane	0.33
120-83-2	2,4-Dichlorophenol	0.33
120-82-1	1,2,4-Trichlorobenzene	0.33
91-20-3	Naphthalene	0.33
106-47-8	4-Chloroaniline	NPTC 0.33
87-68-3	Hexachlorobutadiene	0.33
59-50-7	4-Chloro-3-Methylphenol	0.33
91-57-6	2-Methylnaphthalene	NPTC 0.33
77-47-4	Hexachlorocyclopentadiene	0.33
88-06-2	2,4,6-Trichlorophenol	0.33
95-95-4	2,4,5-Trichlorophenol	NPTC 1.67
91-58-7	2-Chloronaphthalene	0.33
88-74-4	2-Nitroaniline	NPTC 1.67
131-11-3	Dimethylphthalate	0.33
208-96-8	Acenaphthylene	0.33

CAS	ANALYTE	NQL
99-09-2	3-Nitroaniline	NPTC 1.67
83-32-9	Acenaphthene	0.33
51-28-5	2,4-Dinitrophenol	1.67
100-02-7	4-Nitrophenol	1.67
132-64-9	Dibenzofuran	NPTC 0.33
606-20-2	2,6-Dinitrotoluene	0.33
121-14-2	2,6-Dinitrotoluene	0.33
84-66-2	Diethylphthalate	0.33
7005-72-3	4-Chlorophenylphenylether	0.33
86-73-7	Fluorene	0.33
100-01-6	4-Nitroaniline	NPTC 1.67
86-30-6	N-Nitrosodiphenylamine (1)	0.33
534-52-1	4,6-Dinitro-2-Methylphenol	1.67
101-55-3	4-Bromophenylphenylether	0.33
118-74-1	Hexachlorobenzene	0.33
87-86-5	Pentachlorophenol	1.67
85-01-8	Phenanthrene	0.33
120-12-7	Anthracene	0.33
86-74-8	Carbazole	NPTC 0.33
84-74-2	Di-n-Butylphthalate	0.33
206-44-0	Fluoranthene	0.33
92-87-5	Benzidine	1.67
129-00-0	Pyrene	0.33
85-68-7	Butylbenzylphthalate	0.33
91-94-1	3,3'-Dichlorobenzidine	0.67
56-55-3	Benzo(a)Anthracene	0.33
117-81-7	bis(2-Ethylhexyl) Phthalate	0.33
218-01-9	Chrysene	0.33
117-84-0	Di-n-Octylphthalate	0.33
205-99-2	Benzo(b)Fluoranthene	0.33
207-08-9	Benzo(k)Fluoranthene	0.33
50-32-8	Benzo(a)Pyrene	0.33
193-39-5	Indeno(1,2,3-cd)Pyrene	0.33
53-70-3	Dibenzo(a,h)Anthracene	0.33
191-24-2	Benzo(g,h,i)Perylene	0.33

(1) N-nitrosodiphenylamine decomposes in the gas chromatographic inlet and cannot be separated from diphenylamine.

The "Nominal Quantitation Limit" factor is an overall correction factor applied to the method's NQLs for analytical adjustments made during the analysis (i.e., for extractions of more or less than the ideal 30 grams for soil samples, for sample extracts not concentrated to 1.00 ml due to excessive foaming/darkness of the extract, and for sample extract dilutions prior to analyses).

For example, the typical NQL factor for an OASQA soil sample is 1.5. Therefore, the estimated Actual Quantitation Limit for phenol would be 0.50 mg/Kg (i.e., 1.5 x .33 mg/Kg).

Rev. October 16, 1996

AR100042

Volatile Organic Analysis by GC/MS

Analyst:

Douglas Gardner
LMES Chemist

Method:

Eight aqueous and six soil samples from the WILMINGTON DRUM site (REQ99016) were analyzed for the presence of volatile organic compounds amenable to purge and trap and identifiable by mass spectrometry. The samples were collected October 19, 1998 and analyzed October 22, 23, 26, 27, 1998 and November 9, 1998, using a consolidated method derived from EPA methods SDWA 524.2, NPDES 624, RCRA 8260 and the Superfund CLP Statement of Work.

All samples were initially analyzed within the 14 day technical holding time, except sample 981021-07. Due to a sample tracking error, this sample was analyzed 7 days beyond the holding time. No target analytes were detected in this sample and no data were qualified for this sample.

For the aqueous samples analyzed using the initial calibration on October 15, 1998, the nominal quantitation limit (NQL) for each compound is as listed in the attached table, except the NQLs for dichlorodifluoromethane, chloromethane, vinyl chloride, chloroethane, 4-methyl-2-pentanone, 2-hexanone, 1,2-dibromo-3-chloropropane, naphthalene, and 1,2,3-trichlorobenzene are 10. In addition, 2-chloroethylvinyl ether was not amenable to purge and trap and was not identifiable by mass spectrometry. For the aqueous samples analyzed using the initial calibration on November 9, 1998, the NQL for each compound is as listed in the attached table, except the NQLs for acetone, 4-methyl-2-pentanone, and 2-hexanone are 10 and the NQLs for vinyl acetate, 2 butanone and 2-chloroethylvinyl ether are 50.

For the soil samples, the NQL for each compound is as listed in the attached table, except the NQLs for dichlorodifluoromethane, chloromethane, vinyl chloride, chloroethane, and 2,2-dichloropropane are 10 and the NQL for 2-chloroethylvinyl ether is 50.

Only detected results are reported. The results for the soil samples are reported on a WET weight basis. No tentatively identified compounds (TICs) were detected in the samples.

AR100043

Tentatively Identified Compounds

Site: Wilmington Drum

Program: Superfund Removal/Remedial

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/K.g)
981021-01		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/K.g)
981021-02		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/K.g)
981021-03		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/K.g)
981021-04		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/K.g)
981021-05		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/K.g)
981021-06		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/L)
981021-07		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/L)
981021-08		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/L)
981021-09		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/L)
981021-10		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/L)
981021-11		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC (ug/L)
981021-12		None Detected		

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Sample No.	CAS#	COMPOUND NAME	R.T.	CONC.(ug/L)
981021-13		None Detected		

Sample No.	CAS#	COMPOUND NAME	R.T.	CONC.(ug/L)
981021-14		None Detected		

Pesticide/PCB Analysis By Gas Chromatography

Analyst:

Timothy M. Casey
LMES Chemist

Method:

Thirteen mixed matrix samples from WILMINGTON DRUM site (REQ99016) were analyzed for polychlorinated biphenyls and/or pesticides by capillary column gas chromatography. All extractions and analyses were performed according to SOP R3-QA207.0. This SOP is a consolidated method derived from EPA methods SDWA 508, NPDES 608, RCRA 8081 & 8082, and the Superfund CLP Statement of Work. The samples were collected October 19, 1998; extracted October 22 and 26, 1998; and analyzed November 3, 4, 5, and 6, 1998. All soil results are reported on a WET WEIGHT basis.

Quality Control:

Six of the twenty-six surrogate recoveries were outside the acceptance limit criteria (60-150% recovery). One of the twenty-four matrix spike recoveries was outside the acceptance limits and seven of the twelve matrix spike relative percent difference (RPD) values were outside the acceptance limits. These values have been flagged with an "A".

Continuing calibration criteria were monitored for target pesticides. All sample results were bracketed by valid calibration check standards.

Due to the complex nature of the sample matrix, non-target interference peaks may be eluting within pesticide retention time windows. Target analyte results with relative percent difference greater than 25% between the two analytical columns may be considered suspect and have been flagged with an "R".

Due to the presence of polychlorinated biphenyls (PCBs) in several samples, PCB congener peaks may be eluting within pesticide retention time windows. Pesticide results attributed to PCB interference have been flagged with an "I".

The nominal quantitation limit factors for these samples are listed below.

Sample	Nominal Quantitation Limit Factor
981021-01	2.0 Pesticides 1.0 PCBs/Surrogates
981021-02	2.0 Pesticides 1.0 PCBs/Surrogates
981021-03	5.0 Pesticides 1.0 PCBs/Surrogates
981021-04	2.0 Pesticides 1.0 PCBs/Surrogates
981021-05	5.0 Pesticides 1.0 PCBs/Surrogates
981021-06	2.0 Pesticides 1.0 PCBs/Surrogates
981021-08	1.0 All Compounds
981021-09	1.0 All Compounds
981021-10	1.0 All Compounds
981021-11	1.0 All Compounds
981021-12	1.0 All Compounds
981021-13	1.0 All Compounds
981021-14	1.0 All Compounds

Semi-volatile Analysis by GC/MS

Analyst:

Hoang Nguyen
LMES Chemist

Method:

Six soil and seven aqueous samples from WILMINGTON DRUM site (REQ99016) were analyzed for the presence of semi-volatile organic compounds amenable to soxhlet and continuous liquid/liquid extraction and identifiable by mass spectrometry. The samples were collected October 19, 1998 and were extracted October 22 and 26, 1998. These samples were analyzed October 29, November 02 and 03, 1998 by a consolidated method derived from EPA methods SDWA 525.2, NPDES 625, RCRA 8270 and the Superfund CLP Statement of Work.

Only detected results are reported. The results for soil samples are reported on a WET weight basis. The NQL factor for all samples is equal to 1. Tentatively identified compounds (TICs) have been corrected for any blank contamination. The TICs are reported below.

TENTATIVELY IDENTIFIED COMPOUNDS

Site: Wilmington Drum '
Program: Superfund Removal/Remedial

UNITS: mg/Kg

<u>SAMPLE NO.</u>	<u>CAS #</u>	<u>TIC NAME</u>	<u>RT</u>	<u>CONC.</u>
981021-01	*****	Unknown m/z = 43	05.94	0.3 T
	*****	Unknown m/z = 41	06.37	0.6 T
	*****	Unknown m/z = 85	08.92	2.3 T
	*****	Unknown m/z = 43	10.59	0.8 T
	*****	Unknown m/z = 43	12.04	0.4 T
	*****	Unknown m/z = 43	12.50	0.8 T
	*****	Unknown m/z = 55	24.78	0.8 T
	057103	Hexadecanoic acid	24.95	0.5 T
	*****	Unknown m/z = 79	28.12	0.3 T

AR100048

UNITS: mg/Kg

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-02	*****	Unknown m/z = 85	08.91	2.5 T
	*****	Unknown m/z = 43	10.58	0.5 T
	*****	Unknown m/z = 43	12.43	0.4 T
	*****	Unknown m/z = 55	24.75	0.3 T
	057103	Hexadecanoic acid	24.95	0.6 T
	057114	Octadecanoic acid	26.81	0.3 T
	*****	Unknown m/z = 69	32.06	0.3 T
	*****	Unknown alkane m/z = 57	32.48	0.6 T
	*****	Unknown alkane m/z = 57	32.97	0.4 T
	*****	Unknown m/z = 59	34.05	0.3 T
	*****	Unknown m/z = 59	35.90	0.2 T
	*****	Unknown m/z = 81	36.22	0.2 T

UNITS: mg/Kg

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-03	*****	Unknown m/z = 43	05.92	0.4 T
	*****	Unknown m/z = 85	08.92	2.4 T
	*****	Unknown m/z = 43	10.59	0.4 T
	*****	Unknown m/z = 43	12.53	0.7 T
	*****	Unknown m/z = 55	24.79	0.4 T
	*****	Unknown m/z = 43	31.35	0.5 T
	*****	Unknown m/z = 43	31.66	0.6 T
	*****	Unknown alkane m/z = 57	32.51	1.0 T
	*****	Unknown m/z = 57	32.97	0.6 T
	*****	Unknown m/z = 59	35.95	0.4 T

UNITS: mg/Kg

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-04	*****	Unknown m/z = 43	05.90	0.3 T
	*****	Unknown m/z = 85	08.92	2.7 T
	*****	Unknown m/z = 43	10.59	0.6 T
	*****	Unknown m/z = 43	12.53	0.6 T
	*****	Unknown m/z = 55	24.79	0.5 T
	057103	Hexadecanoic acid	24.97	0.4 T
	*****	Unknown alkane m/z = 57	32.51	0.8 T
	*****	Unknown m/z = 95	36.27	0.4 T

UNITS: mg/Kg

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-05	*****	Unknown m/z = 85	08.91	2.0 T
	*****	Unknown m/z = 43	12.49	0.4 T
	*****	Unknown m/z = 55	24.80	0.5 T
	057103	Hexadecanoic acid	25.01	0.5 T
	*****	Unknown alkane m/z = 57	31.13	0.4 T
	*****	Unknown alkane m/z = 57	31.85	0.5 T
	*****	Unknown alkane m/z = 57	32.54	1.8 T
	*****	Unknown m/z = 191	32.99	1.1 T
	*****	Unknown m/z = 59	33.35	0.4 T
	*****	Unknown m/z = 43	33.51	0.6 T
	*****	Unknown m/z = 191	34.65	0.5 T
	*****	Unknown m/z = 59	35.98	0.8 T

UNITS: mg/Kg

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-06	*****	Unknown m/z = 85	08.91	2.0 T
	*****	Unknown m/z = 43	10.58	0.3 T
	*****	Unknown m/z = 43	12.42	0.3 T
	057103	Hexadecanoic acid	24.96	0.3 T
	*****	Unknown alkane m/z = 57	31.10	0.2 T
	*****	Unknown m/z = 59	31.58	0.2 T
	*****	Unknown alkane m/z = 57	32.50	0.5 T
	538238	Glycerol tricaprylate	32.85	0.4 T
	*****	Unknown m/z = 59	35.94	0.2 T
	*****	Unknown m/z = 81	36.29	0.3 T

UNITS: ug/L

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-08		None Detected		

UNITS: ug/L

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-09	*****	Unknown m/z = 69	05.27	8 T
	*****	Unknown m/z = 41	06.16	6 T
	*****	Unknown m/z = 109	09.25	1 T
	134623	Diethyltoluamide	20.93	1 T
	*****	Unknown m/z = 55	24.77	1 T
	*****	Unknown m/z = 79	25.88	1 T

UNITS: ug/L

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-10	*****	Unknown m/z = 69	05.29	4 T
	*****	Unknown m/z = 41	06.18	3 T

UNITS: ug/L

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-11	*****	Unknown m/z = 69	05.22	3 T
	*****	Unknown m/z = 45	05.43	2 T
	*****	Unknown m/z = 41	06.11	2 T

UNITS: ug/L

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-12	*****	Unknown m/z = 69	05.23	12 T
	*****	Unknown m/z = 45	05.44	2 T
	*****	Unknown m/z = 41	06.11	11 T

UNITS: ug/L

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-13	*****	Unknown m/z = 69	05.26	12 T
	*****	Unknown m/z = 41	06.13	10 T
	*****	Unknown m/z = 79	25.85	1 T

UNITS: ug/L

SAMPLE NO.	CAS #	TIC NAME	RT	CONC.
981021-14	*****	Unknown m/z = 69	05.23	10 T
	*****	Unknown m/z = 45	05.44	2 T
	*****	Unknown m/z = 41	06.11	9 T

PROJ. NO.	PROJECT NAME	NO. OF CONTAINERS	REMARKS
4645	W/41419100 OTCM		
SAMPLERS (Signature)			
<i>Allen m. Brown</i>			

STA. NO.	DATE	TIME	COMP.	GRAB	STATION LOCATION	NO. OF CONTAINERS	REMARKS
SD-01	10/21/98	1540	X		SD-01 98102101	3	1-22234418 to 3-22234420
SD-02	11/00		X		SD-02 98102102	3	3-2215304 to 3-2215311
SD-03	11/30		X		SD-03 98102103	3	3-2215312 to 3-2215314
SD-06	11/00		X		SD-06 98102104	3	3-2215321 to 3-2215323
SD-05	10/00		X		SD-05 98102105	3	3-2215318 to 3-2215320
SD-04	12/30		X		SD-04 98102106	3	3-2215306 to 3-2215308
SD-04	12/30		X		SD-04 98102107	3	3-2215324 to 3-2215326
SD-04/10/48	12/30		X		SD-04 98102108	3	3-2215315 to 3-2215317
Temperature in water							

Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time
<i>Allen m. Brown</i>	10/21/98 1500		
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time

Relinquished by: (Signature)	Date / Time	Received for Laboratory by: (Signature)	Date / Time
		<i>Allen m. Brown</i>	10/21/98 7:05

Distribution: Original Accompanies Shipment; Copy to Coordinator Field Files
 Remarks: AIR 0111 # 80865624903

Office of Enforcement

REGION 3
841 Chestnut St.
Philadelphia, Pennsylvania 19107

Distribution: Original Accompanies Shipment, Copy to Coordinator Field Files

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Office of Enforcement

REGION 3
841 Chestnut St.
Philadelphia, Pennsylvania 19107

CHAIN OF CUSTODY RECORD

[illegible]

Distribution: Original Accompanies Shipment; Copy to Coordinator Field File.

AR 100054

CHAIN OF CUSTODY RECORD

RECEIVED
841 Chestnut St.
Philadelphia, Pennsylvania 19107

PROJ. NO. PROJECT NAME

4645

SAMPLERS (Signature)

Philadelphia Area

NO.

OF

CON
TAINERS

REMARKS

STA NO.	DATE	TIME	COMP.	GRAB	STATION LOCATION
Sk-01	10/1/80	1540	X		Sk-01 9810211540
Sk-02	11/00		X		Sk-02 9810211100
Sk-03	11/30		X		Sk-03 9810211130
Sk-05	10/1/80	1000	X		Sk-05 9810211000
Temperature Indicator					

NO.

UPA (HCL PH2O)
ROA (ROA) SUMA
MODULA (MOD) PALER
L-PA (L-PA) PH2O

T46 2/24/80-1

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Date / Time

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Date / Time

Received for Laboratory by: (Signature)

Date / Time

Remarks

AIR 511 IF

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Notice to Laboratory Personnel

Background

Under the authority of Section 104 of the Comprehensive Environmental Response, Compensation, and Liability Act (CERCLA) or Superfund of 1980, Section 311 of the Clean Water Act, and Section 1 of the Resource Conservation and Recovery Act (RCRA), EPA has been delegated the responsibility to undertake emergency actions with respect to the release or potential release of oil, petroleum, or hazardous substances that pose a substantial threat to human health or welfare, or the environment. In addition, EPA provides technical assistance to help anticipate management of the public health, welfare or environment during other emergencies and control situations.

EPA's successful implementation of these emergency response action responsibilities requires that technical support capabilities be provided in the form of on-site Technical Assistance Teams (TAT) for each EPA Region. The WESTON TAT Concept 61-370-0054 provides support to EPA Regions I, II, III, IV, ERT - Edison, and Headquarters - Washington, DC.

Hazard Communication

The samples which accompany this notice have been shipped to your laboratory for analysis in accordance with applicable D.O.T. or IATA Regulations and were collected by the WESTON TAT and were chemically designated by the field response team as either environmental or hazardous material samples.

In general, Environmental Samples are collected from streams, from ponds, small lakes, wells, and off-shore wells that are not immediately expected to be contaminated with hazardous materials. Samples of on-site soils or sludge, and materials collected from drums, bulk storage tanks, obviously contaminated ponds, impoundments, lagoons, pools, and basins from hazardous waste sites are considered Hazardous Samples. Samples which are obtained from a known radioactive material contamination site or which demonstrate beta or gamma activity greater than three times average background as counted with a Geiger-Mueller radiation survey meter are considered Radioactive Samples.

The samples which accompany this notice have been chemically classified by the field response team as:

☒ Environmental ☐ Hazardous ☐ Corrosive (Environ. & Hazardous) ☐ Radioactive

The field team which collected the samples used the following Level(s) of personal protection as designated by EPA and OSHA, precautions to provide protection against possible radiological or chemical exposure:

☐ Level A ☐ Level B ☐ Level C ☒ Level D

This information is intended for use as a guide for the safe handling of these laboratory samples in accordance with EPA and OSHA regulations. The sample classification(s) and Level(s) of personal protection used by the WESTON TAT are not represented to be, nor are they adequate or applicable in all situations, nor are they intended to serve as a substitute for professional/personal judgement.

This form was prepared by: Karin B. R. 11/18/98

Analytical Services TID No. 41111111111111111111 Date 10/20/98

WESTON Office Dallas, TX Phone 800-461-4003 FAX 800-461-4114

Laboratory Name USEPA Region III LAB

AR100056

FedEx USA Airbill

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