

## Carbon Sampling Results

Note:

- Bay West lab report 5-3567 and Aptus/Triangle Labs project P012744 include carbon sample results from 1993 that are not included in **Table 3-2**. Analytical results from the most-recent carbon sampling event more accurately represent analyte concentrations under current operating conditions.

June 7, 1993

Mr. Maurice McClish  
Strebor, Inc.  
2305 Superior Avenue  
Kalamazoo, MI 49001

RE: Spent Activated Carbon Chemical Analyses

Dear Mr. McClish:

Enclosed is a copy of the analytical report for the spent activated carbon samples collected on April 22, 1993. The Total 2,3,7,8-TCDD Equivalents detected in the sample were approximately 58 parts per billion.

Please call me at 1-800-279-0456 if you have any questions regarding the above.

Sincerely,



Martin W. Wangensteen  
Project Manager

enclosure

c: David Cosgriff, Champion (Libby, Montana)  
Philip Grashoff, Jr., Honigman Miller Schwartz & Cohn  
Melinda Kemp, Champion (Stamford, Connecticut)  
Robert Morse, SmithKline Beecham

June 2, 1993

Bay West Environmental Services  
5 Empire Drive  
St. Paul, MN 55103

Attn: Mr. Martin Wangensteen

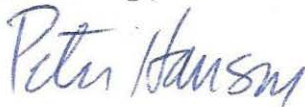
Bay West Environmental Services Project No.: 7095  
(COC: SI-287)  
Bay West Laboratory Project ID: 5-3567  
Samples Collected: April 22, 1993

The following are results from the samples you submitted for analysis on April 23, 1993. As requested, the 4 samples submitted were composited in equal portions by the laboratory and analyzed.

The data from Bay West Laboratory is reported in Tables 1 through 3, followed by the data from Twin City Testing.

Please contact me if you have any questions or comments.

Sincerely,



Peter Hanson  
Laboratory Manager

PH/ly

encl.

Table 1

Bay West Environmental Services Project No.: **7095**  
 Bay West Laboratory Project ID: **5-3567**

**Purgeable Halogenated and Non-Halogenated Compounds**  
**8010/8020 Target Compound List**

<u>Compound Name</u>	<u>Quantitation Limit</u> mg/Kg	<u>Composite Sample</u> (Items 1-4) (35681) mg/Kg
Benzene	0.050	0.055
Bromodichloromethane	0.050	ND
Bromoform	0.050	ND
Bromomethane	0.050	ND
Carbon Tetrachloride	0.050	ND
Chlorobenzene	0.050	ND
Chloroethane	0.050	ND
Chloroform	0.050	ND
Chloromethane	0.050	ND
Dibromochloromethane	0.050	ND
Dichlorobenzene, 1,2-	0.050	ND
Dichlorobenzene, 1,3-	0.050	ND
Dichlorobenzene, 1,4-	0.050	ND
Dichlorodifluoromethane	0.050	ND
Dichloroethane, 1,1-	0.050	ND
Dichloroethane, 1,2-	0.050	ND
Dichloroethene, 1,1-	0.050	ND
Dichloroethene, t-1,2-	0.050	ND
Dichloropropane, 1,2-	0.050	ND
Dichloropropene, c-1,3-	0.050	ND
Dichloropropene, t-1,3-	0.050	ND
Ethylbenzene	0.050	ND
Methylene Chloride	0.050	ND
Tetrachloroethane, 1,1,2,2-	0.050	ND
Tetrachloroethene	0.050	ND
Trichloroethane, 1,1,1-	0.050	ND
Trichloroethane, 1,1,2-	0.050	ND
Trichloroethene	0.050	ND
Trichlorofluoromethane	0.050	ND
Toluene	0.050	1.2
Vinyl Chloride	0.050	ND
Styrene	0.050	ND
Xylenes	0.050	7.8

Analyzed: May 2 & 3, 1993  
 Method: EPA 5030/8010/8020 Modified

ND = Not Detected, concentration less than Quantitation Limit.

Table 2

Bay West Environmental Services Project No: **7095**  
Bay West Laboratory Project ID: **5-3567**

Pentachlorophenol

Parameter	Quantitation Limit mg/Kg	Composite Sample (Items 1-4) (35683) mg/Kg
Pentachlorophenol	0.033	150
Dilution Factor		100

Date Extracted: May 5, 1993  
Date Analyzed: May 11, 1993

Method: EPA 604 Modified

Table 3

Bay West Environmental Services Project No.: **7095**  
Bay West Laboratory Project ID: **5-3567**

## Phthalates

Parameter	Quantitation Limit mg/Kg	Composite Sample (Items 1-4) (35683) mg/Kg
Di-Methyl Phthalate	0.033	0.044
Di-Ethyl Phthalate	0.033	ND
Di-N-Butyl Phthalate	0.033	ND
Butyl Benzyl Phthalate	0.033	ND
Bis (2-Ethylhexyl) Phthalate	0.033	ND
Di-N-Octyl Phthalate	0.033	1.0

Date Extracted: May 5, 1993  
Date Analyzed: May 13, 1993

Method: EPA 606 Modified

ND = Not Detected, concentration less than Quantitation  
Limit.



662 CROMWELL AVENUE  
ST. PAUL, MN 55114  
PHONE 612/645-3601

**REPORT OF: CHEMICAL ANALYSES**

**PROJECT:** BAY WEST LABS, 16824, PRJ #3567

**DATE:** May 20, 1993

**REPORTED TO:** Bay West, Inc.  
Attn: Peter Hanson  
Five Empire Drive  
St. Paul, MN 55103-1867

**LABORATORY NO:** 4410 93-1556

**INTRODUCTION**

This report presents the results of the analyses of one sample received on April 26, 1993, from a representative of Bay West, Inc. The scope of our services was limited to the parameters listed in the attached tables.

**METHODOLOGY**

Analyses are performed according to Twin City Testing Standard Operating Procedures. The procedures are based on the references stated in the analytical results tables.

**RESULTS**

The results are listed in the attached tables.

**DISCUSSION**

The recoveries of the isotopically-labeled PCDD/PCDF internal standards in the sample generally ranged from 52-95%, indicating a level of efficiency through the extraction and enrichment steps that is considered typical for this sample matrix type. Since the quantifications of the native 2,3,7,8-substituted isomers are based on isotope dilution, the data are automatically corrected for variations in recovery and accurate values are obtained.

The sample was found to contain high levels of selected isomers from the hexa, hepta, and octa-chlorinated congener classes which saturated the detector signal of the mass spectrometer, even after dilution of the final sample extract. The affected values have been flagged on the results table and should be regarded as minimum possible concentrations.

A laboratory method blank was prepared and analyzed with the PCDD/PCDF sample extraction batch as part of our routine quality control procedures. The results show the blank to be free of PCDDs and PCDFs, with the exceptions of trace background levels of HxCDF (3 picograms) and OCDD (26 picograms). It should be noted that these levels were below the method calibration ranges. The levels reported for these isomers in the actual sample were two or more orders of magnitude higher than the corresponding blank levels, indicating that the sample processing steps did not contribute significantly to the levels reported for the sample.

**PROJECT:** BAY WEST LABS, 16824, PRJ #3567

**DATE:** May 20, 1993

**LABORATORY NO:** 4410 93-1556


**PAGE:** 2

A laboratory quality control PCDD/PCDF spike sample was also prepared with the sample batch by extracting clean sand that had been fortified with native standard materials. The results show that the spiked native compounds were recovered at levels ranging from 70-100%, indicating a high degree of accuracy for these determinations.

**REMARKS**

The sample was collected on April 22, 1993. If the sample is not consumed in the analysis, it is held for two months from the date of sample receipt and then disposed, unless written instructions to the contrary are received.

**TWIN CITY TESTING CORPORATION**

  
Deneen Walker  
Project Manager

Stephanie Kidder  
Laboratory Manager

DW/SK/jd



**POLYNUCLEAR AROMATIC HYDROCARBON RESULTS**  
**EPA METHOD 8310**

(All values are in  $\mu\text{g}/\text{Kg}$  which is equal to parts-per-billion)

**Client ID:** Blank 3567<sup>1</sup>

**TCT ID:** 315649

<u>Parameter:</u>			<u>PQL</u>
Naphthalene	ND	14,000	360
1-methylnaphthalene	ND	8,900	460
2-methylnaphthalene	ND	9,400	460
Acenaphthylene	ND	4,100	460
Acenaphthene	ND	ND	360
Fluorene	ND	83	42
Phenanthrene	ND	760	130
Anthracene	ND	ND	140
Fluoranthene	ND	ND	42
Pyrene	ND	130	20
Benzo (a) anthracene	ND	ND	2.6
Chrysene	ND	ND	30
Benzo (b) fluoranthene	ND	ND	3.6
Benzo (k) fluoranthene	ND	ND	3.4
Benzo (a) pyrene	ND	ND	4.6
Dibenzo (a,h) anthracene	ND	ND	6.0
Benzo (ghi) perylene	ND	ND	15
Indeno (1,2,3 cd) pyrene	ND	ND	8.8
% Surrogate Recovery	79%	0% <sup>2</sup>	
<b>Date Extracted:</b>	4/28/93	4/28/93	
<b>Date Analyzed:</b>	5/1/93	5/1/93	

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

<sup>1</sup> The sample chromatogram contained many unresolved peaks. It is likely that in addition to the listed PAHs, other PAHs or similar compounds were also present. Since the quantitative values reported here represent the total contribution from all compounds which gave a positive response at the column retention time of a particular PAH, the actual levels may be lower than the reported values. Confirmation of the levels of specific PAHs will require an alternative analytical technique which can isolate and quantify the separate components of the mixture.

<sup>2</sup> No surrogate recovered.

**Reference:** EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.

**POLYNUCLEAR AROMATIC HYDROCARBON RESULTS**  
**EPA METHOD 8310**

(All values are in  $\mu\text{g}/\text{Kg}$  which is equal to parts-per-billion)

**Client ID:** 3567<sup>1</sup>

**TCT ID:** 315649 Re-Extraction

<u>Parameter:</u>		<u>PQL</u>
Naphthalene	40,000	360
1-methylnaphthalene	24,000	460
2-methylnaphthalene	3,100	460
Acenaphthylene	6,600	460
Acenaphthene	ND	360
Fluorene	280	42
Phenanthrene	1,300	130
Anthracene	ND	140
Fluoranthene	ND	42
Pyrene	460	20
Benzo (a) anthracene	30	2.6
Chrysene	40	30
Benzo (b) fluoranthene	ND	3.6
Benzo (k) fluoranthene	ND	3.4
Benzo (a) pyrene	ND	4.6
Dibenzo (a,h) anthracene	ND	6.0
Benzo (ghi) perylene	ND	15
Indeno (1,2,3 cd) pyrene	ND	8.8
% Surrogate Recovery	0% <sup>2</sup>	
<b>Date Extracted:</b>	5/5/93	
<b>Date Analyzed:</b>	5/6/93	

All results are reported on a dry weight basis.

PQL = Practical Quantitation Limit

ND = Not Detected

<sup>1</sup> The sample chromatogram contained many unresolved peaks. It is likely that in addition to the listed PAHs, other PAHs or similar compounds were also present. Since the quantitative values reported here represent the total contribution from all compounds which gave a positive response at the column retention time of a particular PAH, the actual levels may be lower than the reported values. Confirmation of the levels of specific PAHs will require an alternative analytical technique which can isolate and quantify the separate components of the mixture.

<sup>2</sup> No surrogate recovery caused by matrix effects.

**Reference:** EPA Test Methods for Evaluating Solid Waste, SW-846, November 1986, 3rd Edition.

\*\*\*\*\*  
 \*TWIN CITY TESTING CORPORATION\*  
 \*METHOD 8290 ANALYSIS RESULTS \*  
 \*\*\*\*\*  
 Client....BAYWEST

TCT Sample ID.....BLANK-511A  
 Analysis Date.....5/17/93 16:27  
 Filename.....V30517E  
 Injected By.....MCH  
 Total Amount Extracted...0.0102 kg  
 % Moisture.....NA %  
 ICAL Date.....05/01/93  
 CCAL Filename.....V30517B  
 Method Blank ID.....NA  
 Extraction Date.....5/11/93

NATIVE ISOMERS	CONC. ng/kg	LOD ng/kg	INTERNAL STANDARDS	ng's ADDED	PERCENT RECOVERY
2378-TCDF	ND	0.31	2378-TCDF-13C....	2.00	90
TOTAL TCDF	ND	-----	2378-TCDD-13C....	2.00	71
			12378-PeCDF-13C..	2.00	81
2378-TCDD	ND	0.56	23478-PeCDF-13C..	2.00	99
TOTAL TCDD	ND	-----	12378-PeCDD-13C..	2.00	76
			123478-HxCDF-13C.	2.00	113
12378-PeCDF	ND	0.14	123678-HxCDF-13C.	2.00	100
23478-PeCDF	ND	0.12	234678-HxCDF-13C.	2.00	99
TOTAL PeCDF	ND	-----	123789-HxCDF-13C.	2.00	81
			123478-HxCDD-13C.	2.00	83
12378-PeCDD	ND	0.24	123678-HxCDD-13C.	2.00	76
TOTAL PeCDD	ND	-----	1234678-HpCDF-13C	2.00	69
			1234789-HpCDF-13C	2.00	68
123478-HxCDF	ND	0.18	1234678-HpCDD-13C	2.00	46
123678-HxCDF	ND	0.14	OCDD-13C.....	4.00	50
234678-HxCDF	0.32	-----			
123789-HxCDF	ND	0.20	1234-TCDD-13C....	2.00	NA
TOTAL HxCDF	0.32	-----	123789-HxCDD-13C.	2.00	NA
123478-HxCDD	ND	0.31	2378-TCDD-37C14..	0.20	61
123678-HxCDD	ND	0.68			
123789-HxCDD	ND	0.35			
TOTAL HxCDD	ND	-----			
			Total 2378-TCDD		
1234678-HpCDF	ND	0.27	Equivalence:	0.035	ng/kg
1234789-HpCDF	ND	0.38	(Using ITE Factors/DB-5	Data)	
TOTAL HpCDF	ND	-----			
1234678-HpCDD	ND	0.73			
TOTAL HpCDD	ND	-----			
OCDF	ND	0.60			
OCDD	2.60	-----			

All values are expressed on a total (as received) weight basis.

CONC= Concentration (Totals include 2378-substituted isomers.)  
 LOD = Limit of Detection  
 ND = Not Detected  
 NA = Not Applicable

TCT Invoice Number....4410 93-1556

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 \*TWIN CITY TESTING CORPORATION\*  
 \*METHOD 8290 ANALYSIS RESULTS \*  
 \*\*\*\*\*  
 Client....BAYWEST

Client's Sample ID.....3567  
 TCT Sample ID.....315649  
 Analysis Date.....5/17/93 20:25  
 Filename.....V30517I  
 Injected By.....MCH  
 Total Amount Extracted...0.0011 kg  
 % Moisture..... 30.3 %  
 ICAL Date.....05/01/93  
 CCAL Filename.....V30517B  
 Method Blank ID.....BLANK-511A  
 Extraction Date.....5/11/93

NATIVE ISOMERS	CONC. ng/kg	LOD ng/kg	INTERNAL STANDARDS	ng's ADDED	PERCENT RECOVERY
2378-TCDF	270 *	-----	2378-TCDF-13C....	2.00	78
TOTAL TCDF	4300	-----	2378-TCDD-13C....	2.00	74
2378-TCDD	41	-----	12378-PeCDF-13C..	2.00	64
TOTAL TCDD	1700	-----	23478-PeCDF-13C..	2.00	83
12378-PeCDF	890	-----	12378-PeCDD-13C..	2.00	35
23478-PeCDF	1100	-----	123478-HxCDF-13C.	2.00	52
TOTAL PeCDF	12000	-----	123678-HxCDF-13C.	2.00	95
12378-PeCDD	1800	-----	234678-HxCDF-13C.	2.00	86
TOTAL PeCDD	59000	-----	123789-HxCDF-13C.	2.00	65
123478-HxCDF	4700	-----	123478-HxCDD-13C.	2.00	81
123678-HxCDF	1900	-----	123678-HxCDD-13C.	2.00	78
234678-HxCDF	2000	-----	1234678-HpCDF-13C	2.00	69
123789-HxCDF	1500	-----	1234789-HpCDF-13C	2.00	66
TOTAL HxCDF	170000	-----	1234678-HpCDD-13C	2.00	56
123478-HxCDD	130000	-----	OCDD-13C.....	4.00	63
123678-HxCDD	130000	-----	1234-TCDD-13C....	2.00	NA
123789-HxCDD	150000	-----	123789-HxCDD-13C.	2.00	NA
TOTAL HxCDD	1200000 **	-----	2378-TCDD-37C14..	0.20	69
1234678-HpCDF	68000	-----	Total 2378-TCDD		
1234789-HpCDF	7500	-----	Equivalence:	58248 ng/kg	
TOTAL HpCDF	440000 **	-----	(Using ITE Factors/DB-5 Data)		
1234678-HpCDD	1100000 **	-----			
TOTAL HpCDD	2200000 **	-----			
OCDF	520000	-----			
OCDD	2400000 **	-----			

\* Value may include contributions from other TCDF isomers.  
 \*\*Saturated signal.  
 All values are expressed on a dry weight basis.

CONC= Concentration (Totals include 2378-substituted isomers.)  
 LOD = Limit of Detection  
 ND = Not Detected  
 NA = Not Applicable

TCT Invoice Number....4410 93-1556

\*\*\*\*\*  
 \*TWIN CITY TESTING CORPORATION\*  
 \* PCDD/PCDF SPIKE RESULTS \*  
 \*\*\*\*\*  
 Client....BAYWEST

TCT Sample ID.....SPIKE-511A  
 Analysis Date.....5/18/93 11:20  
 Filename.....V30518C  
 Sample Injected By.....CS  
 Total Amount Extracted.....0.0102 kg  
 % Moisture.....NA  
 ICAL Date.....05/01/93  
 CCAL Filename.....V30518B  
 Method Blank ID.....BLANK-511A  
 Extraction Date.....5/11/93

NATIVE ISOMERS	Qs (ng)	Qm (ng)	% REC	INTERNAL STANDARD	ng's ADDED	PERCENT RECOVERY
2378-TCDF	0.20	0.15	77	2378-TCDF-13C....	2.00	69
TOTAL TCDF	0.20	0.15	77	2378-TCDD-13C....	2.00	58
				12378-PeCDF-13C..	2.00	62
2378-TCDD	0.20	0.17	85	23478-PeCDF-13C..	2.00	79
TOTAL TCDD	0.20	0.17	85	12378-PeCDD-13C..	2.00	63
				123478-HxCDF-13C.	2.00	86
12378-PeCDF	1.00	0.81	81	123678-HxCDF-13C.	2.00	84
23478-PeCDF	1.00	0.73	73	234678-HxCDF-13C.	2.00	81
TOTAL PeCDF	2.00	1.53	77	123789-HxCDF-13C.	2.00	73
				123478-HxCDD-13C.	2.00	69
12378-PeCDD	1.00	0.81	81	123678-HxCDD-13C.	2.00	62
TOTAL PeCDD	1.00	0.81	81	1234678-HpCDF-13C	2.00	56
				1234789-HpCDF-13C	2.00	56
123478-HxCDF	1.00	0.75	75	1234678-HpCDD-13C	2.00	39
123678-HxCDF	1.00	0.70	70	OCDD-13C.....	4.00	40
234678-HxCDF	1.00	0.74	74			
123789-HxCDF	1.00	0.83	83	1234-TCDD-13C....	2.00	NA
TOTAL HxCDF	4.00	3.02	75	123789-HxCDD-13C.	2.00	NA
123478-HxCDD	1.00	0.86	86	2378-TCDD-37C14..	0.20	46
123678-HxCDD	1.00	0.86	86			
123789-HxCDD	1.00	0.88	88			
TOTAL HxCDD	3.00	2.60	87			
1234678-HpCDF	1.00	0.72	72			
1234789-HpCDF	1.00	0.71	71			
TOTAL HpCDF	2.00	1.42	71			
1234678-HpCDD	1.00	1.00	100			
TOTAL HpCDD	1.00	1.00	100			
OCDF	2.00	1.79	90			
OCDD	2.00	1.57	79			

Qs = Quantity Spiked  
 Qm = Quantity Measured  
 REC = Recovery (Expressed as Percent)  
 NA = Not Applicable

TCT Invoice Number....4410 93-1556

# GROUND WATER CHAIN-OF-CUSTODY RECORD FOR STREBOR, INC.

3507 BW-ST: 3/91

	LAB: <b>Bay West</b>	SEND RESULTS TO: <b>Martin Wangensteen</b>	CHAIN-OF-CUSTODY NO:
	PROJECT NUMBER <b>7095-</b>	TURNAROUND REQUEST <b>STANDARD</b>	<b>SI- 287</b>

ITEM NO.	SAMPLE NUMBER (PROJECT NO. - SAMPLE ID)	TIME	MATRIX	NUMBER & TYPE OF CONTAINER	ANALYSIS CODE(S)	COMMENTS/BAR CODE	ANALYSIS CODES	
1	7095 - 1B	11:45 A		2x40ml 3x500ml	15-1	Sample from Box # 2	EPA 601/602 (include Xylenes and Styrene)	01
2	7095 - 2	11:00 A		↓ ↓	15-1	↓ ↓ # 41	EPA 604 Phenols	02
3	7095 - 3	11:15 A		↓ ↓	15-1	↓ ↓ # 66	EPA 604 Pentachlorophenol	03
4	7095 - 4	12:30 P		2x40mL 2x500mL	15-1	↓ ↓ # 52	Total Suspended Solids	04
5	7095 -						EPA 610 PAHs	05
6	7095 -						EPA 606 Phthalate Esters	06
7	7095 -						EPA 607 Nitrosamines	07
8	7095 -						EPA 609 Nitroaromatics and Isophorone	08
9	7095 -						EPA 611 Haloethers	09
10	7095 -						EPA 612 Chlorinated Hydrocarbons	10
							CDDs/CDFs (HiRes GC/MS)	11
							Tri-butyl tin oxide (TBTO)	12
							PCBs	13
							Cd, Cr, Cu, Pb, Ni, Zn, Hg, total CN	14
							Other	15

Packed in polystyrene "peanuts"

PERSON RESPONSIBLE FOR SAMPLE COLLECTION <b>Tim Lindgren</b>		AFFILIATION <b>BAY WEST</b>		TRANS NO.	ITEM NO.	RELINQUISHED BY	ACCEPTED BY	DATE	TIME
DATE	<b>4-22-93</b>	TIME	<b>5:00 pm</b>	1	1-4	<b>Tim Lindgren</b>	<b>John Paulson</b>	<b>4-23-93</b>	<b>2:55</b>

ADDITIONAL ANALYSIS (USE BACK OF FRONT SHEET IF NECESSARY)									
15-1.	Composite samples and ANALYZE for the following groups of compounds:								
15-2.	→ Dioxins and Furans - High Resolution GC/MS								
15-3.	→ Phenols - EPA method 8040								
15-4.	→ PAHs - EPA method 8100								
15-5.	→ Phthalates - EPA method 8060								
	→ VOCs - EPA method 8010/8020 + Styrene and Xylenes								

what method analysis

**PCDD/PCDF SUMMARY REPORT**

SAMPLE: L9331402-001 (U11803) *Spent Carbon*

PROJECT ID: NA

SPECIFIC ANALYTES	CONC (PPB)	DL (PPB)	BLANK (PPB)	Definitions:
2,3,7,8-TCDD	ND	3.5	ND	<b>CONC</b> – The concentration, given in parts per billion (ppb) or parts per trillion (ppt).  <b>DL</b> – The detection limit, given in parts per billion (ppb), parts per trillion (ppt), or in nanograms (ng).  <b>BLANK</b> – The concentration of the method blank.  <b>ND</b> – (Non-Detect) The concentration of the analyte is less than the detection limit.
1,2,3,7,8-PeCDD	ND	5.1	ND	
1,2,3,4,7,8-HxCDD	ND	5.9	ND	
1,2,3,6,7,8-HxCDD	30.9	5.7	ND	
1,2,3,7,8,9-HxCDD	ND	5.5	ND	
1,2,3,4,6,7,8-HpCDD	2930	7.4	ND	
OCDD	27500 –	11	ND	
2,3,7,8-TCDF	ND	2.4	ND	
1,2,3,7,8-PeCDF	ND	3.2	ND	
2,3,4,7,8-PeCDF	ND	3.1	ND	
1,2,3,4,7,8-HxCDF	ND	4.2	ND	
1,2,3,6,7,8-HxCDF	17.2	3.6	ND	
2,3,4,6,7,8-HxCDF	ND	4	ND	
1,2,3,7,8,9-HxCDF	ND	4.7	ND	
1,2,3,4,6,7,8-HpCDF	ND	5.3	ND	
1,2,3,4,7,8,9-HpCDF	21.4	6.8	ND	
OCDF	2980	7.9	ND	

TOTAL ANALYTES	CONC (PPB)	DL (PPB)	BLANK (PPB)
TOTAL TCDD	ND	3.50	ND
TOTAL PeCDD	ND	5.10	ND
TOTAL HxCDD	46.4 –	5.90	ND
TOTAL HpCDD	3850 –	7.40	ND
TOTAL TCDF	2.5	2.40	ND
TOTAL PeCDF	23.3	3.20	ND
TOTAL HxCDF	291	4.70	ND
TOTAL HpCDF	1950	6.80	ND

**TOTAL DIOXINS/FURANS: 36643.2 PPB**

**TOTAL 2,3,7,8-TCDD TOXICITY (1989 ITEF) EQUIVALENTS: 65 PPB**

For information, please reference the following when contacting our Technical Services Department:

TLH Project: P012744  
 TLH Batch: B000645T  
 TLH File: MA01291

December 19, 2012

Mr. Paul Walz  
Bay West, Inc.  
5 Empire Drive  
St. Paul, MN 55103

RE: Project: J007095/10 Strebtor Inc  
Pace Project No.: 10214461

Dear Mr. Walz:

Enclosed are the analytical results for sample(s) received by the laboratory on December 06, 2012. The results relate only to the samples included in this report. Results reported herein conform to the most current TNI standards and the laboratory's Quality Assurance Manual, where applicable, unless otherwise noted in the body of the report.

Some analyses have been subcontracted outside of the Pace Network. The subcontracted laboratory report has been attached.

If you have any questions concerning this report, please feel free to contact me.

Sincerely,



Carolynne Trout

carolynne.trout@pacelabs.com  
Project Manager

Enclosures



## REPORT OF LABORATORY ANALYSIS

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Page 1 of 3



### SAMPLE SUMMARY

Project: J007095/10 Strebtor Inc  
Pace Project No.: 10214461

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Lab ID	Sample ID	Matrix	Date Collected	Date Received
10214461001	Spent Carbon	Solid	12/05/12 11:05	12/06/12 09:20

### REPORT OF LABORATORY ANALYSIS

Page 2 of 3

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## PROJECT NARRATIVE

Project:

Pace Project No.:

---

**Method:**

**Description:**

**Client:**

**Date:**

This data package has been reviewed for quality and completeness and is approved for release.

## REPORT OF LABORATORY ANALYSIS

Page 3 of 3

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December 18, 2012

Pace Analytical  
Attn: Carolynne Trout  
1700 Elm Street Suite 200  
Minneapolis, MN 55414

**Project: J007095/10 Strebor Inc.**

Dear Carolynne Trout,

Enclosed is a copy of the laboratory report for the following work order(s) received by TriMatrix Laboratories:

<b>Work Order</b>	<b>Received</b>	<b>Description</b>
1212156	12/07/2012	10214461

This report relates only to the sample(s) as received. Test results are in compliance with the requirements of the National Environmental Laboratory Accreditation Program (NELAP) and/or one of the following certification programs:

ACLASS DoD-ELAP/ISO17025 (#ADE-1542); Arkansas DEP (#12-056-0); Florida DEP (#E87622-24); Georgia EPD (#E87622-24); Illinois DEP (#002841); Kansas DPH (#E-10302); Kentucky DEP (#0021); Louisiana DEP (#03068); Michigan DPH (#0034); Minnesota DPH (#367345); New York ELAP (#46503); North Carolina DNRE (#659); Texas CEQ (#T104704495-12-2); Virginia DCLS (#1622); Wisconsin DNR (#999472650); USDA Soil Import Permit (#P330-09-00163).

Any qualification or narration of results, including sample acceptance requirements and test exceptions to the above referenced programs, is presented in the Statement of Data Qualifications section of this report. Estimates of analytical uncertainties and certification documents for the test results contained within this report are available upon request.

If you have any questions or require further information, please do not hesitate to contact me.

Sincerely,



James D. McFadden  
Project Chemist

cc: Carolynne Trout



### ANALYTICAL REPORT

Client: **Pace Analytical**  
 Project: J007095/10 Strebor Inc.  
 Client Sample ID: **Spent Carbon**  
 Lab Sample ID: **1212156-01**  
 Matrix: Soil  
 Unit: mg/kg dry  
 Dilution Factor: 2  
 QC Batch: 1215313  
 Percent Solids: 55

Work Order: **1212156**  
 Description: 10214461  
 Sampled: 12/05/12 11:05  
 Sampled By: Client  
 Received: 12/07/12 08:15  
 Prepared: 12/13/2012 By: JTS  
 Analyzed: 12/14/12 By: DWJ  
 Analytical Batch: 2L17075

#### \*Semivolatile Organic Compounds by EPA Method 8270C

CAS Number	Analyte	Analytical Result	RL
83-32-9	Acenaphthene	<0.060	0.060
208-96-8	Acenaphthylene	<0.060	0.060
98-86-2	Acetophenone	<0.060	0.060
62-53-3	Aniline	<0.24	0.24
120-12-7	Anthracene	<0.060	0.060
1912-24-9	Atrazine	<0.060	0.060
100-52-7	Benzaldehyde	<b>0.48</b>	0.12
92-87-5	Benzidine	<2.4	2.4
56-55-3	Benzo(a)anthracene	<0.060	0.060
50-32-8	Benzo(a)pyrene	<0.060	0.060
205-99-2	Benzo(b)fluoranthene	<0.060	0.060
207-08-9	Benzo(k)fluoranthene	<0.060	0.060
191-24-2	Benzo(g,h,i)perylene	<0.12	0.12
65-85-0	Benzoic Acid	<1.2	1.2
100-51-6	Benzyl Alcohol	<0.060	0.060
92-52-4	1,1'-Biphenyl	<0.060	0.060
101-55-3	4-Bromophenyl Phenyl Ether	<0.060	0.060
85-68-7	Butyl Benzyl Phthalate	<0.12	0.12
105-60-2	Caprolactam	<0.12	0.12
86-74-8	Carbazole	<0.60	0.60
59-50-7	4-Chloro-3-methylphenol	<0.060	0.060
95-51-2	2-Chloroaniline	<0.060	0.060
106-47-8	4-Chloroaniline	<0.24	0.24
111-91-1	Bis(2-chloroethoxy)methane	<0.060	0.060
111-44-4	Bis(2-chloroethyl) Ether	<0.060	0.060
108-60-1	Bis(2-chloroisopropyl) Ether	<0.060	0.060
91-58-7	2-Chloronaphthalene	<0.060	0.060
95-57-8	2-Chlorophenol	<0.060	0.060
7005-72-3	4-Chlorophenyl Phenyl Ether	<0.060	0.060
218-01-9	Chrysene	<0.060	0.060
53-70-3	Dibenz(a,h)anthracene	<0.12	0.12

Continued on next page

\*See Statement of Data Qualifications

### ANALYTICAL REPORT

Client: **Pace Analytical**  
 Project: J007095/10 Strebtor Inc.  
 Client Sample ID: **Spent Carbon**  
 Lab Sample ID: **1212156-01**  
 Matrix: Soil  
 Unit: mg/kg dry  
 Dilution Factor: 2  
 QC Batch: 1215313  
 Percent Solids: 55

Work Order: **1212156**  
 Description: 10214461  
 Sampled: 12/05/12 11:05  
 Sampled By: Client  
 Received: 12/07/12 08:15  
 Prepared: 12/13/2012 By: JTS  
 Analyzed: 12/14/12 By: DWJ  
 Analytical Batch: 2L17075

#### \*Semivolatile Organic Compounds by EPA Method 8270C (Continued)

CAS Number	Analyte	Analytical Result	RL
132-64-9	Dibenzofuran	<0.060	0.060
84-74-2	Di-n-butyl Phthalate	<0.24	0.24
106-46-7	1,4-Dichlorobenzene	<b>0.22</b>	0.060
95-50-1	1,2-Dichlorobenzene	<0.060	0.060
541-73-1	1,3-Dichlorobenzene	<0.060	0.060
91-94-1	3,3'-Dichlorobenzidine	<3.0	3.0
120-83-2	2,4-Dichlorophenol	<0.12	0.12
87-65-0	2,6-Dichlorophenol	<0.12	0.12
84-66-2	Diethyl Phthalate	<0.060	0.060
105-67-9	2,4-Dimethylphenol	<0.60	0.60
131-11-3	Dimethyl Phthalate	<0.060	0.060
534-52-1	4,6-Dinitro-2-methylphenol	<0.60	0.60
51-28-5	2,4-Dinitrophenol	<0.60	0.60
606-20-2	2,6-Dinitrotoluene	<0.060	0.060
121-14-2	2,4-Dinitrotoluene	<0.12	0.12
117-84-0	Di-n-octyl Phthalate	<0.060	0.060
122-66-7	1,2-Diphenylhydrazine	<0.060	0.060
117-81-7	Bis(2-ethylhexyl) Phthalate	<b>0.68</b>	0.12
206-44-0	Fluoranthene	<0.060	0.060
86-73-7	Fluorene	<0.12	0.12
118-74-1	Hexachlorobenzene	<0.060	0.060
87-68-3	Hexachlorobutadiene	<0.060	0.060
77-47-4	Hexachlorocyclopentadiene	<0.060	0.060
67-72-1	Hexachloroethane	<0.060	0.060
193-39-5	Indeno(1,2,3-cd)pyrene	<0.12	0.12
78-59-1	Isophorone	<0.060	0.060
91-57-6	2-Methylnaphthalene	<0.060	0.060
90-12-0	1-Methylnaphthalene	<0.060	0.060
106-44-5	4-Methylphenol	<0.060	0.060
108-39-4	3-Methylphenol	<0.12	0.12
95-48-7	2-Methylphenol	<0.060	0.060

Continued on next page

\*See Statement of Data Qualifications

**ANALYTICAL REPORT**

Client: **Pace Analytical**  
 Project: J007095/10 Strebtor Inc.  
 Client Sample ID: **Spent Carbon**  
 Lab Sample ID: **1212156-01**  
 Matrix: Soil  
 Unit: mg/kg dry  
 Dilution Factor: 2  
 QC Batch: 1215313  
 Percent Solids: 55

Work Order: **1212156**  
 Description: 10214461  
 Sampled: 12/05/12 11:05  
 Sampled By: Client  
 Received: 12/07/12 08:15  
 Prepared: 12/13/2012 By: JTS  
 Analyzed: 12/14/12 By: DWJ  
 Analytical Batch: 2L17075

**\*Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
91-20-3	Naphthalene	<b>0.094</b>	0.060
100-01-6	4-Nitroaniline	<0.12	0.12
88-74-4	2-Nitroaniline	<0.060	0.060
99-09-2	3-Nitroaniline	<0.12	0.12
98-95-3	Nitrobenzene	<0.060	0.060
88-75-5	2-Nitrophenol	<0.060	0.060
100-02-7	4-Nitrophenol	<2.4	2.4
62-75-9	N-Nitroso-dimethylamine	<0.12	0.12
86-30-6	N-Nitroso-diphenylamine	<0.060	0.060
621-64-7	N-Nitroso-di-n-propylamine	<0.060	0.060
87-86-5	Pentachlorophenol	<b>2.0</b>	0.60
85-01-8	Phenanthrene	<0.060	0.060
108-95-2	Phenol	<0.60	0.60
129-00-0	Pyrene	<0.060	0.060
110-86-1	Pyridine	<0.12	0.12
95-94-3	1,2,4,5-Tetrachlorobenzene	<0.12	0.12
58-90-2	2,3,4,6-Tetrachlorophenol	<b>0.22</b>	0.12
120-82-1	1,2,4-Trichlorobenzene	<0.060	0.060
95-95-4	2,4,5-Trichlorophenol	<0.060	0.060
88-06-2	2,4,6-Trichlorophenol	<0.060	0.060

<i>Surrogates:</i>	<i>% Recovery</i>	<i>Control Limits</i>
<i>2-Fluorophenol</i>	<i>49</i>	<i>33-113</i>
<i>Phenol-d6</i>	<i>49</i>	<i>30-115</i>
<i>Nitrobenzene-d5</i>	<i>58</i>	<i>33-131</i>

Continued on next page

\*See Statement of Data Qualifications

**ANALYTICAL REPORT**

Client: **Pace Analytical**  
 Project: J007095/10 Strebor Inc.  
 Client Sample ID: **Spent Carbon**  
 Lab Sample ID: **1212156-01**  
 Matrix: Soil  
 Unit: mg/kg dry  
 Dilution Factor: 2  
 QC Batch: 1215313  
 Percent Solids: 55

Work Order: **1212156**  
 Description: 10214461  
 Sampled: 12/05/12 11:05  
 Sampled By: Client  
 Received: 12/07/12 08:15  
 Prepared: 12/13/2012 By: JTS  
 Analyzed: 12/14/12 By: DWJ  
 Analytical Batch: 2L17075

**\*Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

<i>Surrogates (Continued):</i>	<i>% Recovery</i>	<i>Control Limits</i>
<i>2-Fluorobiphenyl</i>	<i>70</i>	<i>46-122</i>
<i>2,4,6-Tribromophenol</i>	<i>31</i>	<i>12-124</i>
<i>o-Terphenyl</i>	<i>36</i>	<i>20-155</i>

\*See Statement of Data Qualifications

**ANALYTICAL REPORT**

Client: **Pace Analytical**  
 Project: J007095/10 Strebtor Inc.  
 Client Sample ID: **Spent Carbon**  
 Lab Sample ID: **1212156-01**  
 Matrix: Soil

Work Order: **1212156**  
 Description: 10214461  
 Sampled: 12/05/12 11:05  
 Sampled By: Client  
 Received: 12/07/12 08:15

**Physical/Chemical Parameters by EPA/APHA/ASTM Methods**

Analyte	Analytical Result	RL	Unit	Dilution Factor	Method	Date Time Analyzed	By	QC Batch
<b>Percent Solids</b>	<b>55</b>	0.1	%	1	USEPA-3550C	12/11/12 15:45	SKA	1215229



### QUALITY CONTROL REPORT

#### Semivolatile Organic Compounds by EPA Method 8270C

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1215313** 3550C Sonication Extraction/USEPA-8270C

**Method Blank**

Unit: mg/kg wet

Analyzed: 12/13/2012 By: DWJ  
Analytical Batch: 2L17069

Acenaphthene	<0.017		0.017
Acenaphthylene	<0.017		0.017
Acetophenone	<0.017	--	0.017
Aniline	<0.067		0.067
Anthracene	<0.017		0.017
Atrazine	<0.017		0.017
Benzaldehyde	<0.033		0.033
Benzidine	<0.67		0.67
Benzo(a)anthracene	<0.017	--	0.017
Benzo(a)pyrene	<0.017		0.017
Benzo(b)fluoranthene	<0.017		0.017
Benzo(k)fluoranthene	<0.017		0.017
Benzo(g,h,i)perylene	<0.033		0.033
Benzoic Acid	<0.33	--	0.33
Benzyl Alcohol	<0.017	--	0.017
1,1'-Biphenyl	<0.017	--	0.017
4-Bromophenyl Phenyl Ether	<0.017		0.017
Butyl Benzyl Phthalate	<0.033		0.033
Caprolactam	<0.033	--	0.033
Carbazole	<0.17	--	0.17
4-Chloro-3-methylphenol	<0.017	--	0.017
2-Chloroaniline	<0.017		0.017
4-Chloroaniline	<0.067	--	0.067
Bis(2-chloroethoxy)methane	<0.017		0.017
Bis(2-chloroethyl) Ether	<0.017		0.017
Bis(2-chloroisopropyl) Ether	<0.017		0.017
2-Chloronaphthalene	<0.017		0.017
2-Chlorophenol	<0.017	--	0.017
4-Chlorophenyl Phenyl Ether	<0.017		0.017
Dibenz(a,h)anthracene	<0.033		0.033
Dibenzofuran	<0.017		0.017
Di-n-butyl Phthalate	<0.067	--	0.067
1,4-Dichlorobenzene	<0.017	--	0.017
1,2-Dichlorobenzene	<0.017	--	0.017
1,3-Dichlorobenzene	<0.017	--	0.017
3,3'-Dichlorobenzidine	<0.83		0.83

Continued on next page

### QUALITY CONTROL REPORT

#### Semivolatile Organic Compounds by EPA Method 8270C (Continued)

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1215313 (Continued)** 3550C Sonication Extraction/USEPA-8270C

**Method Blank (Continued)**

Analyzed: 12/13/2012 By: DWJ  
 Analytical Batch: 2L17069

Unit: mg/kg wet

2,4-Dichlorophenol			<0.033					0.033
2,6-Dichlorophenol			<0.033					0.033
Diethyl Phthalate			<0.017			--		0.017
2,4-Dimethylphenol			<0.17					0.17
Dimethyl Phthalate			<0.017					0.017
4,6-Dinitro-2-methylphenol			<0.17			--		0.17
2,4-Dinitrophenol			<0.17					0.17
2,6-Dinitrotoluene			<0.017			--		0.017
2,4-Dinitrotoluene			<0.033			--		0.033
Di-n-octyl Phthalate			<0.017					0.017
1,2-Diphenylhydrazine			<0.017			--		0.017
Bis(2-ethylhexyl) Phthalate			<0.033			--		0.033
Fluoranthene			<0.017					0.017
Fluorene			<0.033					0.033
Hexachlorobenzene			<0.017					0.017
Hexachlorobutadiene			<0.017					0.017
Hexachlorocyclopentadiene			<0.017					0.017
Hexachloroethane			<0.017					0.017
Indeno(1,2,3-cd)pyrene			<0.033					0.033
Isophorone			<0.017			--		0.017
2-Methylnaphthalene			<0.017					0.017
1-Methylnaphthalene			<0.017					0.017
4-Methylphenol			<0.017					0.017
3-Methylphenol			<0.033					0.033
2-Methylphenol			<0.017					0.017
Naphthalene			<0.017			--		0.017
4-Nitroaniline			<0.033			--		0.033
2-Nitroaniline			<0.017					0.017
3-Nitroaniline			<0.033			--		0.033
Nitrobenzene			<0.017					0.017
2-Nitrophenol			<0.017					0.017
4-Nitrophenol			<0.67			--		0.67
N-Nitroso-dimethylamine			<0.033					0.033
N-Nitroso-diphenylamine			<0.017					0.017
N-Nitroso-di-n-propylamine			<0.017			--		0.017
Pentachlorophenol			<0.17			--		0.17

Continued on next page

**QUALITY CONTROL REPORT**
**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1215313 (Continued)** 3550C Sonication Extraction/USEPA-8270C

**Method Blank (Continued)**

Unit: mg/kg wet

 Analyzed: 12/13/2012 By: DWJ  
 Analytical Batch: 2L17069

Phenanthrene			<0.017			--		0.017
Phenol			<0.17			--		0.17
Pyrene			<0.017					0.017
Pyridine			<0.033			--		0.033
1,2,4,5-Tetrachlorobenzene			<0.033					0.033
2,3,4,6-Tetrachlorophenol			<0.033					0.033
1,2,4-Trichlorobenzene			<0.017			--		0.017
2,4,5-Trichlorophenol			<0.017					0.017
2,4,6-Trichlorophenol			<0.017					0.017

**Surrogates:**

<i>2-Fluorophenol</i>				70	33-113			
<i>Phenol-d6</i>				73	30-115			
<i>Nitrobenzene-d5</i>				68	33-131			
<i>2-Fluorobiphenyl</i>				79	46-122			
<i>2,4,6-Tribromophenol</i>				62	12-124			
<i>o-Terphenyl</i>				79	20-155			

**Laboratory Control Sample**

Unit: mg/kg wet

 Analyzed: 12/14/2012 By: DWJ  
 Analytical Batch: 2L17069

Acenaphthene	0.348	<b>0.247</b>		71	55-113	--		0.017
4-Chloro-3-methylphenol	0.348	<b>0.231</b>		66	57-124	--		0.017
2-Chlorophenol	0.348	<b>0.229</b>		66	62-118	--		0.017
1,4-Dichlorobenzene	0.348	<b>0.237</b>		68	61-111	--		0.017
2,4-Dinitrotoluene	0.348	<b>0.253</b>		73	51-128	--		0.033
Naphthalene	0.348	<b>0.240</b>		69	52-128	--		0.017
4-Nitrophenol	0.348	<b>0.267</b>		77	36-131	--		0.67
N-Nitroso-di-n-propylamine	0.348	<b>0.239</b>		69	48-127	--		0.017
Pentachlorophenol	0.348	<b>0.204</b>		58	19-117	--		0.17
Phenol	0.348	<b>0.214</b>		61	53-120	--		0.17
Pyrene	0.348	<b>0.276</b>		79	60-132	--		0.017

Continued on next page

**QUALITY CONTROL REPORT**
**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1215313 (Continued)** 3550C Sonication Extraction/USEPA-8270C

**Laboratory Control Sample (Continued)**

Unit: mg/kg wet

 Analyzed: 12/14/2012 By: DWJ  
 Analytical Batch: 2L17069

 1,2,4-Trichlorobenzene 0.348 **0.229** 66 51-110 -- 0.017

**Surrogates:**

<i>2-Fluorophenol</i>	66	33-113
<i>Phenol-d6</i>	68	30-115
<i>Nitrobenzene-d5</i>	62	33-131
<i>2-Fluorobiphenyl</i>	72	46-122
<i>2,4,6-Tribromophenol</i>	70	12-124
<i>o-Terphenyl</i>	75	20-155

**Laboratory Control Sample Duplicate**

Unit: mg/kg wet

 Analyzed: 12/14/2012 By: DWJ  
 Analytical Batch: 2L17069

Acenaphthene	0.348	<b>0.245</b>	70	55-113	1	20	0.017
4-Chloro-3-methylphenol	0.348	<b>0.236</b>	68	57-124	2	20	0.017
2-Chlorophenol	0.348	<b>0.231</b>	66	62-118	0.9	20	0.017
1,4-Dichlorobenzene	0.348	<b>0.235</b>	67	61-111	0.7	20	0.017
2,4-Dinitrotoluene	0.348	<b>0.263</b>	76	51-128	4	20	0.033
Naphthalene	0.348	<b>0.239</b>	69	52-128	0.1	20	0.017
4-Nitrophenol	0.348	<b>0.268</b>	77	36-131	0.4	20	0.67
N-Nitroso-di-n-propylamine	0.348	<b>0.238</b>	68	48-127	0.7	20	0.017
Pentachlorophenol	0.348	<b>0.176</b>	50	19-117	15	20	0.17
Phenol	0.348	<b>0.216</b>	62	53-120	0.8	20	0.17
Pyrene	0.348	<b>0.282</b>	81	60-132	2	20	0.017
1,2,4-Trichlorobenzene	0.348	<b>0.227</b>	65	51-110	0.6	20	0.017

**Surrogates:**

<i>2-Fluorophenol</i>	65	33-113
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Continued on next page

\*See Statement of Data Qualifications

Page 10 of 16

**QUALITY CONTROL REPORT**
**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
---------	--------------	------------	--------	--------------	----------------	-----	------------	----

**QC Batch: 1215313 (Continued)** 3550C Sonication Extraction/USEPA-8270C

**Laboratory Control Sample Duplicate (Continued)**

Analyzed: 12/14/2012 By: DWJ

Unit: mg/kg wet

Analytical Batch: 2L17069

**Surrogates (Continued):**

<i>Phenol-d6</i>	67	30-115
<i>Nitrobenzene-d5</i>	61	33-131
<i>2-Fluorobiphenyl</i>	71	46-122
<i>2,4,6-Tribromophenol</i>	66	12-124
<i>o-Terphenyl</i>	72	20-155

**QUALITY CONTROL REPORT**
**Physical/Chemical Parameters by EPA/APHA/ASTM Methods**

QC Type	Sample Conc.	Spike Qty.	Result	Unit	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
<b>Analyte: Percent Solids/USEPA-3550C</b>									
QC Batch: 1215229 (General Inorganic Prep)						Analyzed: 12/11/2012 By: SKA			
Method Blank			<0.1	%				0.1	



**PRETREATMENT SUMMARY PAGE**

Client: **Pace Analytical**  
Project: **J007095/10 Strebor Inc.**

<b>Pretreatment</b>	<b>Lab Sample ID</b>	<b>QC Batch</b>	<b>By</b>	<b>Date &amp; Time Prepared</b>
USEPA-3550C Ultrasonic Extraction	1212156-01	1215313	JTS	12/13/2012 7:56

**STATEMENT OF DATA QUALIFICATIONS****Semivolatile Organic Compounds by EPA Method 8270C**

**Qualification:** The quality control batch(s), associated with the following samples and analyses, do not contain an MS/MSD or MS/DUP due to insufficient sample volumes. An LCS and LCSD were analyzed as the measure of batch precision and accuracy.

Analysis: USEPA-8270C

Sample: 1212156-01      Spent Carbon

**Qualification:** Manual integration was required on the analytes listed below. All manual integrations were performed and reviewed in accordance with TriMatrix laboratory policy.

Analysis: USEPA-8270C

Sample/Analyte: 1215313-BS1

Chrysene-d12



Chain of Custody

#11095



Workorder: 10214461

Workorder Name: J007095/10 Strebor Inc

Results Requested 12/19/2012

Report / Invoice To		Subcontract To		Requested Analysis																
Carolynne Trout Pace Analytical Minnesota 1700 Elm Street Suite 200 Minneapolis, MN 55414 Phone (612)607-1700 Email: carlyne.trout@pacelabs.com		PO		MS (10/21/12) ID 24, 5 Trichlorophenol 2, 4, 6-Trichlorophenol 2, 4, 6-Trichlorophenol 2, 4, 6-Trichlorophenol 2, 4, 6-Trichlorophenol																
Item	Sample ID	Collect Date/Time	Lab ID	Matrix	Preserved Containers															
1	Spent Carbon	12/5/2012 11:05	10214461001	Solid																
2																				
3																				
4																				
5																				
												Comments E-1212156 LAB USE ONLY -01								
Transfers	Released By	Date/Time	Received By	Date/Time	Comments															
1		12/19/12 09:15	[Signature]	12/19/12 09:15	SOL T work, 100 minutes on the sample															
2																				
3																				
Cooler Temperature on Receipt		54 °C	Custody Seal		Y or N	Received on Ice		Y or N	Samples Intact											Y or N

### SAMPLE RECEIVING / LOG-IN CHECKLIST

		Client: <u>Pace Analytical</u> Report Record Page/Line #: <u>47-5</u>	Work Order #: <u>12/2156</u> Project Client: <u>JONE</u> Samples:
Recorded by (Initials/Date): <u>LR 12/7/12</u>		<input type="checkbox"/> Cooler <input type="checkbox"/> Box <input type="checkbox"/> Other: <u>1</u>	<input type="checkbox"/> IR Gun (#202) <input type="checkbox"/> Digital Thermometer (#54) <input type="checkbox"/> Other (#):
See Additional Cooler Information Form:			

Cooler #	Time	Cooler #	Time	Cooler #	Time
<u>Pace</u>	<u>0855</u>				

Custody Seals: <input type="checkbox"/> None <input type="checkbox"/> Present / Intact <input type="checkbox"/> Present / Not Intact	Custody Seals: <input type="checkbox"/> None <input type="checkbox"/> Present / Intact <input type="checkbox"/> Present / Not Intact	Custody Seals: <input type="checkbox"/> None <input type="checkbox"/> Present / Intact <input type="checkbox"/> Present / Not Intact	Custody Seals: <input type="checkbox"/> None <input type="checkbox"/> Present / Intact <input type="checkbox"/> Present / Not Intact
Coolant Location: <input type="checkbox"/> Dispersed / Top / Middle / Bottom	Coolant Location: <input type="checkbox"/> Dispersed / Top / Middle / Bottom	Coolant Location: <input type="checkbox"/> Dispersed / Top / Middle / Bottom	Coolant Location: <input type="checkbox"/> Dispersed / Top / Middle / Bottom
Coolant/Temperature Taken Via: <input type="checkbox"/> Loose Ice / Avg 2-3 containers <input checked="" type="checkbox"/> Bagged Ice / Avg 2-3 containers <input type="checkbox"/> Blue Ice / Avg 2-3 containers <input type="checkbox"/> None / Avg 2-3 containers	Coolant/Temperature Taken Via: <input type="checkbox"/> Loose Ice / Avg 2-3 containers <input type="checkbox"/> Bagged Ice / Avg 2-3 containers <input type="checkbox"/> Blue Ice / Avg 2-3 containers <input type="checkbox"/> None / Avg 2-3 containers	Coolant/Temperature Taken Via: <input type="checkbox"/> Loose Ice / Avg 2-3 containers <input type="checkbox"/> Bagged Ice / Avg 2-3 containers <input type="checkbox"/> Blue Ice / Avg 2-3 containers <input type="checkbox"/> None / Avg 2-3 containers	Coolant/Temperature Taken Via: <input type="checkbox"/> Loose Ice / Avg 2-3 containers <input type="checkbox"/> Bagged Ice / Avg 2-3 containers <input type="checkbox"/> Blue Ice / Avg 2-3 containers <input type="checkbox"/> None / Avg 2-3 containers
Alternate Temperature Taken Via: <input checked="" type="checkbox"/> Temperature Blank (TB) <input type="checkbox"/> 1 Container	Alternate Temperature Taken Via: <input type="checkbox"/> Temperature Blank (TB) <input type="checkbox"/> 1 Container	Alternate Temperature Taken Via: <input type="checkbox"/> Temperature Blank (TB) <input type="checkbox"/> 1 Container	Alternate Temperature Taken Via: <input type="checkbox"/> Temperature Blank (TB) <input type="checkbox"/> 1 Container
Recorded °C    Correction Factor °C    Actual °C Temp Blank:    -    - <u>4.5</u>	Recorded °C    Correction Factor °C    Actual °C Temp Blank:    -    -	Recorded °C    Correction Factor °C    Actual °C Temp Blank:    -    -	Recorded °C    Correction Factor °C    Actual °C Temp Blank:    -    -
TB location: <u>Representative</u> / Not Representative 1: <u>5.4</u> - <u>5.4</u> 2: 3:	TB location:    Representative / Not Representative 1: 2: 3:	TB location:    Representative / Not Representative 1: 2: 3:	TB location:    Representative / Not Representative 1: 2: 3:
Average °C <input type="checkbox"/> Cooler ID on COC? <u>5.4</u> <input type="checkbox"/> VOC Trip Blank received?	Average °C <input type="checkbox"/> Cooler ID on COC? <input type="checkbox"/> VOC Trip Blank received?	Average °C <input type="checkbox"/> Cooler ID on COC? <input type="checkbox"/> VOC Trip Blank received?	Average °C <input type="checkbox"/> Cooler ID on COC? <input type="checkbox"/> VOC Trip Blank received?

**If any shaded areas checked, complete Sample Receiving Non-Conformance and/or Inventory Form**

<b>Paperwork Received</b> Yes    No <input type="checkbox"/> Chain of Custody record(s)? If No, Initiated By _____ <input type="checkbox"/> Received for Lab Signed/Date/Time? <input type="checkbox"/> Shipping document? <input type="checkbox"/> Other: _____ <b>COC Information</b> <input type="checkbox"/> TriMatrix COC <input checked="" type="checkbox"/> Other: _____ COC ID Numbers:	<b>Check Sample Preservation</b> N/A    Yes    No <input checked="" type="checkbox"/> Average sample temperature (±0.1 °C) <input checked="" type="checkbox"/> Was thermal preservation required? If "No", Project Chemical Approval Initials: _____ If "Yes", Completed Non-Con Cooler - Cont. Inventory Form? <input type="checkbox"/> Completed Sample Preservation Verification Form? <input checked="" type="checkbox"/> Samples chemically preserved correctly? If "No", added orange tag? <input checked="" type="checkbox"/> Received pre-preserved VOC soils? <input type="checkbox"/> MeOH <input type="checkbox"/> Na <sub>2</sub> SO <sub>4</sub>
<b>Check COC for Accuracy</b> Yes    No <input type="checkbox"/> Analysis Requested? <input type="checkbox"/> Sample ID matches COC? <input checked="" type="checkbox"/> Sample Date and Time matches COC? <input type="checkbox"/> Container type completed on COC? <input checked="" type="checkbox"/> All container types indicated are received?	<b>Check for Short Hold-Time Prep/Analyses</b> <input type="checkbox"/> Bacteriology <input type="checkbox"/> Air Bags <input type="checkbox"/> EnCore / Methanol Pre-Preserved <input type="checkbox"/> Formaldehyde/Aldehyde <input type="checkbox"/> Green-tagged containers <input type="checkbox"/> Yellow/White-tagged Labbers (SV Prep-Lab) <div style="border: 1px solid black; padding: 5px; margin-top: 10px;"> <b>AFTER HOURS ONLY:</b>          COPIES OF COC TO LAB AREA(S)  <input checked="" type="checkbox"/> NONE RECEIVED  <input type="checkbox"/> RECEIVED COCs TO LAB(S)       </div>
<b>Sample Condition Summary</b> N/A    Yes    No <input type="checkbox"/> Broken containers/bds? <input type="checkbox"/> Missing or incomplete labels? <input type="checkbox"/> Illegible information on labels? <input type="checkbox"/> Low volume received? <input type="checkbox"/> Inappropriate or non-TriMatrix containers received? <input type="checkbox"/> VOC Vials / TOX containers have headspace? <input type="checkbox"/> Extra sample locations / containers not listed on COC?	<b>Notes</b> <input type="checkbox"/> Trip Blank received <input type="checkbox"/> Trip Blank not listed on COC Cooler Received (Date/Time)    Paperwork Delivered (Date/Time)    <1 Hour Goal Met? <u>12/7/12 0815</u> <u>12/7/12 0901</u> (Yes / No)

**CHAIN-OF-CUSTODY / Analytical Request Document**

The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

1021446

Page: 1 of 1  
1582752

<b>Section A</b> Required Client Information:		<b>Section B</b> Required Project Information:		<b>Section C</b> Invoice Information:		<b>REGULATORY AGENCY</b>	
Company: Bay West Inc.		Report To: Paul Walz		Attention: Accounts Payable		NPDES <input type="checkbox"/> GROUND WATER <input type="checkbox"/> DRINKING WATER <input type="checkbox"/>	
Address: 5 Empire Drive ST. PAUL, MN 55103		Copy To:		Company Name: Bay West Inc.		UST <input type="checkbox"/> RCRA <input type="checkbox"/> OTHER <input type="checkbox"/>	
Email To: paulw@baywest.com		Purchase Order No.: 36287		Address: 5 Empire Drive		Site Location	
Phone: 651-291-3491 Fax:		Project Name: STrebor Inc.		Pace Quote Reference:		STATE: MI	
Requested Due Date/TAT:		Project Number: J007095/10		Pace Project Manager:		Pace Profile #:	

ITEM #	Section D Required Client Information	Matrix Codes MATRIX / CODE	MATRIX CODE (see valid codes to left)	SAMPLE TYPE (G=GRAB C=COMP)	COLLECTED				SAMPLE TEMP AT COLLECTION	# OF CONTAINERS	Preservatives								Analysis Test ↓	Residual Chlorine (Y/N)	Pace Project No./ Lab I.D.				
					COMPOSITE START	COMPOSITE END/GRAB	DATE	TIME			DATE	TIME	Unpreserved	H <sub>2</sub> SO <sub>4</sub>	HNO <sub>3</sub>	HCl	NaOH	Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub>				Methanol	Other	Y/N	N
1	Spent Carbon		SL	G			12-5	1105	60	2	X														
2																									
3																									
4																									
5																									
6																									
7																									
8																									
9																									
10																									
11																									
12																									

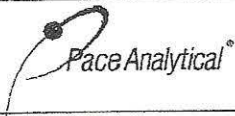
ADDITIONAL COMMENTS	RELINQUISHED BY / AFFILIATION	DATE	TIME	ACCEPTED BY / AFFILIATION	DATE	TIME	SAMPLE CONDITIONS			
Item #1 Analyze for F027 parameters (list attached) CDraw 12/6/12 per Paul Walz	Michael E. McElish / Bay West	12-5-12	1500	TN / Pace	12/6/12	920	0.8	Y	Y	Y

<b>SAMPLER NAME AND SIGNATURE</b>		Temp in °C	Received on Ice (Y/N)	Custody Sealed Cooler (Y/N)	Samples Intact (Y/N)
PRINT Name of SAMPLER: Michael E. McElish	DATE Signed (MM/DD/YY): 12-05-12				
SIGNATURE of SAMPLER: <i>Michael E. McElish</i>					

ORIGINAL

20 of 23

\*Important Note: By signing this form you are accepting Pace's NET 30 day payment terms and agreeing to late charges of 1.5% per month for any invoices not paid within 30 days.



Document Name:  
Sample Condition Upon Receipt Form

Document Revised: 13Nov2012  
Page 1 of 1

Document No.:  
F-MN-L-213-rev.05

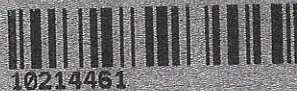
Issuing Authority:  
Pace Minnesota Quality Office

Sample Condition  
Upon Receipt

Client Name:

Project #:

WO#: 10214461



Courier:  Fed Ex  UPS  USPS  Client  
 Commercial  Pace  Other:

Tracking Number: 8009 7975 6127

Custody Seal on Cooler/Box Present?  Yes  No      Seals Intact?  Yes  No

Optional: Proj. Due Date:      Proj. Name:

Packing Material:  Bubble Wrap  Bubble Bags  None  Other:      Temp Blank?  Yes  No

Thermometer Used:  B88A912167504  80512447      Type of Ice:  Wet  Blue  None  Samples on ice, cooling process has begun

Cooler Temp Read (°C): 0.3      Cooler Temp Corrected (°C): 0.8      Biological Tissue Frozen?  Yes  No  
Temp should be above freezing to 6°C      Date and Initials of Person Examining Contents: 12/6/12 TN

Comments:

Chain of Custody Present?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	1.
Chain of Custody Filled Out?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	2.
Chain of Custody Relinquished?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	3.
Sampler Name and/or Signature on COC?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	4.
Samples Arrived within Hold Time?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	5.
Short Hold Time Analysis (<72 hr)?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A	6.
Rush Turn Around Time Requested?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A	7.
Sufficient Volume?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	8.
Correct Containers Used?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	9.
-Pace Containers Used?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	
Containers Intact?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	10.
Filtered Volume Received for Dissolved Tests?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	11.
Sample Labels Match COC?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	12.
-Includes Date/Time/ID/Analysis Matrix: SL		
All containers needing acid/base preservation have been checked? Noncompliances are noted in 13.	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	13.
All containers needing preservation are found to be in compliance with EPA recommendation? (HNO <sub>3</sub> , H <sub>2</sub> SO <sub>4</sub> , HCl<2; NaOH>12)	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	Sample #
Exceptions: VOA, Coliform, TOC, Oil and Grease, WI-DRO (water)-	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No	Initial when completed:      Lot # of added preservative:
Headspace in VOA Vials (>6mm)?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	14.
Trip Blank Present?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	15.
Trip Blank Custody Seals Present?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	
Pace Trip Blank Lot # (if purchased):		

CLIENT NOTIFICATION/RESOLUTION

Field Data Required?  Yes  No

Person Contacted: \_\_\_\_\_ Date/Time: \_\_\_\_\_

Comments/Resolution: \_\_\_\_\_

Project Manager Review: CTM

Date: 12/6/12

Note: Whenever there is a discrepancy affecting North Carolina compliance samples, a copy of this form will be sent to the North Carolina DEHNR Certification Office ( i.e out of hold, incorrect preservative, out of temp, incorrect containers)

TREATMENT STANDARDS FOR HAZARDOUS WASTES

Waste Code	Waste Description and Treatment/Regulatory Subcategory <sup>1</sup>	REGULATED HAZARDOUS CONSTITUENT		WASTEWATERS	NONWASTEWATERS
		Common Name	CAS <sup>2</sup> Number	Concentration mg/l <sup>3</sup> ; or Technology Code <sup>4</sup>	Concentration in mg/kg <sup>5</sup> unless noted as "mg/l TCLP" or Technology Code
		Pentachlorophenol	87-86-5	0.089	7.4
F027	Discarded unused formulations containing tri-, tetra-, or pentachlorophenol or discarded unused formulations containing compounds derived from these chlorophenols. (This listing does not include formulations containing hexachlorophene synthesized from prepurified 2,4,5-trichlorophenol as the sole component.).	HxCDDs (All Hexachlorodibenzop-dioxins)	NA	0.000063	0.001
		HxCDFs (All Hexachlorodibenzofurans)	NA	0.000063	0.001
		PeCDDs (All Pentachlorodibenzop-dioxins)	NA	0.000063	0.001
		PeCDFs (All Pentachlorodibenzofurans)	NA	0.000035	0.001
		TCDDs (All Tetrachlorodibenzop-dioxins)	NA	0.000063	0.001
		TCDFs (All Tetrachlorodibenzofurans)	NA	0.000063	0.001
		2,4,5-Trichlorophenol	95-95-4	0.18	7.4
		2,4,6-Trichlorophenol	88-06-2	0.035	7.4

TREATMENT STANDARDS FOR HAZARDOUS WASTES

Waste Code	Waste Description and Treatment/Regulatory Subcategory <sup>1</sup>	REGULATED HAZARDOUS CONSTITUENT		WASTEWATERS	NONWASTEWATERS
		Common Name	CAS <sup>2</sup> Number	Concentration mg/l <sup>3</sup> ; or Technology Code <sup>4</sup>	Concentration in mg/kg <sup>5</sup> unless noted as "mg/l TCLP" or Technology Code
		2,3,4,6-Tetrachlorophenol	58-90-2	0.030	7.4
		Pentachlorophenol	87-86-5	0.089	7.4
F028	Residues resulting from the incineration or thermal treatment of soil contaminated with EPA Hazardous Wastes Nos. F020, F021, F023, F026, and F027.	HxCDDs (All Hexachlorodibenzo-p-dioxins)	NA	0.000063	0.001
		HxCDFs (All Hexachlorodibenzofurans)	NA	0.000063	0.001
		PeCDDs (All Pentachlorodibenzo-p-dioxins)	NA	0.000063	0.001
		PeCDFs (All Pentachlorodibenzofurans)	NA	0.000035	0.001
		TCDDs (All Tetrachlorodibenzo-p-dioxins)	NA	0.000063	0.001
		TCDFs (All Tetrachlorodibenzofurans)	NA	0.000063	0.001
		2,4,5-Trichlorophenol	95-95-4	0.18	7.4

67 RL in mg/kg

**Report Prepared for:**

Rick VanAllen  
Bay West, Inc.  
5 Empire Drive  
Saint Paul MN 55103

**REPORT OF  
LABORATORY  
ANALYSIS FOR  
PCDD/PCDF**

**Report Prepared Date:**  
December 19, 2012

**Report Information:**

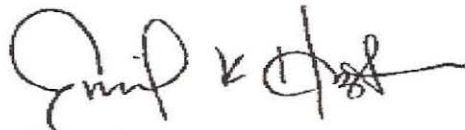
**Pace Project #: 10214459**  
**Sample Receipt Date: 12/06/2012**  
**Client Project #: J007095/10**  
**Client Sub PO #: 36287**  
**State Cert #: 9909**

**Invoicing & Reporting Options:**

The report provided has been invoiced as a Level 2 PCDD/PCDF Report. If an upgrade of this report package is requested, an additional charge may be applied.

Please review the attached invoice for accuracy and forward any questions to Scott Unze, your Pace Project Manager.

**This report has been reviewed by:**



December 19, 2012

Emily Hazelroth, Project Manager  
(612) 607-6407  
(612) 607-6444 (fax)  
emily.hazelroth@pacelabs.com



**Report of Laboratory Analysis**

This report should not be reproduced, except in full, without the written consent of Pace Analytical Services, Inc.

The results relate only to the samples included in this report.



## **DISCUSSION**

This report presents the results from the analysis performed on one sample submitted by a representative of BayWest, Inc. The sample was analyzed for the presence or absence of polychlorodibenzo-p-dioxins (PCDDs) and polychlorodibenzofurans (PCDFs) using a modified version of USEPA Method 8290. The reporting limits were based on signal-to-noise measurements.

The recoveries of the isotopically-labeled PCDD/PCDF internal standards in the sample extract ranged from 66-104%. All of the labeled standard recoveries obtained for this project were within the 40-135% target range specified in Method 8290. Also, since the quantification of the native 2,3,7,8-substituted congeners was based on isotope dilution, the data were automatically corrected for variation in recovery and accurate values were obtained.

In one case, an interfering substance impacted the determination of a PCDD congener; the affected value was flagged "I" due to an incorrect isotope ratio. Concentrations below the calibration range were flagged "J" and should be regarded as estimates. Concentrations above the calibration range were flagged "E" and also should be regarded as estimates.

A laboratory method blank was prepared and analyzed with the sample batch as part of our routine quality control procedures. The results show the blank to contain trace levels of selected congeners. These levels were below the calibration range of the method. The levels reported for the affected congeners in the field sample were higher than the corresponding blank levels by one or more orders of magnitude. These results indicate that the sample processing steps did not contribute significantly to the levels reported for the field sample.

A laboratory spike sample was also prepared with the sample batch using clean sand that had been fortified with native standard materials. The results show that the spiked native compounds were recovered at 80-110%, indicating a high degree of accuracy for these determinations. Matrix spikes were prepared with the sample batch using sample material from a separate project; results from these analyses will be provided upon request.

## **REPORT OF LABORATORY ANALYSIS**

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## Minnesota Laboratory Certifications

Authority	Certificate #	Authority	Certificate #
Alabama	40770	Montana	92
Alaska	MN00064	Nebraska	
Arizona	AZ0014	Nevada	MN_00064_200
Arkansas	88-0680	New Jersey (NE)	MN002
California	01155CA	New Mexico	MN00064
Colorado	MN00064	New York (NEL)	11647
Connecticut	PH-0256	North Carolina	27700
EPA Region 5	WD-15J	North Dakota	R-036
EPA Region 8	8TMS-Q	Ohio	4150
Florida (NELAP)	E87605	Ohio VAP	CL101 9507
Georgia (DNR)	959	Oklahoma	D9922
Guam	959	Oregon (ELAP)	MN200001-005
Hawaii	SLD	Oregon (OREL)	MN300001-001
Idaho	MN00064	Pennsylvania	68-00563
Illinois	200012	Saipan	MP0003
Indiana	C-MN-01	South Carolina	74003001
Indiana	C-MN-01	Tennessee	2818
Iowa	368	Tennessee	02818
Kansas	E-10167	Texas	T104704192-08
Kentucky	90062	Utah (NELAP)	PAM
Louisiana	03086	Virginia	00251
Maine	2007029	Washington	C755
Maryland	322	West Virginia	9952C
Michigan	9909	Wisconsin	999407970
Minnesota	027-053-137	Wyoming	8TMS-Q
Mississippi	MN00064		

## REPORT OF LABORATORY ANALYSIS

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# Appendix A

## Sample Management



# CHAIN-OF-CUSTODY / Analytical Request Document

The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

Page: 1 of 1  
**10214459**  
**1582752**

Report No. .... 10214459\_8290

Section A Required Client Information:		Section B Required Project Information:		Section C Invoice Information:	
Company: <u>Bay West Inc.</u>		Report To: <u>Paul Walz</u>		Attention: <u>Accounts Payable</u>	
Address: <u>5 Empire Drive ST. PAUL, MN 55103</u>		Copy To:		Company Name: <u>Bay West Inc.</u>	
Email To: <u>paulw@baywest.com</u>		Purchase Order No.: <u>36287</u>		Address: <u>5 Empire Drive</u>	
Phone: <u>651-291-3491</u> Fax:		Project Name: <u>STrebor Inc.</u>		Pace Quote Reference:	
Requested Due Date/TAT:		Project Number: <u>5007095/10</u>		Pace Project Manager:	
				Pace Profile #:	

REGULATORY AGENCY		
<input type="checkbox"/> NPDES	<input type="checkbox"/> GROUND WATER	<input type="checkbox"/> DRINKING WATER
<input type="checkbox"/> UST	<input type="checkbox"/> RCRA	<input type="checkbox"/> OTHER _____
Site Location	STATE: <u>MI</u>	

ITEM #	Section D Required Client Information	Matrix Codes MATRIX / CODE	MATRIX CODE (see valid codes to left)	SAMPLE TYPE (G=GRAB C=COMP)	COLLECTED				SAMPLE TEMP AT COLLECTION	# OF CONTAINERS	Requested Analysis Filtered (Y/N)											
					COMPOSITE START		COMPOSITE END/GRAB				Unpreserved	Preservatives									Analysis Test ↓	Residual Chlorine (Y/N)
					DATE	TIME	DATE	TIME				H <sub>2</sub> SO <sub>4</sub>	HNO <sub>3</sub>	HCl	NaOH	Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub>	Methanol	Other				
1	<u>Spent Carbon</u>		<u>SL G</u>		<u>12-5</u>	<u>1105</u>	<u>60</u>	<u>2</u>	<u>X</u>													
2																						
3																						
4																						
5																						
6																						
7																						
8																						
9																						
10																						
11																						
12																						

ADDITIONAL COMMENTS	RELINQUISHED BY / AFFILIATION	DATE	TIME	ACCEPTED BY / AFFILIATION	DATE	TIME	SAMPLE CONDITIONS			
<u>Item #1</u>	<u>Michael E. McClish / Bay West</u>	<u>12-5-12</u>	<u>1500</u>	<u>TN/Pace</u>	<u>12/6/12</u>	<u>920</u>	<u>0.8</u>	<u>Y</u>	<u>Y</u>	<u>Y</u>

SAMPLER NAME AND SIGNATURE  
 PRINT Name of SAMPLER: Michael E. McClish  
 SIGNATURE of SAMPLER: [Signature] DATE Signed (MM/DD/YY): 12-05-12

Temp in °C	Received on Ice (Y/N)	Custody Sealed Cooler (Y/N)	Samples Intact (Y/N)

ORIGINAL

Page 5 of 11

**Sample Condition Upon Receipt**

**Client Name:** Bay West      **Project #:** \_\_\_\_\_

**WO# : 10214459**



**Courier:**  Fed Ex     UPS     USPS     Client  
 Commercial     Pace     Other: \_\_\_\_\_

**Tracking Number:** 8009 7975 6127

**Custody Seal on Cooler/Box Present?**  Yes     No      **Seals Intact?**  Yes     No      **Optional:** Proj. Due Date: \_\_\_\_\_ Proj. Name: \_\_\_\_\_

**Packing Material:**  Bubble Wrap     Bubble Bags     None     Other: \_\_\_\_\_      **Temp Blank?**  Yes     No

**Thermometer Used:**  B88A912167504     80512447      **Type of Ice:**  Wet     Blue     None     Samples on ice, cooling process has begun

**Cooler Temp Read (°C):** 0.3      **Cooler Temp Corrected (°C):** 0.8      **Biological Tissue Frozen?**  Yes     No  
**Temp should be above freezing to 6°C**      **Date and Initials of Person Examining Contents:** 12/6/12 TN

**Comments:**

Chain of Custody Present?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	1.	
Chain of Custody Filled Out?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	2.	
Chain of Custody Relinquished?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	3.	
Sampler Name and/or Signature on COC?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	4.	
Samples Arrived within Hold Time?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	5.	
Short Hold Time Analysis (<72 hr)?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A	6.	
Rush Turn Around Time Requested?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A	7.	
Sufficient Volume?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	8.	
Correct Containers Used?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	9.	
-Pace Containers Used?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A		
Containers Intact?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	10.	
Filtered Volume Received for Dissolved Tests?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	11.	
Sample Labels Match COC?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	12.	
-Includes Date/Time/ID/Analysis Matrix: <u>SL</u>			
All containers needing acid/base preservation have been checked? Noncompliances are noted in 13.	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	13.	<input type="checkbox"/> HNO <sub>3</sub> <input type="checkbox"/> H <sub>2</sub> SO <sub>4</sub> <input type="checkbox"/> NaOH <input type="checkbox"/> HCl
All containers needing preservation are found to be in compliance with EPA recommendation? (HNO <sub>3</sub> , H <sub>2</sub> SO <sub>4</sub> , HCl<2; NaOH>12)	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A		Sample # _____
Exceptions: VOA, Coliform, TOC, Oil and Grease, WI-DRO (water)	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No		Initial when completed: _____ Lot # of added preservative: _____
Headspace in VOA Vials (>6mm)?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	14.	
Trip Blank Present?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	15.	
Trip Blank Custody Seals Present?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A		
Pace Trip Blank Lot # (if purchased):			

**CLIENT NOTIFICATION/RESOLUTION**

**Field Data Required?**  Yes     No

**Person Contacted:** \_\_\_\_\_ **Date/Time:** \_\_\_\_\_

**Comments/Resolution:** \_\_\_\_\_

**Project Manager Review:** \_\_\_\_\_

**Date:** 12/7/12

Note: Whenever there is a discrepancy affecting North Carolina compliance samples, a copy of this form will be sent to the North Carolina DEHNR Certification Office (i.e. out of hold, incorrect preservative, out of temp, incorrect containers)

## Reporting Flags

- A = Reporting Limit based on signal to noise
- B = Less than 10x higher than method blank level
- C = Result obtained from confirmation analysis
- D = Result obtained from analysis of diluted sample
- E = Exceeds calibration range
- I = Interference present
- J = Estimated value
- Nn = Value obtained from additional analysis
- P = PCDE Interference
- R = Recovery outside target range
- S = Peak saturated
- U = Analyte not detected
- V = Result verified by confirmation analysis
- X = %D Exceeds limits
- Y = Calculated using average of daily RFs
- \* = See Discussion

### REPORT OF LABORATORY ANALYSIS

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## **Appendix B**

### Sample Analysis Summary



**Method 8290 Sample Analysis Results**

Client - Bay West, Inc.

Client's Sample ID	Spent Carbon		
Lab Sample ID	10214459001		
Filename	F121216B_10		
Injected By	BAL		
Total Amount Extracted	1.10 g	Matrix	Solid
% Moisture	44.1	Dilution	NA
Dry Weight Extracted	0.615 g	Collected	12/05/2012 11:05
ICAL ID	F121120	Received	12/06/2012 09:20
CCal Filename(s)	F121216B_08 & F121216B_21	Extracted	12/12/2012 19:30
Method Blank ID	BLANK-34872	Analyzed	12/17/2012 00:59

Native Isomers	Conc ng/Kg	EMPC ng/Kg	RL ng/Kg	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	22	—	1.2	2,3,7,8-TCDF-13C	2.00	77
Total TCDF	420	—	1.2	2,3,7,8-TCDD-13C	2.00	92
				1,2,3,7,8-PeCDF-13C	2.00	78
2,3,7,8-TCDD	220	—	1.1	2,3,4,7,8-PeCDF-13C	2.00	81
Total TCDD	23000	—	1.1 E	1,2,3,7,8-PeCDD-13C	2.00	91
				1,2,3,4,7,8-HxCDF-13C	2.00	77
1,2,3,7,8-PeCDF	130	—	2.5	1,2,3,6,7,8-HxCDF-13C	2.00	79
2,3,4,7,8-PeCDF	71	—	2.2 J	2,3,4,6,7,8-HxCDF-13C	2.00	78
Total PeCDF	780	—	2.4	1,2,3,7,8,9-HxCDF-13C	2.00	75
				1,2,3,4,7,8-HxCDD-13C	2.00	85
1,2,3,7,8-PeCDD	1300	—	2.1	1,2,3,6,7,8-HxCDD-13C	2.00	77
Total PeCDD	130000	—	2.1 E	1,2,3,4,6,7,8-HpCDF-13C	2.00	80
				1,2,3,4,7,8,9-HpCDF-13C	2.00	81
1,2,3,4,7,8-HxCDF	270	—	4.2	1,2,3,4,6,7,8-HpCDD-13C	2.00	104
1,2,3,6,7,8-HxCDF	67	—	4.5 J	OCDD-13C	4.00	66
2,3,4,6,7,8-HxCDF	48	—	5.3 J			
1,2,3,7,8,9-HxCDF	58	—	4.3 J	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	3100	—	4.6	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	74000	—	1.3 E	2,3,7,8-TCDD-37Cl4	0.20	85
1,2,3,6,7,8-HxCDD	3000	—	1.3			
1,2,3,7,8,9-HxCDD	1100	—	1.2			
Total HxCDD	150000	—	1.3 E			
1,2,3,4,6,7,8-HpCDF	2300	—	4.1	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	140	—	4.9	Equivalence: 11000 ng/Kg		
Total HpCDF	8100	—	4.5	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	110000	—	2.8 E			
Total HpCDD	150000	—	2.8 E			
OCDF	7600	—	2.1 Y			
OCDD	410000	—	2.8 E			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
J = Estimated value  
E = Exceeds calibration range  
Y = Calculated using average of daily RFs

**REPORT OF LABORATORY ANALYSIS**

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### Method 8290 Blank Analysis Results

Lab Sample ID	BLANK-34872	Matrix	Solid
Filename	F121215B_05	Dilution	NA
Total Amount Extracted	20.9 g	Extracted	12/12/2012 19:30
ICAL ID	F121120	Analyzed	12/15/2012 20:04
CCal Filename(s)	F121215B_01 & F121215B_18	Injected By	BAL

Native Isomers	Conc ng/Kg	EMPC ng/Kg	RL ng/Kg	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.051	—	0.036 J	2,3,7,8-TCDF-13C	2.00	74
Total TCDF	0.097	—	0.036 J	2,3,7,8-TCDD-13C	2.00	89
				1,2,3,7,8-PeCDF-13C	2.00	77
2,3,7,8-TCDD	ND	—	0.037	2,3,4,7,8-PeCDF-13C	2.00	77
Total TCDD	ND	—	0.037	1,2,3,7,8-PeCDD-13C	2.00	86
				1,2,3,4,7,8-HxCDF-13C	2.00	82
1,2,3,7,8-PeCDF	ND	—	0.056	1,2,3,6,7,8-HxCDF-13C	2.00	82
2,3,4,7,8-PeCDF	ND	—	0.035	2,3,4,6,7,8-HxCDF-13C	2.00	84
Total PeCDF	ND	—	0.045	1,2,3,7,8,9-HxCDF-13C	2.00	80
				1,2,3,4,7,8-HxCDD-13C	2.00	88
1,2,3,7,8-PeCDD	ND	—	0.044	1,2,3,6,7,8-HxCDD-13C	2.00	79
Total PeCDD	ND	—	0.044	1,2,3,4,6,7,8-HpCDF-13C	2.00	79
				1,2,3,4,7,8,9-HpCDF-13C	2.00	77
1,2,3,4,7,8-HxCDF	ND	—	0.030	1,2,3,4,6,7,8-HpCDD-13C	2.00	90
1,2,3,6,7,8-HxCDF	ND	—	0.029	OCDD-13C	4.00	60
2,3,4,6,7,8-HxCDF	ND	—	0.026			
1,2,3,7,8,9-HxCDF	ND	—	0.030	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	—	0.029	1,2,3,7,8,9-HxCDD-13C	2.00	NA
				2,3,7,8-TCDD-37Cl4	0.20	81
1,2,3,4,7,8-HxCDD	ND	—	0.037			
1,2,3,6,7,8-HxCDD	ND	—	0.047			
1,2,3,7,8,9-HxCDD	ND	—	0.045			
Total HxCDD	ND	—	0.043			
				Total 2,3,7,8-TCDD		
1,2,3,4,6,7,8-HpCDF	ND	—	0.023	Equivalence: 0.065 ng/Kg		
1,2,3,4,7,8,9-HpCDF	ND	—	0.027	(Using 2005 WHO Factors - Using PRL/2 where ND)		
Total HpCDF	ND	—	0.025			
1,2,3,4,6,7,8-HpCDD	—	0.056	0.037 I			
Total HpCDD	ND	—	0.037			
OCDF	ND	—	0.040			
OCDD	0.230	—	0.070 J			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).

EMPC = Estimated Maximum Possible Concentration

RL = Reporting Limit

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

I = Interference present

## REPORT OF LABORATORY ANALYSIS

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**Method 8290 Laboratory Control Spike Results**

Lab Sample ID	LCS-34873	Matrix	Solid
Filename	F121215B_02	Dilution	NA
Total Amount Extracted	22.6 g	Extracted	12/12/2012 19:30
ICAL ID	F121120	Analyzed	12/15/2012 17:52
CCal Filename(s)	F121215B_01 & F121215B_18	Injected By	BAL
Method Blank ID	BLANK-34872		

Native Isomers	Qs (ng)	Qm (ng)	% Rec.	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.20	0.21	104	2,3,7,8-TCDF-13C	2.0	78
Total TCDF				2,3,7,8-TCDD-13C	2.0	93
				1,2,3,7,8-PeCDF-13C	2.0	78
2,3,7,8-TCDD	0.20	0.16	80	2,3,4,7,8-PeCDF-13C	2.0	79
Total TCDD				1,2,3,7,8-PeCDD-13C	2.0	87
				1,2,3,4,7,8-HxCDF-13C	2.0	85
1,2,3,7,8-PeCDF	1.0	1.0	103	1,2,3,6,7,8-HxCDF-13C	2.0	87
2,3,4,7,8-PeCDF	1.0	0.99	99	2,3,4,6,7,8-HxCDF-13C	2.0	86
Total PeCDF				1,2,3,7,8,9-HxCDF-13C	2.0	82
				1,2,3,4,7,8-HxCDD-13C	2.0	89
1,2,3,7,8-PeCDD	1.0	0.88	88	1,2,3,6,7,8-HxCDD-13C	2.0	80
Total PeCDD				1,2,3,4,6,7,8-HpCDF-13C	2.0	78
				1,2,3,4,7,8,9-HpCDF-13C	2.0	74
1,2,3,4,7,8-HxCDF	1.0	1.1	105	1,2,3,4,6,7,8-HpCDD-13C	2.0	89
1,2,3,6,7,8-HxCDF	1.0	0.97	97	OCDD-13C	4.0	60
2,3,4,6,7,8-HxCDF	1.0	1.0	100			
1,2,3,7,8,9-HxCDF	1.0	0.99	99	1,2,3,4-TCDD-13C	2.0	NA
Total HxCDF				1,2,3,7,8,9-HxCDD-13C	2.0	NA
1,2,3,4,7,8-HxCDD	1.0	0.99	99	2,3,7,8-TCDD-37Cl4	0.20	84
1,2,3,6,7,8-HxCDD	1.0	1.1	110			
1,2,3,7,8,9-HxCDD	1.0	1.0	105			
Total HxCDD						
1,2,3,4,6,7,8-HpCDF	1.0	1.0	104			
1,2,3,4,7,8,9-HpCDF	1.0	0.92	92			
Total HpCDF						
1,2,3,4,6,7,8-HpCDD	1.0	0.87	87			
Total HpCDD						
OCDF	2.0	1.8	91			
OCDD	2.0	2.1	106			

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

**REPORT OF LABORATORY ANALYSIS**

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**Report Prepared for:**

Paul Walz  
Bay West, Inc.  
5 Empire Drive  
Saint Paul MN 55103

**REPORT OF  
LABORATORY  
ANALYSIS FOR  
PCDD/PCDF**

**Report Prepared Date:**

March 27, 2013

**Report Information:**

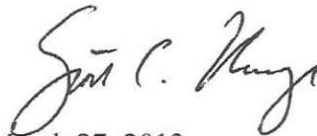
**Pace Project #: 10222116**  
**Sample Receipt Date: 03/09/2013**  
**Client Project #: FMC ESD 00439**  
**Client Sub PO #: N/A**  
**State Cert #: N/A**

**Invoicing & Reporting Options:**

The report provided has been invoiced as a Level 2 PCDD/PCDF Report. If an upgrade of this report package is requested, an additional charge may be applied.

Please review the attached invoice for accuracy and forward any questions to Scott Unze, your Pace Project Manager.

**This report has been reviewed by:**



March 27, 2013

Scott Unze, Project Manager  
(612) 607-6383  
(612) 607-6444 (fax)  
scott.unze@pacelabs.com



**Report of Laboratory Analysis**

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The results relate only to the samples included in this report.



## **DISCUSSION**

This report presents the results from the analyses performed on six samples submitted by a representative of BayWest, Inc. The samples were analyzed for the presence or absence of polychlorodibenzo-p-dioxins (PCDDs) and polychlorodibenzofurans (PCDFs) using a modified version of USEPA Method 8290. The reporting limits were based on signal-to-noise measurements.

Second column confirmation analyses of 2,3,7,8-TCDF values obtained from the primary (DB5-MS) column are performed only when specifically requested for a project and only when the values are above the concentration of the lowest calibration standard. Typical resolution for this isomer using the DB5-MS column ranges from 25-30%.

The recoveries of the isotopically-labeled PCDD/PCDF internal standards in the sample extracts ranged from 42-97%. All of the labeled standard recoveries obtained for this project were within the 40-135% target range specified in Method 8290. Also, since the quantification of the native 2,3,7,8-substituted congeners was based on isotope dilution, the data were automatically corrected for variation in recovery and accurate values were obtained.

In some cases, interfering substances impacted the determinations of PCDD or PCDF congeners; the affected values were flagged "I" where incorrect isotope ratios were obtained. Concentrations below the calibration range were flagged "J" and should be regarded as estimates. Concentrations above the calibration range were flagged "E" and should also be regarded as estimates. Values obtained from dilutions of the sample extracts were flagged "D". Saturated signals were flagged "S" and should be regarded as minimum possible concentrations.

A laboratory method blank was prepared and analyzed with each sample batch as part of our routine quality control procedures. The results show the blanks to contain trace levels of selected congeners. These levels were below the calibration range of the method. The levels reported for the affected congeners in the field samples were higher than the corresponding blank levels by one or more orders of magnitude. These results indicate that the sample processing steps did not contribute significantly to the levels reported for the field samples.

Laboratory spike samples were also prepared with the sample batches using clean sand or water that had been fortified with native standard materials. The results show that the spiked native compounds were recovered at 91-121% with relative percent differences of 0.0-6.8%. These results indicate high degrees of accuracy and precision for these determinations. Matrix spikes were prepared with the solid sample batch using sample material from a separate project; results from these analyses will be provided upon request. Matrix spikes were not prepared with the water sample batch.

## **REPORT OF LABORATORY ANALYSIS**

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## Minnesota Laboratory Certifications

Authority	Certificate #	Authority	Certificate #
Alabama	40770	Montana	92
Alaska	MN00064	Nebraska	
Arizona	AZ0014	Nevada	MN_00064_200
Arkansas	88-0680	New Jersey (NE)	MN002
California	01155CA	New Mexico	MN00064
Colorado	MN00064	New York (NEL)	11647
Connecticut	PH-0256	North Carolina	27700
EPA Region 5	WD-15J	North Dakota	R-036
EPA Region 8	8TMS-Q	Ohio	4150
Florida (NELAP)	E87605	Ohio VAP	CL101 9507
Georgia (DNR)	959	Oklahoma	D9922
Guam	959	Oregon (ELAP)	MN200001-005
Hawaii	SLD	Oregon (OREL)	MN300001-001
Idaho	MN00064	Pennsylvania	68-00563
Illinois	200012	Saipan	MP0003
Indiana	C-MN-01	South Carolina	74003001
Indiana	C-MN-01	Tennessee	2818
Iowa	368	Tennessee	02818
Kansas	E-10167	Texas	T104704192-08
Kentucky	90062	Utah (NELAP)	PAM
Louisiana	03086	Virginia	00251
Maine	2007029	Washington	C755
Maryland	322	West Virginia	9952C
Michigan	9909	Wisconsin	999407970
Minnesota	027-053-137	Wyoming	8TMS-Q
Mississippi	MN00064		

## REPORT OF LABORATORY ANALYSIS

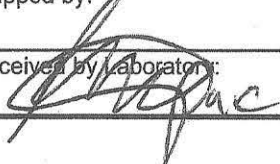
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# Appendix A

## Sample Management

### CHAIN OF CUSTODY RECORD

10222116

Project No.: <b>FMC ESD 00439</b>		Laboratory: <b>PACE ANALYTICAL</b>		Laboratory Contact: <b>SCOTT UNZE</b>		Adventus Remediation Technologies 1345 Fewster Drive Mississauga, Ontario Canada L4W 2A5 Tel: (905) 273-5374 Fax: (905) 273-4367		Analytical Laboratory to Complete												
P.O. #		Adventus Remediation Tech. Contact (Name/Tel.): <b>EVA JANZEN EXT. 232</b>						Submission Number												
Date	Time	Type		Matrix		No. of Containers	Adventus Remediation Tech. Sample Number				VOCs	SVOCs	Pesticides	Total Organic Carbon	Metals	TPH	CP	DIOXINS/FURANS (METHOD 8190)	Remarks	Lab Sample #
		Composite	Grab	Soil	Water															
<b>MAR. 8/13</b>		X		X		2	5	3	1	8	3							X		001
		X		X		2	5	3	1	8	4							X		002
		X		X		2	5	3	1	8	5							X		003
		X			X	1	5	3	1	8	6							X		004
		X			X	1	5	3	1	8	7							X		005
		X			X	1	5	3	1	8	8							X		006
Sampled by: (print name & initial) <b>SOMVEN S.O. + E. JANZEN 39</b>		Date <b>MAR. 8/13</b>		Received by: (signature) 		Date		<b>NOTES</b> Send analytical results to appropriate ART contact person.												
Relinquished by: (Signature) <b>Sandra Owen</b>		Date <b>MAR. 8/13</b>		Shipped by:		Shipping Bill														
Relinquished by: (Signature)		Date		Received by Laboratory:		Date/Time <b>3913 0914</b>														

T = 4.1



Document Name:  
**Sample Condition Upon Receipt Form**  
 Document No.:  
**F-MN-L-213-rev.06**

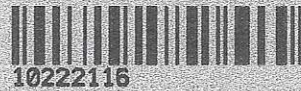
Document Revised: 28Jan2013  
 Page 1 of 1  
 Issuing Authority:  
 Pace Minnesota Quality Office

Sample Condition  
 Upon Receipt

Client Name:  
FMC

Project #:

WO#: **10222116**



Courier:  Fed Ex  UPS  USPS  Client  
 Commercial  Pace  Other: \_\_\_\_\_

Tracking Number: 794927784451

Custody Seal on Cooler/Box Present?  Yes  No Seals Intact?  Yes  No

Optional: Proj. Due Date: \_\_\_\_\_ Proj. Name: \_\_\_\_\_

Packing Material:  Bubble Wrap  Bubble Bags  None  Other: \_\_\_\_\_

Temp Blank?  Yes  No

Thermom. Used:  888A912167504  80512447  72337080 Type of Ice:  Wet  Blue  None  Samples on Ice, cooling process has begun

Cooler Temp Read (°C): 3.4 Cooler Temp Corrected (°C): 4.1 Biological Tissue Frozen?  Yes  No

Temp should be above freezing to 6°C Correction Factor: 6.7 Date and Initials of Person Examining Contents: 3913

Comments:

Chain of Custody Present?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	1.
Chain of Custody Filled Out?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	2.
Chain of Custody Relinquished?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	3.
Sampler Name and/or Signature on COC?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	4.
Samples Arrived within Hold Time?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	5.
Short Hold Time Analysis (<72 hr)?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A	6.
Rush Turn Around Time Requested?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A	7.
Sufficient Volume?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	8.
Correct Containers Used?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	9.
-Pace Containers Used?	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A	10.
Containers Intact?	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	10.
Filtered Volume Received for Dissolved Tests?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A	11.
Sample Labels Match COC? -Includes Date/Time/ID/Analysis Matrix: <u>WT/SL</u>	<input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A	12. <u>No lids on any COC put bottles in order</u>
All containers needing acid/base preservation have been checked? Noncompliances are noted in 13. All containers needing preservation are found to be in compliance with EPA recommendation? (HNO <sub>3</sub> , H <sub>2</sub> SO <sub>4</sub> , HCl<2; NaOH>12) Exceptions: VOA, Coliform, TOC, Oil and Grease, WI-DRO (water)	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A <input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No	13. <input type="checkbox"/> HNO <sub>3</sub> <input type="checkbox"/> H <sub>2</sub> SO <sub>4</sub> <input type="checkbox"/> NaOH <input type="checkbox"/> HCl Sample # Initial when completed: _____ Lot # of added preservative: _____
Headspace in VOA Vials (>6mm)?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	14.
Trip Blank Present?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	15.
Trip Blank Custody Seals Present?	<input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A	
Pace Trip Blank Lot # (if purchased):		

CLIENT NOTIFICATION/RESOLUTION

Field Data Required?  Yes  No

Person Contacted: \_\_\_\_\_

Date/Time: \_\_\_\_\_

Comments/Resolution: dis 53183, 53184 53185, 53186, 53187, 53188

Project Manager Review: \_\_\_\_\_

Date: 03/12/13

Note: Whenever there is a discrepancy affecting North Carolina compliance samples, a copy of this form will be sent to the North Carolina DEHNR Certification Office (i.e. out of hold, incorrect preservative, out of temp, incorrect containers)

## Reporting Flags

- A = Reporting Limit based on signal to noise
- B = Less than 10x higher than method blank level
- C = Result obtained from confirmation analysis
- D = Result obtained from analysis of diluted sample
- E = Exceeds calibration range
- I = Interference present
- J = Estimated value
- Nn = Value obtained from additional analysis
- P = PCDE Interference
- R = Recovery outside target range
- S = Peak saturated
- U = Analyte not detected
- V = Result verified by confirmation analysis
- X = %D Exceeds limits
- Y = Calculated using average of daily RFs
- \* = See Discussion

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## **Appendix B**

### **Sample Analysis Summary**



**Method 8290 Sample Analysis Results**

Client - Bay West, Inc.

Client's Sample ID	53183		
Lab Sample ID	10222116001		
Filename	F130318B_06		
Injected By	SMT		
Total Amount Extracted	961 mL	Matrix	Water
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	03/08/2013
ICAL ID	F130315	Received	03/09/2013 09:14
CCal Filename(s)	F130318A_18 & F130318B_10	Extracted	03/14/2013 16:00
Method Blank ID	BLANK-35740	Analyzed	03/18/2013 21:06

Native Isomers	Conc pg/L	EMPC pg/L	RL pg/L	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	—	1.70	2,3,7,8-TCDF-13C	2.00	53
Total TCDF	ND	—	1.70	2,3,7,8-TCDD-13C	2.00	65
				1,2,3,7,8-PeCDF-13C	2.00	66
2,3,7,8-TCDD	ND	—	1.40	2,3,4,7,8-PeCDF-13C	2.00	71
Total TCDD	ND	—	1.40	1,2,3,7,8-PeCDD-13C	2.00	84
				1,2,3,4,7,8-HxCDF-13C	2.00	67
1,2,3,7,8-PeCDF	ND	—	2.70	1,2,3,6,7,8-HxCDF-13C	2.00	75
2,3,4,7,8-PeCDF	ND	—	1.70	2,3,4,6,7,8-HxCDF-13C	2.00	75
Total PeCDF	2.2	—	2.20 J	1,2,3,7,8,9-HxCDF-13C	2.00	75
				1,2,3,4,7,8-HxCDD-13C	2.00	73
1,2,3,7,8-PeCDD	ND	—	1.20	1,2,3,6,7,8-HxCDD-13C	2.00	74
Total PeCDD	ND	—	1.20	1,2,3,4,6,7,8-HpCDF-13C	2.00	78
				1,2,3,4,7,8,9-HpCDF-13C	2.00	85
1,2,3,4,7,8-HxCDF	ND	—	1.90	1,2,3,4,6,7,8-HpCDD-13C	2.00	88
1,2,3,6,7,8-HxCDF	ND	—	1.90	OCDD-13C	4.00	75
2,3,4,6,7,8-HxCDF	ND	—	1.60			
1,2,3,7,8,9-HxCDF	ND	—	1.10	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	36.0	—	1.60 J	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	3.2	—	0.88 J	2,3,7,8-TCDD-37Cl4	0.20	62
1,2,3,6,7,8-HxCDD	4.9	—	0.94 J			
1,2,3,7,8,9-HxCDD	ND	—	0.95			
Total HxCDD	9.3	—	0.92 J			
1,2,3,4,6,7,8-HpCDF	140.0	—	1.90	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	6.6	—	2.10 J	Equivalence: 10 pg/L		
Total HpCDF	410.0	—	2.00	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	340.0	—	0.67			
Total HpCDD	500.0	—	0.67			
OCDF	990.0	—	1.90			
OCDD	7600.0	—	1.20			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.  
J = Estimated value

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

**REPORT OF LABORATORY ANALYSIS**

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### Method 8290 Sample Analysis Results

Client - Bay West, Inc.

Client's Sample ID	53184		
Lab Sample ID	10222116002		
Filename	F130318B_07		
Injected By	SMT		
Total Amount Extracted	956 mL	Matrix	Water
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	03/08/2013
ICAL ID	F130315	Received	03/09/2013 09:14
CCal Filename(s)	F130318A_18 & F130318B_10	Extracted	03/14/2013 16:00
Method Blank ID	BLANK-35740	Analyzed	03/18/2013 21:49

Native Isomers	Conc pg/L	EMPC pg/L	RL pg/L	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	—	1.60	2,3,7,8-TCDF-13C	2.00	42
Total TCDF	1.60	—	1.60 J	2,3,7,8-TCDD-13C	2.00	50
				1,2,3,7,8-PeCDF-13C	2.00	53
2,3,7,8-TCDD	ND	—	1.20	2,3,4,7,8-PeCDF-13C	2.00	57
Total TCDD	ND	—	1.20	1,2,3,7,8-PeCDD-13C	2.00	66
				1,2,3,4,7,8-HxCDF-13C	2.00	53
1,2,3,7,8-PeCDF	ND	—	1.40	1,2,3,6,7,8-HxCDF-13C	2.00	61
2,3,4,7,8-PeCDF	ND	—	1.20	2,3,4,6,7,8-HxCDF-13C	2.00	59
Total PeCDF	ND	—	1.30	1,2,3,7,8,9-HxCDF-13C	2.00	59
				1,2,3,4,7,8-HxCDD-13C	2.00	60
1,2,3,7,8-PeCDD	ND	—	1.20	1,2,3,6,7,8-HxCDD-13C	2.00	56
Total PeCDD	ND	—	1.20	1,2,3,4,6,7,8-HpCDF-13C	2.00	60
				1,2,3,4,7,8,9-HpCDF-13C	2.00	66
1,2,3,4,7,8-HxCDF	ND	—	1.00	1,2,3,4,6,7,8-HpCDD-13C	2.00	69
1,2,3,6,7,8-HxCDF	ND	—	1.10	OCDD-13C	4.00	59
2,3,4,6,7,8-HxCDF	ND	—	1.20			
1,2,3,7,8,9-HxCDF	ND	—	1.40	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	—	1.20	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	—	1.2	0.66 I	2,3,7,8-TCDD-37Cl4	0.20	50
1,2,3,6,7,8-HxCDD	—	1.1	0.74 I			
1,2,3,7,8,9-HxCDD	ND	—	0.74			
Total HxCDD	0.87	—	0.71 J			
1,2,3,4,6,7,8-HpCDF	8.60	—	1.00 J	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	—	1.20	Equivalence: 2.3 pg/L		
Total HpCDF	27.00	—	1.10 J	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	26.00	—	1.10 J			
Total HpCDD	41.00	—	1.10 J			
OCDF	52.00	—	1.10 J			
OCDD	480.00	—	1.00			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

J = Estimated value  
I = Interference present

## REPORT OF LABORATORY ANALYSIS

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**Method 8290 Sample Analysis Results**

Client - Bay West, Inc.

Client's Sample ID	53185		
Lab Sample ID	10222116003		
Filename	F130318B_08		
Injected By	SMT		
Total Amount Extracted	974 mL	Matrix	Water
% Moisture	NA	Dilution	NA
Dry Weight Extracted	NA	Collected	03/08/2013
ICAL ID	F130315	Received	03/09/2013 09:14
CCal Filename(s)	F130318A_18 & F130318B_10	Extracted	03/14/2013 16:00
Method Blank ID	BLANK-35740	Analyzed	03/18/2013 22:33

Native Isomers	Conc pg/L	EMPC pg/L	RL pg/L	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	—	1.40	2,3,7,8-TCDF-13C	2.00	43
Total TCDF	ND	—	1.40	2,3,7,8-TCDD-13C	2.00	53
				1,2,3,7,8-PeCDF-13C	2.00	55
2,3,7,8-TCDD	ND	—	0.95	2,3,4,7,8-PeCDF-13C	2.00	60
Total TCDD	ND	—	0.95	1,2,3,7,8-PeCDD-13C	2.00	69
				1,2,3,4,7,8-HxCDF-13C	2.00	55
1,2,3,7,8-PeCDF	ND	—	1.60	1,2,3,6,7,8-HxCDF-13C	2.00	62
2,3,4,7,8-PeCDF	ND	—	1.00	2,3,4,6,7,8-HxCDF-13C	2.00	61
Total PeCDF	ND	—	1.30	1,2,3,7,8,9-HxCDF-13C	2.00	62
				1,2,3,4,7,8-HxCDD-13C	2.00	57
1,2,3,7,8-PeCDD	ND	—	1.00	1,2,3,6,7,8-HxCDD-13C	2.00	62
Total PeCDD	ND	—	1.00	1,2,3,4,6,7,8-HpCDF-13C	2.00	63
				1,2,3,4,7,8,9-HpCDF-13C	2.00	69
1,2,3,4,7,8-HxCDF	ND	—	1.20	1,2,3,4,6,7,8-HpCDD-13C	2.00	72
1,2,3,6,7,8-HxCDF	ND	—	1.30	OCDD-13C	4.00	59
2,3,4,6,7,8-HxCDF	ND	—	1.10			
1,2,3,7,8,9-HxCDF	ND	—	1.00	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	2.8	—	1.20 J	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	—	1.10	2,3,7,8-TCDD-37Cl4	0.20	53
1,2,3,6,7,8-HxCDD	ND	—	1.10			
1,2,3,7,8,9-HxCDD	ND	—	0.92			
Total HxCDD	ND	—	1.00			
1,2,3,4,6,7,8-HpCDF	9.7	—	0.90 J	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	—	1.00	Equivalence: 2.1 pg/L		
Total HpCDF	28.0	—	0.96 J	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	24.0	—	1.60 J			
Total HpCDD	37.0	—	1.60 J			
OCDF	43.0	—	1.20 J			
OCDD	400.0	—	1.20			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.  
J = Estimated value

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

**REPORT OF LABORATORY ANALYSIS**

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### Method 8290 Sample Analysis Results

Client - Bay West, Inc.

Client's Sample ID	53186				
Lab Sample ID	10222116004				
Filename	F130320A_13				
Injected By	SMT				
Total Amount Extracted	14.7 g	Matrix	Solid		
% Moisture	60.3	Dilution	10		
Dry Weight Extracted	5.84 g	Collected	03/08/2013		
ICAL ID	F130315	Received	03/09/2013 09:14		
CCal Filename(s)	F130320A_03 & F130320A_18	Extracted	03/13/2013 15:30		
Method Blank ID	BLANK-35733	Analyzed	03/20/2013 18:43		

Native Isomers	Conc ng/Kg	EMPC ng/Kg	RL ng/Kg		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	4.7	—	2.50	JD	2,3,7,8-TCDF-13C	2.00	65 D
Total TCDF	120.0	—	2.50	D	2,3,7,8-TCDD-13C	2.00	78 D
					1,2,3,7,8-PeCDF-13C	2.00	71 D
2,3,7,8-TCDD	46.0	—	1.90	D	2,3,4,7,8-PeCDF-13C	2.00	74 D
Total TCDD	3700.0	—	1.90	D	1,2,3,7,8-PeCDD-13C	2.00	86 D
					1,2,3,4,7,8-HxCDF-13C	2.00	63 D
1,2,3,7,8-PeCDF	13.0	—	1.70	JD	1,2,3,6,7,8-HxCDF-13C	2.00	70 D
2,3,4,7,8-PeCDF	25.0	—	2.50	JD	2,3,4,6,7,8-HxCDF-13C	2.00	64 D
Total PeCDF	220.0	—	2.10	D	1,2,3,7,8,9-HxCDF-13C	2.00	63 D
					1,2,3,4,7,8-HxCDD-13C	2.00	70 D
1,2,3,7,8-PeCDD	140.0	—	0.81	D	1,2,3,6,7,8-HxCDD-13C	2.00	70 D
Total PeCDD	20000.0	—	0.81	D	1,2,3,4,6,7,8-HpCDF-13C	2.00	78 D
					1,2,3,4,7,8,9-HpCDF-13C	2.00	73 D
1,2,3,4,7,8-HxCDF	89.0	—	3.90	D	1,2,3,4,6,7,8-HpCDD-13C	2.00	84 D
1,2,3,6,7,8-HxCDF	25.0	—	2.80	JD	OCDD-13C	4.00	67 D
2,3,4,6,7,8-HxCDF	20.0	—	2.90	JD			
1,2,3,7,8,9-HxCDF	16.0	—	4.30	JD	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	920.0	—	3.40	D	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	12000.0	—	6.80	D	2,3,7,8-TCDD-37Cl4	0.20	78 D
1,2,3,6,7,8-HxCDD	1200.0	—	6.10	D			
1,2,3,7,8,9-HxCDD	480.0	—	9.90	D			
Total HxCDD	26000.0	—	7.60	D			
1,2,3,4,6,7,8-HpCDF	850.0	—	0.69	D	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	64.0	—	0.59	JD	Equivalence: 1900 ng/Kg		
Total HpCDF	2800.0	—	0.64	D	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	30000.0	—	0.77	D			
Total HpCDD	42000.0	—	0.77	D			
OCDF	3200.0	—	1.30	D			
OCDD	170000.0	—	1.50	ED			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

E = Exceeds calibration range

D = Result obtained from analysis of diluted sample

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**Method 8290 Sample Analysis Results**

Client - Bay West, Inc.

Client's Sample ID	53187			
Lab Sample ID	10222116005			
Filename	F130320A_15			
Injected By	SMT			
Total Amount Extracted	14.4 g	Matrix	Solid	
% Moisture	55.8	Dilution	50	
Dry Weight Extracted	6.36 g	Collected	03/08/2013	
ICAL ID	F130315	Received	03/09/2013 09:14	
CCal Filename(s)	F130320A_03 & F130320A_18	Extracted	03/13/2013 15:30	
Method Blank ID	BLANK-35733	Analyzed	03/20/2013 20:09	

Native Isomers	Conc ng/Kg	EMPC ng/Kg	RL ng/Kg		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	13	---	11.0	JD	2,3,7,8-TCDF-13C	2.00	69 D
Total TCDF	200	---	11.0	D	2,3,7,8-TCDD-13C	2.00	83 D
					1,2,3,7,8-PeCDF-13C	2.00	73 D
2,3,7,8-TCDD	220	---	14.0	D	2,3,4,7,8-PeCDF-13C	2.00	79 D
Total TCDD	16000	---	14.0	D	1,2,3,7,8-PeCDD-13C	2.00	93 D
					1,2,3,4,7,8-HxCDF-13C	2.00	67 D
1,2,3,7,8-PeCDF	140	---	11.0	JD	1,2,3,6,7,8-HxCDF-13C	2.00	74 D
2,3,4,7,8-PeCDF	84	---	11.0	JD	2,3,4,6,7,8-HxCDF-13C	2.00	71 D
Total PeCDF	660	---	11.0	D	1,2,3,7,8,9-HxCDF-13C	2.00	70 D
					1,2,3,4,7,8-HxCDD-13C	2.00	74 D
1,2,3,7,8-PeCDD	770	---	9.2	D	1,2,3,6,7,8-HxCDD-13C	2.00	74 D
Total PeCDD	98000	---	9.2	D	1,2,3,4,6,7,8-HpCDF-13C	2.00	77 D
					1,2,3,4,7,8,9-HpCDF-13C	2.00	83 D
1,2,3,4,7,8-HxCDF	430	---	22.0	D	1,2,3,4,6,7,8-HpCDD-13C	2.00	91 D
1,2,3,6,7,8-HxCDF	110	---	15.0	JD	OCDD-13C	4.00	62 D
2,3,4,6,7,8-HxCDF	61	---	14.0	JD			
1,2,3,7,8,9-HxCDF	67	---	14.0	JD	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	2400	---	16.0	D	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	67000	---	6.2	D	2,3,7,8-TCDD-37Cl4	0.20	92 D
1,2,3,6,7,8-HxCDD	6100	---	6.7	D			
1,2,3,7,8,9-HxCDD	2700	---	6.1	D			
Total HxCDD	140000	---	6.3	D			
1,2,3,4,6,7,8-HpCDF	1800	---	16.0	D	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	200	---	16.0	JD	Equivalence: 11000 ng/Kg		
Total HpCDF	5900	---	16.0	D	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	170000	---	9.3	ED			
Total HpCDD	230000	---	9.3	ED			
OCDF	6700	---	7.4	D			
OCDD	2700000	---	24.0	ESD			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

E = Exceeds calibration range

S = Peak saturated

D = Result obtained from analysis of diluted sample

**REPORT OF LABORATORY ANALYSIS**

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### Method 8290 Sample Analysis Results

Client - Bay West, Inc.

Client's Sample ID	53188			
Lab Sample ID	10222116006			
Filename	F130320A_14			
Injected By	SMT			
Total Amount Extracted	14.7 g	Matrix	Solid	
% Moisture	50.7	Dilution	10	
Dry Weight Extracted	7.25 g	Collected	03/08/2013	
ICAL ID	F130315	Received	03/09/2013 09:14	
CCal Filename(s)	F130320A_03 & F130320A_18	Extracted	03/13/2013 15:30	
Method Blank ID	BLANK-35733	Analyzed	03/20/2013 19:26	

Native Isomers	Conc ng/Kg	EMPC ng/Kg	RL ng/Kg		Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	4.8	—	2.50	JD	2,3,7,8-TCDF-13C	2.00	71 D
Total TCDF	83.0	—	2.50	D	2,3,7,8-TCDD-13C	2.00	83 D
					1,2,3,7,8-PeCDF-13C	2.00	75 D
2,3,7,8-TCDD	54.0	—	1.60	D	2,3,4,7,8-PeCDF-13C	2.00	77 D
Total TCDD	3300.0	—	1.60	D	1,2,3,7,8-PeCDD-13C	2.00	91 D
					1,2,3,4,7,8-HxCDF-13C	2.00	65 D
1,2,3,7,8-PeCDF	17.0	—	1.90	JD	1,2,3,6,7,8-HxCDF-13C	2.00	74 D
2,3,4,7,8-PeCDF	22.0	—	1.40	JD	2,3,4,6,7,8-HxCDF-13C	2.00	69 D
Total PeCDF	150.0	—	1.60	D	1,2,3,7,8,9-HxCDF-13C	2.00	69 D
					1,2,3,4,7,8-HxCDD-13C	2.00	73 D
1,2,3,7,8-PeCDD	170.0	—	1.60	D	1,2,3,6,7,8-HxCDD-13C	2.00	71 D
Total PeCDD	17000.0	—	1.60	D	1,2,3,4,6,7,8-HpCDF-13C	2.00	88 D
					1,2,3,4,7,8,9-HpCDF-13C	2.00	86 D
1,2,3,4,7,8-HxCDF	71.0	—	3.70	D	1,2,3,4,6,7,8-HpCDD-13C	2.00	97 D
1,2,3,6,7,8-HxCDF	23.0	—	2.10	JD	OCDD-13C	4.00	91 D
2,3,4,6,7,8-HxCDF	14.0	—	2.50	JD			
1,2,3,7,8,9-HxCDF	14.0	—	1.30	JD	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	580.0	—	2.40	D	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	9900.0	—	1.60	D	2,3,7,8-TCDD-37Cl4	0.20	80 D
1,2,3,6,7,8-HxCDD	1600.0	—	1.20	D			
1,2,3,7,8,9-HxCDD	470.0	—	1.20	D			
Total HxCDD	23000.0	—	1.30	D			
1,2,3,4,6,7,8-HpCDF	470.0	—	1.30	D	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	40.0	—	2.10	JD	Equivalence: 1800 ng/Kg		
Total HpCDF	1500.0	—	1.70	D	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	23000.0	—	1.20	D			
Total HpCDD	30000.0	—	1.20	D			
OCDF	1800.0	—	0.92	D			
OCDD	320000.0	—	1.50	ESD			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value  
E = Exceeds calibration range  
S = Peak saturated  
D = Result obtained from analysis of diluted sample

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**Method 8290 Blank Analysis Results**

Lab Sample ID	BLANK-35740	Matrix	Water
Filename	F130318B_05	Dilution	NA
Total Amount Extracted	916 mL	Extracted	03/14/2013 16:00
ICAL ID	F130315	Analyzed	03/18/2013 20:23
CCal Filename(s)	F130318A_18 & F130318B_10	Injected By	SMT

Native Isomers	Conc pg/L	EMPC pg/L	RL pg/L	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	ND	—	1.20	2,3,7,8-TCDF-13C	2.00	44
Total TCDF	ND	—	1.20	2,3,7,8-TCDD-13C	2.00	55
				1,2,3,7,8-PeCDF-13C	2.00	62
2,3,7,8-TCDD	ND	—	1.10	2,3,4,7,8-PeCDF-13C	2.00	69
Total TCDD	ND	—	1.10	1,2,3,7,8-PeCDD-13C	2.00	81
				1,2,3,4,7,8-HxCDF-13C	2.00	65
1,2,3,7,8-PeCDF	ND	—	1.20	1,2,3,6,7,8-HxCDF-13C	2.00	77
2,3,4,7,8-PeCDF	ND	—	0.81	2,3,4,6,7,8-HxCDF-13C	2.00	74
Total PeCDF	ND	—	1.00	1,2,3,7,8,9-HxCDF-13C	2.00	75
				1,2,3,4,7,8-HxCDD-13C	2.00	75
1,2,3,7,8-PeCDD	ND	—	1.00	1,2,3,6,7,8-HxCDD-13C	2.00	73
Total PeCDD	ND	—	1.00	1,2,3,4,6,7,8-HpCDF-13C	2.00	78
				1,2,3,4,7,8,9-HpCDF-13C	2.00	87
1,2,3,4,7,8-HxCDF	ND	—	0.75	1,2,3,4,6,7,8-HpCDD-13C	2.00	89
1,2,3,6,7,8-HxCDF	ND	—	0.52	OCDD-13C	4.00	77
2,3,4,6,7,8-HxCDF	ND	—	0.50			
1,2,3,7,8,9-HxCDF	ND	—	0.56	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	—	0.58	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	—	0.77	2,3,7,8-TCDD-37Cl4	0.20	55
1,2,3,6,7,8-HxCDD	ND	—	0.65			
1,2,3,7,8,9-HxCDD	ND	—	0.68			
Total HxCDD	ND	—	0.70			
1,2,3,4,6,7,8-HpCDF	ND	—	0.60	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	—	0.59	Equivalence: 1.5 pg/L		
Total HpCDF	ND	—	0.60	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	2.0	—	0.85 J			
Total HpCDD	2.0	—	0.85 J			
OCDF	—	2.0	0.96 I			
OCDD	15.0	—	2.30 J			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).

EMPC = Estimated Maximum Possible Concentration

RL = Reporting Limit

J = Estimated value

I = Interference present

**REPORT OF LABORATORY ANALYSIS**

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**Method 8290 Blank Analysis Results**

Lab Sample ID	BLANK-35733	Matrix	Solid
Filename	F130318B_13	Dilution	NA
Total Amount Extracted	20.9 g	Extracted	03/13/2013 15:30
ICAL ID	F130315	Analyzed	03/19/2013 08:15
CCal Filename(s)	F130318B_10 & F130318B_17	Injected By	SMT

Native Isomers	Conc ng/Kg	EMPC ng/Kg	RL ng/Kg	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	—	0.10	0.044 I	2,3,7,8-TCDF-13C	2.00	68
Total TCDF	0.220	—	0.044 J	2,3,7,8-TCDD-13C	2.00	78
				1,2,3,7,8-PeCDF-13C	2.00	71
2,3,7,8-TCDD	ND	—	0.054	2,3,4,7,8-PeCDF-13C	2.00	66
Total TCDD	ND	—	0.054	1,2,3,7,8-PeCDD-13C	2.00	78
				1,2,3,4,7,8-HxCDF-13C	2.00	68
1,2,3,7,8-PeCDF	ND	—	0.060	1,2,3,6,7,8-HxCDF-13C	2.00	72
2,3,4,7,8-PeCDF	ND	—	0.060	2,3,4,6,7,8-HxCDF-13C	2.00	68
Total PeCDF	ND	—	0.060	1,2,3,7,8,9-HxCDF-13C	2.00	67
				1,2,3,4,7,8-HxCDD-13C	2.00	69
1,2,3,7,8-PeCDD	ND	—	0.056	1,2,3,6,7,8-HxCDD-13C	2.00	70
Total PeCDD	ND	—	0.056	1,2,3,4,6,7,8-HpCDF-13C	2.00	69
				1,2,3,4,7,8,9-HpCDF-13C	2.00	69
1,2,3,4,7,8-HxCDF	ND	—	0.063	1,2,3,4,6,7,8-HpCDD-13C	2.00	76
1,2,3,6,7,8-HxCDF	ND	—	0.051	OCDD-13C	4.00	59
2,3,4,6,7,8-HxCDF	ND	—	0.057			
1,2,3,7,8,9-HxCDF	ND	—	0.060	1,2,3,4-TCDD-13C	2.00	NA
Total HxCDF	ND	—	0.058	1,2,3,7,8,9-HxCDD-13C	2.00	NA
1,2,3,4,7,8-HxCDD	ND	—	0.070	2,3,7,8-TCDD-37Cl4	0.20	85
1,2,3,6,7,8-HxCDD	ND	—	0.081			
1,2,3,7,8,9-HxCDD	ND	—	0.069			
Total HxCDD	0.095	—	0.073 J			
1,2,3,4,6,7,8-HpCDF	ND	—	0.079	Total 2,3,7,8-TCDD		
1,2,3,4,7,8,9-HpCDF	ND	—	0.089	Equivalence: 0.091 ng/Kg		
Total HpCDF	ND	—	0.084	(Using 2005 WHO Factors - Using PRL/2 where ND)		
1,2,3,4,6,7,8-HpCDD	—	0.11	0.094 I			
Total HpCDD	ND	—	0.094			
OCDF	ND	—	0.180			
OCDD	0.620	—	0.170 J			

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).

EMPC = Estimated Maximum Possible Concentration

RL = Reporting Limit

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J = Estimated value

I = Interference present

**REPORT OF LABORATORY ANALYSIS**

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### Method 8290 Laboratory Control Spike Results

Lab Sample ID	LCS-35741	Matrix	Water
Filename	F130318B_01	Dilution	NA
Total Amount Extracted	902 mL	Extracted	03/14/2013 16:00
ICAL ID	F130315	Analyzed	03/18/2013 17:31
CCal Filename(s)	F130318A_18 & F130318B_10	Injected By	SMT
Method Blank ID	BLANK-35740		

Native Isomers	Qs (ng)	Qm (ng)	% Rec.	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.20	0.22	111	2,3,7,8-TCDF-13C	2.0	55
Total TCDF				2,3,7,8-TCDD-13C	2.0	66
				1,2,3,7,8-PeCDF-13C	2.0	68
2,3,7,8-TCDD	0.20	0.18	91	2,3,4,7,8-PeCDF-13C	2.0	76
Total TCDD				1,2,3,7,8-PeCDD-13C	2.0	86
				1,2,3,4,7,8-HxCDF-13C	2.0	75
1,2,3,7,8-PeCDF	1.0	1.2	116	1,2,3,6,7,8-HxCDF-13C	2.0	82
2,3,4,7,8-PeCDF	1.0	1.1	108	2,3,4,6,7,8-HxCDF-13C	2.0	78
Total PeCDF				1,2,3,7,8,9-HxCDF-13C	2.0	79
				1,2,3,4,7,8-HxCDD-13C	2.0	74
1,2,3,7,8-PeCDD	1.0	0.99	99	1,2,3,6,7,8-HxCDD-13C	2.0	78
Total PeCDD				1,2,3,4,6,7,8-HpCDF-13C	2.0	81
				1,2,3,4,7,8,9-HpCDF-13C	2.0	85
1,2,3,4,7,8-HxCDF	1.0	1.1	110	OCDD-13C	4.0	74
1,2,3,6,7,8-HxCDF	1.0	1.1	106			
2,3,4,6,7,8-HxCDF	1.0	1.1	108			
1,2,3,7,8,9-HxCDF	1.0	1.1	107	1,2,3,4-TCDD-13C	2.0	NA
Total HxCDF				1,2,3,7,8,9-HxCDD-13C	2.0	NA
1,2,3,4,7,8-HxCDD	1.0	1.1	113	2,3,7,8-TCDD-37Cl4	0.20	66
1,2,3,6,7,8-HxCDD	1.0	1.2	119			
1,2,3,7,8,9-HxCDD	1.0	1.2	116			
Total HxCDD						
1,2,3,4,6,7,8-HpCDF	1.0	1.1	113			
1,2,3,4,7,8,9-HpCDF	1.0	1.0	100			
Total HpCDF						
1,2,3,4,6,7,8-HpCDD	1.0	0.98	98			
Total HpCDD						
OCDF	2.0	2.2	109			
OCDD	2.0	2.2	111			

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

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**Method 8290 Laboratory Control Spike Results**

Lab Sample ID	LCS-35734	Matrix	Solid
Filename	F130318B_16	Dilution	NA
Total Amount Extracted	20.5 g	Extracted	03/13/2013 15:30
ICAL ID	F130315	Analyzed	03/19/2013 10:24
CCal Filename(s)	F130318B_10 & F130318B_17	Injected By	SMT
Method Blank ID	BLANK-35733		

Native Isomers	Qs (ng)	Qm (ng)	% Rec.	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.20	0.24	118	2,3,7,8-TCDF-13C	2.0	68
Total TCDF				2,3,7,8-TCDD-13C	2.0	79
				1,2,3,7,8-PeCDF-13C	2.0	70
2,3,7,8-TCDD	0.20	0.18	92	2,3,4,7,8-PeCDF-13C	2.0	65
Total TCDD				1,2,3,7,8-PeCDD-13C	2.0	80
				1,2,3,4,7,8-HxCDF-13C	2.0	66
1,2,3,7,8-PeCDF	1.0	1.1	114	1,2,3,6,7,8-HxCDF-13C	2.0	72
2,3,4,7,8-PeCDF	1.0	1.1	108	2,3,4,6,7,8-HxCDF-13C	2.0	69
Total PeCDF				1,2,3,7,8,9-HxCDF-13C	2.0	67
				1,2,3,4,7,8-HxCDD-13C	2.0	71
1,2,3,7,8-PeCDD	1.0	0.97	97	1,2,3,6,7,8-HxCDD-13C	2.0	67
Total PeCDD				1,2,3,4,6,7,8-HpCDF-13C	2.0	72
				1,2,3,4,7,8,9-HpCDF-13C	2.0	71
1,2,3,4,7,8-HxCDF	1.0	1.1	113	1,2,3,4,6,7,8-HpCDD-13C	2.0	77
1,2,3,6,7,8-HxCDF	1.0	1.1	106	OCDD-13C	4.0	58
2,3,4,6,7,8-HxCDF	1.0	1.1	108			
1,2,3,7,8,9-HxCDF	1.0	1.1	105	1,2,3,4-TCDD-13C	2.0	NA
Total HxCDF				1,2,3,7,8,9-HxCDD-13C	2.0	NA
1,2,3,4,7,8-HxCDD	1.0	1.1	111	2,3,7,8-TCDD-37Cl4	0.20	82
1,2,3,6,7,8-HxCDD	1.0	1.2	116			
1,2,3,7,8,9-HxCDD	1.0	1.2	116			
Total HxCDD						
1,2,3,4,6,7,8-HpCDF	1.0	1.1	113			
1,2,3,4,7,8,9-HpCDF	1.0	1.0	101			
Total HpCDF						
1,2,3,4,6,7,8-HpCDD	1.0	1.00	100			
Total HpCDD						
OCDF	2.0	2.0	102			
OCDD	2.0	2.2	112			

Qs = Quantity Spiked  
 Qm = Quantity Measured  
 Rec. = Recovery (Expressed as Percent)  
 R = Recovery outside of target range

Y = RF averaging used in calculations  
 Nn = Value obtained from additional analysis  
 NA = Not Applicable  
 \* = See Discussion

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### Method 8290 Laboratory Control Spike Results

Lab Sample ID	LCSD-35742	Matrix	Water
Filename	F130318B_02	Dilution	NA
Total Amount Extracted	914 mL	Extracted	03/14/2013 16:00
ICAL ID	F130315	Analyzed	03/18/2013 18:13
CCal Filename(s)	F130318A_18 & F130318B_10	Injected By	SMT
Method Blank ID	BLANK-35740		

Native Isomers	Qs (ng)	Qm (ng)	% Rec.	Internal Standards	ng's Added	Percent Recovery
2,3,7,8-TCDF	0.20	0.24	118	2,3,7,8-TCDF-13C	2.0	66
Total TCDF				2,3,7,8-TCDD-13C	2.0	75
				1,2,3,7,8-PeCDF-13C	2.0	72
2,3,7,8-TCDD	0.20	0.19	94	2,3,4,7,8-PeCDF-13C	2.0	78
Total TCDD				1,2,3,7,8-PeCDD-13C	2.0	89
				1,2,3,4,7,8-HxCDF-13C	2.0	69
1,2,3,7,8-PeCDF	1.0	1.1	114	1,2,3,6,7,8-HxCDF-13C	2.0	78
2,3,4,7,8-PeCDF	1.0	1.1	108	2,3,4,6,7,8-HxCDF-13C	2.0	76
Total PeCDF				1,2,3,7,8,9-HxCDF-13C	2.0	77
				1,2,3,4,7,8-HxCDD-13C	2.0	72
1,2,3,7,8-PeCDD	1.0	0.97	97	1,2,3,6,7,8-HxCDD-13C	2.0	78
Total PeCDD				1,2,3,4,6,7,8-HpCDF-13C	2.0	76
				1,2,3,4,7,8,9-HpCDF-13C	2.0	86
1,2,3,4,7,8-HxCDF	1.0	1.2	117	1,2,3,4,6,7,8-HpCDD-13C	2.0	91
1,2,3,6,7,8-HxCDF	1.0	1.1	107	OCDD-13C	4.0	73
2,3,4,6,7,8-HxCDF	1.0	1.1	110			
1,2,3,7,8,9-HxCDF	1.0	1.1	108	1,2,3,4-TCDD-13C	2.0	NA
Total HxCDF				1,2,3,7,8,9-HxCDD-13C	2.0	NA
1,2,3,4,7,8-HxCDD	1.0	1.2	121	2,3,7,8-TCDD-37Cl4	0.20	75
1,2,3,6,7,8-HxCDD	1.0	1.1	113			
1,2,3,7,8,9-HxCDD	1.0	1.2	118			
Total HxCDD						
1,2,3,4,6,7,8-HpCDF	1.0	1.2	116			
1,2,3,4,7,8,9-HpCDF	1.0	1.0	102			
Total HpCDF						
1,2,3,4,6,7,8-HpCDD	1.0	0.98	98			
Total HpCDD						
OCDF	2.0	2.3	113			
OCDD	2.0	2.2	111			

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

## REPORT OF LABORATORY ANALYSIS

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**Method 8290**

**Spike Recovery Relative Percent Difference (RPD) Results**

Client Bay West, Inc.  
 Spike 1 ID LCS-35741 Spike 2 ID LCSD-35742  
 Spike 1 Filename F130318B\_01 Spike 2 Filename F130318B\_02

Compound	Spike 1 %REC	Spike 2 %REC	%RPD
2,3,7,8-TCDF	111	118	6.1
2,3,7,8-TCDD	91	94	3.2
1,2,3,7,8-PeCDF	116	114	1.7
2,3,4,7,8-PeCDF	108	108	0.0
1,2,3,7,8-PeCDD	99	97	2.0
1,2,3,4,7,8-HxCDF	110	117	6.2
1,2,3,6,7,8-HxCDF	106	107	0.9
2,3,4,6,7,8-HxCDF	108	110	1.8
1,2,3,7,8,9-HxCDF	107	108	0.9
1,2,3,4,7,8-HxCDD	113	121	6.8
1,2,3,6,7,8-HxCDD	119	113	5.2
1,2,3,7,8,9-HxCDD	116	118	1.7
1,2,3,4,6,7,8-HpCDF	113	116	2.6
1,2,3,4,7,8,9-HpCDF	100	102	2.0
1,2,3,4,6,7,8-HpCDD	98	98	0.0
OCDF	109	113	3.6
OCDD	111	111	0.0

%REC = Percent Recovered

RPD = The difference between the two values divided by the mean value

**REPORT OF LABORATORY ANALYSIS**

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March 25, 2013

Bay West  
Attn: Mr. Paul Walz  
5 Empire Dr.  
St. Paul, MN 55103

**Project: Bench Scale Test**

Dear Mr. Paul Walz,

Enclosed is a copy of the laboratory report for the following work order(s) received by TriMatrix Laboratories:

<b>Work Order</b>	<b>Received</b>	<b>Description</b>
1303154	03/09/2013	FMC ESD 00439

This report relates only to the sample(s) as received. Test results are in compliance with the requirements of the National Environmental Laboratory Accreditation Program (NELAP) and/or one of the following certification programs:

ACLASS DoD-ELAP/ISO17025 (#ADE-1542); Arkansas DEP (#12-056-0); Florida DEP (#E87622-24); Georgia EPD (#E87622-24); Illinois DEP (#002841); Kansas DPH (#E-10302); Kentucky DEP (#0021); Louisiana DEP (#03068); Michigan DPH (#0034); Minnesota DPH (#367345); New York ELAP (#46503); North Carolina DNRE (#659); Texas CEQ (#T104704495-12-2); Virginia DCLS (#1622); Wisconsin DNR (#999472650); USDA Soil Import Permit (#P330-09-00163).

Any qualification or narration of results, including sample acceptance requirements and test exceptions to the above referenced programs, is presented in the Statement of Data Qualifications section of this report. Estimates of analytical uncertainties and certification documents for the test results contained within this report are available upon request.

If you have any questions or require further information, please do not hesitate to contact me.

Sincerely,



James D. McFadden  
Project Chemist

**ANALYTICAL REPORT**

Client: **Bay West**  
 Project: Bench Scale Test  
 Client Sample ID: **53183**  
 Lab Sample ID: **1303154-01**  
 Matrix: Water  
 Unit: ug/L  
 Dilution Factor: 1  
 QC Batch: 1302297

Work Order: **1303154**  
 Description: FMC ESD 00439  
 Sampled: 03/08/13 12:00  
 Sampled By: Client  
 Received: 03/09/13 09:00  
 Prepared: 03/15/13 By: SMS9  
 Analyzed: 03/20/13 By: JLB  
 Analytical Batch: 3C21059

**Semivolatile Organic Compounds by EPA Method 8270C**

CAS Number	Analyte	Analytical Result	RL
83-32-9	Acenaphthene	<0.50	0.50
208-96-8	Acenaphthylene	<0.50	0.50
98-86-2	Acetophenone	<0.50	0.50
120-12-7	Anthracene	<0.50	0.50
1912-24-9	Atrazine	<0.50	0.50
100-52-7	Benzaldehyde	<0.50	0.50
56-55-3	Benzo(a)anthracene	<0.50	0.50
50-32-8	Benzo(a)pyrene	<0.50	0.50
205-99-2	Benzo(b)fluoranthene	<0.50	0.50
207-08-9	Benzo(k)fluoranthene	<0.50	0.50
191-24-2	Benzo(g,h,i)perylene	<0.50	0.50
92-52-4	1,1'-Biphenyl	<0.50	0.50
101-55-3	4-Bromophenyl Phenyl Ether	<0.50	0.50
85-68-7	Butyl Benzyl Phthalate	<1.0	1.0
105-60-2	Caprolactam	<1.0	1.0
86-74-8	Carbazole	<0.50	0.50
59-50-7	4-Chloro-3-methylphenol	<0.50	0.50
106-47-8	4-Chloroaniline	<1.0	1.0
111-91-1	Bis(2-chloroethoxy)methane	<0.50	0.50
111-44-4	Bis(2-chloroethyl) Ether	<0.50	0.50
108-60-1	Bis(2-chloroisopropyl) Ether	<0.50	0.50
91-58-7	2-Chloronaphthalene	<0.50	0.50
95-57-8	2-Chlorophenol	<0.50	0.50
7005-72-3	4-Chlorophenyl Phenyl Ether	<0.50	0.50
218-01-9	Chrysene	<0.50	0.50
53-70-3	Dibenz(a,h)anthracene	<0.50	0.50
132-64-9	Dibenzofuran	<0.50	0.50
84-74-2	Di-n-butyl Phthalate	<b>4.9</b>	1.0
91-94-1	3,3'-Dichlorobenzidine	<10	10
120-83-2	2,4-Dichlorophenol	<0.50	0.50
84-66-2	Diethyl Phthalate	<0.50	0.50

Continued on next page

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53183</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-01</b>	Sampled By:	Client
Matrix:	Water	Received:	03/09/13 09:00
Unit:	ug/L	Prepared:	03/15/13 By: SMS9
Dilution Factor:	1	Analyzed:	03/20/13 By: JLB
QC Batch:	1302297	Analytical Batch:	3C21059

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
105-67-9	2,4-Dimethylphenol	<1.0	1.0
131-11-3	Dimethyl Phthalate	<0.50	0.50
534-52-1	4,6-Dinitro-2-methylphenol	<5.0	5.0
51-28-5	2,4-Dinitrophenol	<5.0	5.0
121-14-2	2,4-Dinitrotoluene	<0.50	0.50
606-20-2	2,6-Dinitrotoluene	<0.50	0.50
117-84-0	Di-n-octyl Phthalate	<0.50	0.50
117-81-7	Bis(2-ethylhexyl) Phthalate	<b>0.55</b>	0.50
206-44-0	Fluoranthene	<0.50	0.50
86-73-7	Fluorene	<0.50	0.50
118-74-1	Hexachlorobenzene	<0.50	0.50
87-68-3	Hexachlorobutadiene	<0.50	0.50
77-47-4	Hexachlorocyclopentadiene	<0.50	0.50
67-72-1	Hexachloroethane	<0.50	0.50
193-39-5	Indeno(1,2,3-cd)pyrene	<0.50	0.50
78-59-1	Isophorone	<0.50	0.50
91-57-6	2-Methylnaphthalene	<0.50	0.50
95-48-7	2-Methylphenol	<0.50	0.50
106-44-5	4-Methylphenol	<0.50	0.50
91-20-3	Naphthalene	<0.50	0.50
88-74-4	2-Nitroaniline	<0.50	0.50
99-09-2	3-Nitroaniline	<1.0	1.0
100-01-6	4-Nitroaniline	<1.0	1.0
98-95-3	Nitrobenzene	<0.50	0.50
100-02-7	4-Nitrophenol	<5.0	5.0
88-75-5	2-Nitrophenol	<0.50	0.50
86-30-6	N-Nitroso-diphenylamine	<0.50	0.50
621-64-7	N-Nitroso-di-n-propylamine	<0.50	0.50
87-86-5	Pentachlorophenol	<0.50	0.50
85-01-8	Phenanthrene	<0.50	0.50
108-95-2	Phenol	<0.50	0.50

Continued on next page



**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53183</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-01</b>	Sampled By:	Client
Matrix:	Water	Received:	03/09/13 09:00
Unit:	ug/L	Prepared:	03/15/13 By: SMS9
Dilution Factor:	1	Analyzed:	03/20/13 By: JLB
QC Batch:	1302297	Analytical Batch:	3C21059

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
129-00-0	Pyrene	<0.50	0.50
95-94-3	1,2,4,5-Tetrachlorobenzene	<2.0	2.0
58-90-2	2,3,4,6-Tetrachlorophenol	<5.0	5.0
88-06-2	2,4,6-Trichlorophenol	<0.50	0.50
95-95-4	2,4,5-Trichlorophenol	<0.50	0.50
<b>Surrogates:</b>		<b>% Recovery</b>	<b>Control Limits</b>
	<i>2-Fluorophenol</i>	<i>39</i>	<i>20-70</i>
	<i>Phenol-d6</i>	<i>25</i>	<i>18-45</i>
	<i>Nitrobenzene-d5</i>	<i>77</i>	<i>31-123</i>
	<i>2-Fluorobiphenyl</i>	<i>72</i>	<i>25-113</i>
	<i>2,4,6-Tribromophenol</i>	<i>71</i>	<i>30-121</i>
	<i>o-Terphenyl</i>	<i>80</i>	<i>42-125</i>

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53184</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-02</b>	Sampled By:	Client
Matrix:	Water	Received:	03/09/13 09:00
Unit:	ug/L	Prepared:	03/15/13 By: SMS9
Dilution Factor:	1	Analyzed:	03/20/13 By: JLB
QC Batch:	1302297	Analytical Batch:	3C21059

**Semivolatile Organic Compounds by EPA Method 8270C**

CAS Number	Analyte	Analytical Result	RL
83-32-9	Acenaphthene	<0.50	0.50
208-96-8	Acenaphthylene	<0.50	0.50
98-86-2	Acetophenone	<0.50	0.50
120-12-7	Anthracene	<0.50	0.50
1912-24-9	Atrazine	<0.50	0.50
100-52-7	Benzaldehyde	<b>2.1</b>	0.50
56-55-3	Benzo(a)anthracene	<0.50	0.50
50-32-8	Benzo(a)pyrene	<0.50	0.50
205-99-2	Benzo(b)fluoranthene	<0.50	0.50
207-08-9	Benzo(k)fluoranthene	<0.50	0.50
191-24-2	Benzo(g,h,i)perylene	<0.50	0.50
92-52-4	1,1'-Biphenyl	<0.50	0.50
101-55-3	4-Bromophenyl Phenyl Ether	<0.50	0.50
85-68-7	Butyl Benzyl Phthalate	<1.0	1.0
105-60-2	Caprolactam	<1.0	1.0
86-74-8	Carbazole	<0.50	0.50
59-50-7	4-Chloro-3-methylphenol	<0.50	0.50
106-47-8	4-Chloroaniline	<1.0	1.0
111-91-1	Bis(2-chloroethoxy)methane	<0.50	0.50
111-44-4	Bis(2-chloroethyl) Ether	<0.50	0.50
108-60-1	Bis(2-chloroisopropyl) Ether	<0.50	0.50
91-58-7	2-Chloronaphthalene	<0.50	0.50
95-57-8	2-Chlorophenol	<0.50	0.50
7005-72-3	4-Chlorophenyl Phenyl Ether	<0.50	0.50
218-01-9	Chrysene	<0.50	0.50
53-70-3	Dibenz(a,h)anthracene	<0.50	0.50
132-64-9	Dibenzofuran	<0.50	0.50
84-74-2	Di-n-butyl Phthalate	<b>5.7</b>	1.0
91-94-1	3,3'-Dichlorobenzidine	<10	10
120-83-2	2,4-Dichlorophenol	<0.50	0.50
84-66-2	Diethyl Phthalate	<0.50	0.50

Continued on next page

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53184</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-02</b>	Sampled By:	Client
Matrix:	Water	Received:	03/09/13 09:00
Unit:	ug/L	Prepared:	03/15/13 By: SMS9
Dilution Factor:	1	Analyzed:	03/20/13 By: JLB
QC Batch:	1302297	Analytical Batch:	3C21059

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
105-67-9	2,4-Dimethylphenol	<1.0	1.0
131-11-3	Dimethyl Phthalate	<0.50	0.50
534-52-1	4,6-Dinitro-2-methylphenol	<5.0	5.0
51-28-5	2,4-Dinitrophenol	<5.0	5.0
121-14-2	2,4-Dinitrotoluene	<0.50	0.50
606-20-2	2,6-Dinitrotoluene	<0.50	0.50
117-84-0	Di-n-octyl Phthalate	<0.50	0.50
117-81-7	Bis(2-ethylhexyl) Phthalate	<b>2.3</b>	0.50
206-44-0	Fluoranthene	<0.50	0.50
86-73-7	Fluorene	<0.50	0.50
118-74-1	Hexachlorobenzene	<0.50	0.50
87-68-3	Hexachlorobutadiene	<0.50	0.50
77-47-4	Hexachlorocyclopentadiene	<0.50	0.50
67-72-1	Hexachloroethane	<0.50	0.50
193-39-5	Indeno(1,2,3-cd)pyrene	<0.50	0.50
78-59-1	Isophorone	<0.50	0.50
91-57-6	2-Methylnaphthalene	<0.50	0.50
95-48-7	2-Methylphenol	<0.50	0.50
106-44-5	4-Methylphenol	<0.50	0.50
91-20-3	Naphthalene	<0.50	0.50
88-74-4	2-Nitroaniline	<0.50	0.50
99-09-2	3-Nitroaniline	<1.0	1.0
100-01-6	4-Nitroaniline	<1.0	1.0
98-95-3	Nitrobenzene	<0.50	0.50
100-02-7	4-Nitrophenol	<5.0	5.0
88-75-5	2-Nitrophenol	<0.50	0.50
86-30-6	N-Nitroso-diphenylamine	<0.50	0.50
621-64-7	N-Nitroso-di-n-propylamine	<0.50	0.50
87-86-5	Pentachlorophenol	<b>0.87</b>	0.50
85-01-8	Phenanthrene	<0.50	0.50
108-95-2	Phenol	<0.50	0.50

Continued on next page

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53184</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-02</b>	Sampled By:	Client
Matrix:	Water	Received:	03/09/13 09:00
Unit:	ug/L	Prepared:	03/15/13 By: SMS9
Dilution Factor:	1	Analyzed:	03/20/13 By: JLB
QC Batch:	1302297	Analytical Batch:	3C21059

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
129-00-0	Pyrene	<0.50	0.50
95-94-3	1,2,4,5-Tetrachlorobenzene	<2.0	2.0
58-90-2	2,3,4,6-Tetrachlorophenol	<5.0	5.0
88-06-2	2,4,6-Trichlorophenol	<0.50	0.50
95-95-4	2,4,5-Trichlorophenol	<0.50	0.50
<b>Surrogates:</b>			
		<b>% Recovery</b>	<b>Control Limits</b>
	<i>2-Fluorophenol</i>	36	20-70
	<i>Phenol-d6</i>	25	18-45
	<i>Nitrobenzene-d5</i>	50	31-123
	<i>2-Fluorobiphenyl</i>	41	25-113
	<i>2,4,6-Tribromophenol</i>	49	30-121
	<i>o-Terphenyl</i>	53	42-125

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53185</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-03</b>	Sampled By:	Client
Matrix:	Water	Received:	03/09/13 09:00
Unit:	ug/L	Prepared:	03/15/13 By: SMS9
Dilution Factor:	1	Analyzed:	03/20/13 By: JLB
QC Batch:	1302297	Analytical Batch:	3C21059

**Semivolatile Organic Compounds by EPA Method 8270C**

CAS Number	Analyte	Analytical Result	RL
83-32-9	Acenaphthene	<0.50	0.50
208-96-8	Acenaphthylene	<0.50	0.50
98-86-2	Acetophenone	<0.50	0.50
120-12-7	Anthracene	<0.50	0.50
1912-24-9	Atrazine	<0.50	0.50
100-52-7	Benzaldehyde	<b>4.0</b>	0.50
56-55-3	Benzo(a)anthracene	<0.50	0.50
50-32-8	Benzo(a)pyrene	<0.50	0.50
205-99-2	Benzo(b)fluoranthene	<0.50	0.50
207-08-9	Benzo(k)fluoranthene	<0.50	0.50
191-24-2	Benzo(g,h,i)perylene	<0.50	0.50
92-52-4	1,1'-Biphenyl	<0.50	0.50
101-55-3	4-Bromophenyl Phenyl Ether	<0.50	0.50
85-68-7	Butyl Benzyl Phthalate	<1.0	1.0
105-60-2	Caprolactam	<1.0	1.0
86-74-8	Carbazole	<0.50	0.50
59-50-7	4-Chloro-3-methylphenol	<0.50	0.50
106-47-8	4-Chloroaniline	<1.0	1.0
111-91-1	Bis(2-chloroethoxy)methane	<0.50	0.50
111-44-4	Bis(2-chloroethyl) Ether	<0.50	0.50
108-60-1	Bis(2-chloroisopropyl) Ether	<0.50	0.50
91-58-7	2-Chloronaphthalene	<0.50	0.50
95-57-8	2-Chlorophenol	<0.50	0.50
7005-72-3	4-Chlorophenyl Phenyl Ether	<0.50	0.50
218-01-9	Chrysene	<0.50	0.50
53-70-3	Dibenz(a,h)anthracene	<0.50	0.50
132-64-9	Dibenzofuran	<0.50	0.50
84-74-2	Di-n-butyl Phthalate	<b>4.1</b>	1.0
91-94-1	3,3'-Dichlorobenzidine	<10	10
120-83-2	2,4-Dichlorophenol	<0.50	0.50
84-66-2	Diethyl Phthalate	<0.50	0.50

Continued on next page

**ANALYTICAL REPORT**

Client: **Bay West**  
 Project: Bench Scale Test  
 Client Sample ID: **53185**  
 Lab Sample ID: **1303154-03**  
 Matrix: Water  
 Unit: ug/L  
 Dilution Factor: 1  
 QC Batch: 1302297

Work Order: **1303154**  
 Description: FMC ESD 00439  
 Sampled: 03/08/13 12:00  
 Sampled By: Client  
 Received: 03/09/13 09:00  
 Prepared: 03/15/13 By: SMS9  
 Analyzed: 03/20/13 By: JLB  
 Analytical Batch: 3C21059

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
105-67-9	2,4-Dimethylphenol	<1.0	1.0
131-11-3	Dimethyl Phthalate	<0.50	0.50
534-52-1	4,6-Dinitro-2-methylphenol	<5.0	5.0
51-28-5	2,4-Dinitrophenol	<5.0	5.0
121-14-2	2,4-Dinitrotoluene	<0.50	0.50
606-20-2	2,6-Dinitrotoluene	<0.50	0.50
117-84-0	Di-n-octyl Phthalate	<0.50	0.50
117-81-7	Bis(2-ethylhexyl) Phthalate	<b>23</b>	0.50
206-44-0	Fluoranthene	<0.50	0.50
86-73-7	Fluorene	<0.50	0.50
118-74-1	Hexachlorobenzene	<0.50	0.50
87-68-3	Hexachlorobutadiene	<0.50	0.50
77-47-4	Hexachlorocyclopentadiene	<0.50	0.50
67-72-1	Hexachloroethane	<0.50	0.50
193-39-5	Indeno(1,2,3-cd)pyrene	<0.50	0.50
78-59-1	Isophorone	<0.50	0.50
91-57-6	2-Methylnaphthalene	<0.50	0.50
95-48-7	2-Methylphenol	<0.50	0.50
106-44-5	4-Methylphenol	<0.50	0.50
91-20-3	Naphthalene	<0.50	0.50
88-74-4	2-Nitroaniline	<0.50	0.50
99-09-2	3-Nitroaniline	<1.0	1.0
100-01-6	4-Nitroaniline	<1.0	1.0
98-95-3	Nitrobenzene	<0.50	0.50
100-02-7	4-Nitrophenol	<5.0	5.0
88-75-5	2-Nitrophenol	<0.50	0.50
86-30-6	N-Nitroso-diphenylamine	<0.50	0.50
621-64-7	N-Nitroso-di-n-propylamine	<0.50	0.50
87-86-5	Pentachlorophenol	<b>0.77</b>	0.50
85-01-8	Phenanthrene	<0.50	0.50
108-95-2	Phenol	<0.50	0.50

Continued on next page

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53185</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-03</b>	Sampled By:	Client
Matrix:	Water	Received:	03/09/13 09:00
Unit:	ug/L	Prepared:	03/15/13 By: SMS9
Dilution Factor:	1	Analyzed:	03/20/13 By: JLB
QC Batch:	1302297	Analytical Batch:	3C21059

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
129-00-0	Pyrene	<0.50	0.50
95-94-3	1,2,4,5-Tetrachlorobenzene	<2.0	2.0
58-90-2	2,3,4,6-Tetrachlorophenol	<5.0	5.0
88-06-2	2,4,6-Trichlorophenol	<0.50	0.50
95-95-4	2,4,5-Trichlorophenol	<0.50	0.50
<b>Surrogates:</b>		<b>% Recovery</b>	<b>Control Limits</b>
	<i>2-Fluorophenol</i>	<i>31</i>	<i>20-70</i>
	<i>Phenol-d6</i>	<i>23</i>	<i>18-45</i>
	<i>Nitrobenzene-d5</i>	<i>52</i>	<i>31-123</i>
	<i>2-Fluorobiphenyl</i>	<i>40</i>	<i>25-113</i>
	<i>2,4,6-Tribromophenol</i>	<i>52</i>	<i>30-121</i>
	<i>o-Terphenyl</i>	<i>52</i>	<i>42-125</i>

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53186</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-04</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00
Unit:	mg/kg dry	Prepared:	03/19/13 By: JTS
Dilution Factor:	2	Analyzed:	03/21/13 By: JLB
QC Batch:	1302404	Analytical Batch:	3C21065
Percent Solids:	41		

**Semivolatile Organic Compounds by EPA Method 8270C**

CAS Number	Analyte	Analytical Result	RL
83-32-9	Acenaphthene	<0.081	0.081
208-96-8	Acenaphthylene	<0.081	0.081
98-86-2	Acetophenone	<0.081	0.081
120-12-7	Anthracene	<0.081	0.081
1912-24-9	Atrazine	<0.081	0.081
100-52-7	Benzaldehyde	<0.32	0.32
56-55-3	Benzo(a)anthracene	<0.081	0.081
50-32-8	Benzo(a)pyrene	<0.081	0.081
205-99-2	Benzo(b)fluoranthene	<0.081	0.081
207-08-9	Benzo(k)fluoranthene	<0.081	0.081
191-24-2	Benzo(g,h,i)perylene	<0.16	0.16
65-85-0	Benzoic Acid	<3.2	3.2
100-51-6	Benzyl Alcohol	<0.081	0.081
92-52-4	1,1'-Biphenyl	<0.081	0.081
101-55-3	4-Bromophenyl Phenyl Ether	<0.081	0.081
85-68-7	Butyl Benzyl Phthalate	<0.16	0.16
105-60-2	Caprolactam	<0.32	0.32
86-74-8	Carbazole	<0.81	0.81
59-50-7	4-Chloro-3-methylphenol	<0.081	0.081
106-47-8	4-Chloroaniline	<0.32	0.32
111-91-1	Bis(2-chloroethoxy)methane	<0.081	0.081
111-44-4	Bis(2-chloroethyl) Ether	<0.081	0.081
108-60-1	Bis(2-chloroisopropyl) Ether	<0.081	0.081
91-58-7	2-Chloronaphthalene	<0.081	0.081
95-57-8	2-Chlorophenol	<0.081	0.081
7005-72-3	4-Chlorophenyl Phenyl Ether	<0.081	0.081
218-01-9	Chrysene	<0.081	0.081
53-70-3	Dibenz(a,h)anthracene	<0.16	0.16
132-64-9	Dibenzofuran	<0.081	0.081
84-74-2	Di-n-butyl Phthalate	<0.32	0.32
91-94-1	3,3'-Dichlorobenzidine	<4.0	4.0

Continued on next page



### ANALYTICAL REPORT

Client:	Bay West	Work Order:	1303154
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	53186	Sampled:	03/08/13 12:00
Lab Sample ID:	1303154-04	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00
Unit:	mg/kg dry	Prepared:	03/19/13 By: JTS
Dilution Factor:	2	Analyzed:	03/21/13 By: JLB
QC Batch:	1302404	Analytical Batch:	3C21065
Percent Solids:	41		

#### Semivolatile Organic Compounds by EPA Method 8270C (Continued)

CAS Number	Analyte	Analytical Result	RL
120-83-2	2,4-Dichlorophenol	<0.16	0.16
84-66-2	Diethyl Phthalate	<0.081	0.081
105-67-9	2,4-Dimethylphenol	<0.81	0.81
131-11-3	Dimethyl Phthalate	<0.081	0.081
534-52-1	4,6-Dinitro-2-methylphenol	<0.81	0.81
51-28-5	2,4-Dinitrophenol	<0.81	0.81
606-20-2	2,6-Dinitrotoluene	<0.081	0.081
121-14-2	2,4-Dinitrotoluene	<0.16	0.16
117-84-0	Di-n-octyl Phthalate	<0.081	0.081
117-81-7	Bis(2-ethylhexyl) Phthalate	<b>0.41</b>	0.16
206-44-0	Fluoranthene	<0.081	0.081
86-73-7	Fluorene	<0.16	0.16
118-74-1	Hexachlorobenzene	<0.081	0.081
87-68-3	Hexachlorobutadiene	<0.081	0.081
77-47-4	Hexachlorocyclopentadiene	<0.081	0.081
67-72-1	Hexachloroethane	<0.081	0.081
193-39-5	Indeno(1,2,3-cd)pyrene	<0.16	0.16
78-59-1	Isophorone	<0.081	0.081
91-57-6	2-Methylnaphthalene	<0.081	0.081
106-44-5	4-Methylphenol	<0.081	0.081
95-48-7	2-Methylphenol	<0.081	0.081
91-20-3	Naphthalene	<0.081	0.081
100-01-6	4-Nitroaniline	<0.16	0.16
88-74-4	2-Nitroaniline	<0.081	0.081
99-09-2	3-Nitroaniline	<0.16	0.16
98-95-3	Nitrobenzene	<0.081	0.081
88-75-5	2-Nitrophenol	<0.081	0.081
100-02-7	4-Nitrophenol	<3.2	3.2
86-30-6	N-Nitroso-diphenylamine	<0.081	0.081
621-64-7	N-Nitroso-di-n-propylamine	<0.081	0.081
87-86-5	Pentachlorophenol	<b>2.4</b>	0.81

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**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53186</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-04</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00
Unit:	mg/kg dry	Prepared:	03/19/13 By: JTS
Dilution Factor:	2	Analyzed:	03/21/13 By: JLB
QC Batch:	1302404	Analytical Batch:	3C21065
Percent Solids:	41		

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
85-01-8	Phenanthrene	<0.081	0.081
108-95-2	Phenol	<0.81	0.81
129-00-0	Pyrene	<0.081	0.081
95-94-3	1,2,4,5-Tetrachlorobenzene	<0.16	0.16
58-90-2	2,3,4,6-Tetrachlorophenol	<0.16	0.16
120-82-1	1,2,4-Trichlorobenzene	<0.081	0.081
95-95-4	2,4,5-Trichlorophenol	<0.081	0.081
88-06-2	2,4,6-Trichlorophenol	<0.081	0.081
<b>Surrogates:</b>			
		<b>% Recovery</b>	<b>Control Limits</b>
	<i>2-Fluorophenol</i>	71	33-113
	<i>Phenol-d6</i>	72	30-115
	<i>Nitrobenzene-d5</i>	67	33-131
	<i>2-Fluorobiphenyl</i>	74	46-122
	<i>2,4,6-Tribromophenol</i>	70	12-124
	<i>o-Terphenyl</i>	75	20-155

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53186</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-04</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00

**Physical/Chemical Parameters by EPA/APHA/ASTM Methods**

Analyte	Analytical Result	RL	Unit	Dilution Factor	Method	Date Time Analyzed	By	QC Batch
Percent Solids	41	0.1	%	1	USEPA-3550C	03/19/13 11:00	BAR	1302409

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53187</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-05</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00
Unit:	mg/kg dry	Prepared:	03/19/13 By: JTS
Dilution Factor:	4	Analyzed:	03/21/13 By: JLB
QC Batch:	1302404	Analytical Batch:	3C21065
Percent Solids:	45		

**Semivolatile Organic Compounds by EPA Method 8270C**

CAS Number	Analyte	Analytical Result	RL
83-32-9	Acenaphthene	<0.15	0.15
208-96-8	Acenaphthylene	<0.15	0.15
98-86-2	Acetophenone	<0.15	0.15
120-12-7	Anthracene	<0.15	0.15
1912-24-9	Atrazine	<0.15	0.15
100-52-7	Benzaldehyde	<0.59	0.59
56-55-3	Benzo(a)anthracene	<0.15	0.15
50-32-8	Benzo(a)pyrene	<0.15	0.15
205-99-2	Benzo(b)fluoranthene	<0.15	0.15
207-08-9	Benzo(k)fluoranthene	<0.15	0.15
191-24-2	Benzo(g,h,i)perylene	<0.29	0.29
65-85-0	Benzoic Acid	<5.9	5.9
100-51-6	Benzyl Alcohol	<0.15	0.15
92-52-4	1,1'-Biphenyl	<0.15	0.15
101-55-3	4-Bromophenyl Phenyl Ether	<0.15	0.15
85-68-7	Butyl Benzyl Phthalate	<0.29	0.29
105-60-2	Caprolactam	<0.59	0.59
86-74-8	Carbazole	<1.5	1.5
59-50-7	4-Chloro-3-methylphenol	<0.15	0.15
106-47-8	4-Chloroaniline	<0.59	0.59
111-91-1	Bis(2-chloroethoxy)methane	<0.15	0.15
111-44-4	Bis(2-chloroethyl) Ether	<0.15	0.15
108-60-1	Bis(2-chloroisopropyl) Ether	<0.15	0.15
91-58-7	2-Chloronaphthalene	<0.15	0.15
95-57-8	2-Chlorophenol	<0.15	0.15
7005-72-3	4-Chlorophenyl Phenyl Ether	<0.15	0.15
218-01-9	Chrysene	<0.15	0.15
53-70-3	Dibenz(a,h)anthracene	<0.29	0.29
132-64-9	Dibenzofuran	<0.15	0.15
84-74-2	Di-n-butyl Phthalate	<0.59	0.59
91-94-1	3,3'-Dichlorobenzidine	<7.3	7.3

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**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53187</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-05</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00
Unit:	mg/kg dry	Prepared:	03/19/13 By: JTS
Dilution Factor:	4	Analyzed:	03/21/13 By: JLB
QC Batch:	1302404	Analytical Batch:	3C21065
Percent Solids:	45		

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
120-83-2	2,4-Dichlorophenol	<0.29	0.29
84-66-2	Diethyl Phthalate	<0.15	0.15
105-67-9	2,4-Dimethylphenol	<1.5	1.5
131-11-3	Dimethyl Phthalate	<0.15	0.15
534-52-1	4,6-Dinitro-2-methylphenol	<1.5	1.5
51-28-5	2,4-Dinitrophenol	<1.5	1.5
606-20-2	2,6-Dinitrotoluene	<0.15	0.15
121-14-2	2,4-Dinitrotoluene	<0.29	0.29
117-84-0	Di-n-octyl Phthalate	<0.15	0.15
117-81-7	Bis(2-ethylhexyl) Phthalate	<b>2.9</b>	0.29
206-44-0	Fluoranthene	<0.15	0.15
86-73-7	Fluorene	<0.29	0.29
118-74-1	Hexachlorobenzene	<0.15	0.15
87-68-3	Hexachlorobutadiene	<0.15	0.15
77-47-4	Hexachlorocyclopentadiene	<0.15	0.15
67-72-1	Hexachloroethane	<0.15	0.15
193-39-5	Indeno(1,2,3-cd)pyrene	<0.29	0.29
78-59-1	Isophorone	<0.15	0.15
91-57-6	2-Methylnaphthalene	<0.15	0.15
106-44-5	4-Methylphenol	<0.15	0.15
95-48-7	2-Methylphenol	<0.15	0.15
91-20-3	Naphthalene	<0.15	0.15
100-01-6	4-Nitroaniline	<0.29	0.29
88-74-4	2-Nitroaniline	<0.15	0.15
99-09-2	3-Nitroaniline	<0.29	0.29
98-95-3	Nitrobenzene	<0.15	0.15
88-75-5	2-Nitrophenol	<0.15	0.15
100-02-7	4-Nitrophenol	<5.9	5.9
86-30-6	N-Nitroso-diphenylamine	<0.15	0.15
621-64-7	N-Nitroso-di-n-propylamine	<0.15	0.15
87-86-5	Pentachlorophenol	<b>3.4</b>	1.5

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**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53187</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-05</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00
Unit:	mg/kg dry	Prepared:	03/19/13 By: JTS
Dilution Factor:	4	Analyzed:	03/21/13 By: JLB
QC Batch:	1302404	Analytical Batch:	3C21065
Percent Solids:	45		

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
85-01-8	Phenanthrene	<0.15	0.15
108-95-2	Phenol	<1.5	1.5
129-00-0	Pyrene	<0.15	0.15
95-94-3	1,2,4,5-Tetrachlorobenzene	<0.29	0.29
58-90-2	2,3,4,6-Tetrachlorophenol	<0.29	0.29
120-82-1	1,2,4-Trichlorobenzene	<0.15	0.15
95-95-4	2,4,5-Trichlorophenol	<0.15	0.15
88-06-2	2,4,6-Trichlorophenol	<0.15	0.15
<b>Surrogates:</b>			
		<b>% Recovery</b>	<b>Control Limits</b>
	<i>2-Fluorophenol</i>	<i>60</i>	<i>33-113</i>
	<i>Phenol-d6</i>	<i>63</i>	<i>30-115</i>
	<i>Nitrobenzene-d5</i>	<i>60</i>	<i>33-131</i>
	<i>2-Fluorobiphenyl</i>	<i>62</i>	<i>46-122</i>
	<i>2,4,6-Tribromophenol</i>	<i>64</i>	<i>12-124</i>
	<i>o-Terphenyl</i>	<i>78</i>	<i>20-155</i>

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53187</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-05</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00

**Physical/Chemical Parameters by EPA/APHA/ASTM Methods**

Analyte	Analytical Result	RL	Unit	Dilution Factor	Method	Date Time Analyzed	By	QC Batch
Percent Solids	45	0.1	%	1	USEPA-3550C	03/19/13 11:00	BAR	1302409

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53188</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-06</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00
Unit:	mg/kg dry	Prepared:	03/19/13 By: JTS
Dilution Factor:	4	Analyzed:	03/21/13 By: JLB
QC Batch:	1302404	Analytical Batch:	3C21065
Percent Solids:	52		

**Semivolatile Organic Compounds by EPA Method 8270C**

CAS Number	Analyte	Analytical Result	RL
83-32-9	Acenaphthene	<0.13	0.13
208-96-8	Acenaphthylene	<0.13	0.13
98-86-2	Acetophenone	<0.13	0.13
120-12-7	Anthracene	<0.13	0.13
1912-24-9	Atrazine	<0.13	0.13
100-52-7	Benzaldehyde	<0.51	0.51
56-55-3	Benzo(a)anthracene	<0.13	0.13
50-32-8	Benzo(a)pyrene	<0.13	0.13
205-99-2	Benzo(b)fluoranthene	<0.13	0.13
207-08-9	Benzo(k)fluoranthene	<0.13	0.13
191-24-2	Benzo(g,h,i)perylene	<0.26	0.26
65-85-0	Benzoic Acid	<5.1	5.1
100-51-6	Benzyl Alcohol	<0.13	0.13
92-52-4	1,1'-Biphenyl	<0.13	0.13
101-55-3	4-Bromophenyl Phenyl Ether	<0.13	0.13
85-68-7	Butyl Benzyl Phthalate	<0.26	0.26
105-60-2	Caprolactam	<0.51	0.51
86-74-8	Carbazole	<1.3	1.3
59-50-7	4-Chloro-3-methylphenol	<0.13	0.13
106-47-8	4-Chloroaniline	<0.51	0.51
111-91-1	Bis(2-chloroethoxy)methane	<0.13	0.13
111-44-4	Bis(2-chloroethyl) Ether	<0.13	0.13
108-60-1	Bis(2-chloroisopropyl) Ether	<0.13	0.13
91-58-7	2-Chloronaphthalene	<0.13	0.13
95-57-8	2-Chlorophenol	<0.13	0.13
7005-72-3	4-Chlorophenyl Phenyl Ether	<0.13	0.13
218-01-9	Chrysene	<0.13	0.13
53-70-3	Dibenz(a,h)anthracene	<0.26	0.26
132-64-9	Dibenzofuran	<0.13	0.13
84-74-2	Di-n-butyl Phthalate	<0.51	0.51
91-94-1	3,3'-Dichlorobenzidine	<6.4	6.4

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**ANALYTICAL REPORT**

Client: **Bay West**  
 Project: Bench Scale Test  
 Client Sample ID: **53188**  
 Lab Sample ID: **1303154-06**  
 Matrix: Soil  
 Unit: mg/kg dry  
 Dilution Factor: 4  
 QC Batch: 1302404  
 Percent Solids: 52

Work Order: **1303154**  
 Description: FMC ESD 00439  
 Sampled: 03/08/13 12:00  
 Sampled By: Dave Poague  
 Received: 03/09/13 09:00  
 Prepared: 03/19/13 By: JTS  
 Analyzed: 03/21/13 By: JLB  
 Analytical Batch: 3C21065

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
120-83-2	2,4-Dichlorophenol	<0.26	0.26
84-66-2	Diethyl Phthalate	<0.13	0.13
105-67-9	2,4-Dimethylphenol	<1.3	1.3
131-11-3	Dimethyl Phthalate	<0.13	0.13
534-52-1	4,6-Dinitro-2-methylphenol	<1.3	1.3
51-28-5	2,4-Dinitrophenol	<1.3	1.3
606-20-2	2,6-Dinitrotoluene	<0.13	0.13
121-14-2	2,4-Dinitrotoluene	<0.26	0.26
117-84-0	Di-n-octyl Phthalate	<0.13	0.13
117-81-7	Bis(2-ethylhexyl) Phthalate	<b>0.84</b>	0.26
206-44-0	Fluoranthene	<0.13	0.13
86-73-7	Fluorene	<0.26	0.26
118-74-1	Hexachlorobenzene	<0.13	0.13
87-68-3	Hexachlorobutadiene	<0.13	0.13
77-47-4	Hexachlorocyclopentadiene	<0.13	0.13
67-72-1	Hexachloroethane	<0.13	0.13
193-39-5	Indeno(1,2,3-cd)pyrene	<0.26	0.26
78-59-1	Isophorone	<0.13	0.13
91-57-6	2-Methylnaphthalene	<0.13	0.13
106-44-5	4-Methylphenol	<0.13	0.13
95-48-7	2-Methylphenol	<0.13	0.13
91-20-3	Naphthalene	<b>0.25</b>	0.13
100-01-6	4-Nitroaniline	<0.26	0.26
88-74-4	2-Nitroaniline	<0.13	0.13
99-09-2	3-Nitroaniline	<0.26	0.26
98-95-3	Nitrobenzene	<0.13	0.13
88-75-5	2-Nitrophenol	<0.13	0.13
100-02-7	4-Nitrophenol	<5.1	5.1
86-30-6	N-Nitroso-diphenylamine	<0.13	0.13
621-64-7	N-Nitroso-di-n-propylamine	<0.13	0.13
87-86-5	Pentachlorophenol	<b>4.3</b>	1.3

Continued on next page

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53188</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-06</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00
Unit:	mg/kg dry	Prepared:	03/19/13 By: JTS
Dilution Factor:	4	Analyzed:	03/21/13 By: JLB
QC Batch:	1302404	Analytical Batch:	3C21065
Percent Solids:	52		

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

CAS Number	Analyte	Analytical Result	RL
85-01-8	Phenanthrene	<0.13	0.13
108-95-2	Phenol	<1.3	1.3
129-00-0	Pyrene	<0.13	0.13
95-94-3	1,2,4,5-Tetrachlorobenzene	<0.26	0.26
58-90-2	2,3,4,6-Tetrachlorophenol	<0.26	0.26
120-82-1	1,2,4-Trichlorobenzene	<0.13	0.13
95-95-4	2,4,5-Trichlorophenol	<0.13	0.13
88-06-2	2,4,6-Trichlorophenol	<0.13	0.13

<i>Surrogates:</i>	<i>% Recovery</i>	<i>Control Limits</i>
<i>2-Fluorophenol</i>	<i>71</i>	<i>33-113</i>
<i>Phenol-d6</i>	<i>76</i>	<i>30-115</i>
<i>Nitrobenzene-d5</i>	<i>70</i>	<i>33-131</i>
<i>2-Fluorobiphenyl</i>	<i>70</i>	<i>46-122</i>
<i>2,4,6-Tribromophenol</i>	<i>71</i>	<i>12-124</i>
<i>o-Terphenyl</i>	<i>80</i>	<i>20-155</i>

**ANALYTICAL REPORT**

Client:	<b>Bay West</b>	Work Order:	<b>1303154</b>
Project:	Bench Scale Test	Description:	FMC ESD 00439
Client Sample ID:	<b>53188</b>	Sampled:	03/08/13 12:00
Lab Sample ID:	<b>1303154-06</b>	Sampled By:	Dave Poague
Matrix:	Soil	Received:	03/09/13 09:00

**Physical/Chemical Parameters by EPA/APHA/ASTM Methods**

Analyte	Analytical Result	RL	Unit	Dilution Factor	Method	Date Time Analyzed	By	QC Batch
Percent Solids	52	0.1	%	1	USEPA-3550C	03/19/13 11:00	BAR	1302409

**QUALITY CONTROL REPORT**
**Semivolatile Organic Compounds by EPA Method 8270C**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1302297** 3510C Liquid-Liquid Extraction/USEPA-8270C

**Method Blank**

Analyzed: 03/20/2013 By: JLB

Unit: ug/L

Analytical Batch: 3C21059

Acenaphthene			<0.50					0.50
Acenaphthylene			<0.50					0.50
Acetophenone			<0.50			--		0.50
Anthracene			<0.50					0.50
Atrazine			<0.50					0.50
Benzaldehyde			<0.50			--		0.50
Benzo(a)anthracene			<0.50			--		0.50
Benzo(a)pyrene			<0.50			--		0.50
Benzo(b)fluoranthene			<0.50					0.50
Benzo(k)fluoranthene			<0.50					0.50
Benzo(g,h,i)perylene			<0.50					0.50
1,1'-Biphenyl			<0.50			--		0.50
4-Bromophenyl Phenyl Ether			<0.50					0.50
Butyl Benzyl Phthalate			<1.0			--		1.0
Caprolactam			<1.0					1.0
Carbazole			<0.50			--		0.50
4-Chloro-3-methylphenol			<0.50					0.50
4-Chloroaniline			<1.0					1.0
Bis(2-chloroethoxy)methane			<0.50					0.50
Bis(2-chloroethyl) Ether			<0.50					0.50
Bis(2-chloroisopropyl) Ether			<0.50					0.50
2-Chloronaphthalene			<0.50					0.50
2-Chlorophenol			<0.50					0.50
4-Chlorophenyl Phenyl Ether			<0.50					0.50
Chrysene			<0.50			--		0.50
Dibenz(a,h)anthracene			<0.50					0.50
Dibenzofuran			<0.50					0.50
Di-n-butyl Phthalate			<1.0			--		1.0
3,3'-Dichlorobenzidine			<10			--		10
2,4-Dichlorophenol			<0.50					0.50
Diethyl Phthalate			<0.50			--		0.50
2,4-Dimethylphenol			<1.0					1.0
Dimethyl Phthalate			<0.50					0.50
4,6-Dinitro-2-methylphenol			<5.0			--		5.0
2,4-Dinitrophenol			<5.0					5.0
2,4-Dinitrotoluene			<0.50					0.50

Continued on next page

**QUALITY CONTROL REPORT**
**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1302297 (Continued)** 3510C Liquid-Liquid Extraction/USEPA-8270C

**Method Blank (Continued)**

 Analyzed: 03/20/2013 By: JLB  
 Analytical Batch: 3C21059

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
<b>Method Blank (Continued)</b>								
Unit: ug/L								
2,6-Dinitrotoluene			<0.50					0.50
Di-n-octyl Phthalate			<0.50			--		0.50
Bis(2-ethylhexyl) Phthalate			<0.50			--		0.50
Fluoranthene			<0.50			--		0.50
Fluorene			<0.50					0.50
Hexachlorobenzene			<0.50			--		0.50
Hexachlorobutadiene			<0.50					0.50
Hexachlorocyclopentadiene			<0.50					0.50
Hexachloroethane			<0.50					0.50
Indeno(1,2,3-cd)pyrene			<0.50					0.50
Isophorone			<0.50			--		0.50
2-Methylnaphthalene			<0.50					0.50
2-Methylphenol			<0.50					0.50
4-Methylphenol			<0.50			--		0.50
Naphthalene			<0.50					0.50
2-Nitroaniline			<0.50					0.50
3-Nitroaniline			<1.0			--		1.0
4-Nitroaniline			<1.0					1.0
Nitrobenzene			<0.50			--		0.50
4-Nitrophenol			<5.0					5.0
2-Nitrophenol			<0.50					0.50
N-Nitroso-diphenylamine			<0.50					0.50
N-Nitroso-di-n-propylamine			<0.50			--		0.50
Pentachlorophenol			<0.50			--		0.50
Phenanthrene			<0.50			--		0.50
Phenol			<0.50			--		0.50
Pyrene			<0.50			--		0.50
1,2,4,5-Tetrachlorobenzene			<2.0					2.0
2,3,4,6-Tetrachlorophenol			<5.0					5.0
2,4,6-Trichlorophenol			<0.50					0.50
2,4,5-Trichlorophenol			<0.50					0.50

Continued on next page

**QUALITY CONTROL REPORT**

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1302297 (Continued)** 3510C Liquid-Liquid Extraction/USEPA-8270C

**Method Blank (Continued)**

Analyzed: 03/20/2013 By: JLB  
 Analytical Batch: 3C21059

Unit: ug/L

**Surrogates:**

<i>2-Fluorophenol</i>	54	20-70
<i>Phenol-d6</i>	36	18-45
<i>Nitrobenzene-d5</i>	90	31-123
<i>2-Fluorobiphenyl</i>	94	25-113
<i>2,4,6-Tribromophenol</i>	73	30-121
<i>o-Terphenyl</i>	101	42-125

**Laboratory Control Sample**

Analyzed: 03/20/2013 By: JLB  
 Analytical Batch: 3C21059

Unit: ug/L

Acenaphthene	10.0	<b>10.0</b>	100	53-126	--	20	0.50
4-Chloro-3-methylphenol	10.0	<b>9.18</b>	92	53-120	--	20	0.50
2-Chlorophenol	10.0	<b>8.58</b>	86	44-121	--	20	0.50
1,4-Dichlorobenzene	10.0	<b>8.39</b>	84	41-124	--	20	0.50
2,4-Dinitrotoluene	10.0	<b>10.9</b>	109	55-131	--	20	0.50
Naphthalene	10.0	<b>9.13</b>	91	50-127	--	20	0.50
4-Nitrophenol	10.0	<b>2.48</b>	25	17-70	--	20	5.0
N-Nitroso-di-n-propylamine	10.0	<b>8.09</b>	81	49-125	--	20	0.50
Pentachlorophenol	10.0	<b>7.95</b>	80	21-124	--	20	0.50
Phenol	10.0	<b>4.25</b>	42	22-60	--	20	0.50
Pyrene	10.0	<b>10.9</b>	109	60-134	--	20	0.50
1,2,4-Trichlorobenzene	10.0	<b>8.54</b>	85	47-123	--	20	0.50

**Surrogates:**

<i>2-Fluorophenol</i>	51	20-70
<i>Phenol-d6</i>	32	18-45
<i>Nitrobenzene-d5</i>	89	31-123
<i>2-Fluorobiphenyl</i>	92	25-113
<i>2,4,6-Tribromophenol</i>	95	30-121
<i>o-Terphenyl</i>	96	42-125

Continued on next page

**QUALITY CONTROL REPORT**
**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1302404** 3550C Sonication Extraction/USEPA-8270C

**Method Blank**

Analyzed: 03/20/2013 By: JLB

Unit: mg/kg wet

Analytical Batch: 3C21061

Acenaphthene			<0.017					0.017
Acenaphthylene			<0.017					0.017
Acetophenone			<0.017					0.017
Anthracene			<0.017					0.017
Atrazine			<0.017					0.017
Benzaldehyde			<0.067			--		0.067
Benzo(a)anthracene			<0.017			--		0.017
Benzo(a)pyrene			<0.017			--		0.017
Benzo(b)fluoranthene			<0.017					0.017
Benzo(k)fluoranthene			<0.017					0.017
Benzo(g,h,i)perylene			<0.033					0.033
Benzoic Acid			<0.67			--		0.67
Benzyl Alcohol			<0.017					0.017
1,1'-Biphenyl			<0.017			--		0.017
4-Bromophenyl Phenyl Ether			<0.017					0.017
Butyl Benzyl Phthalate			<0.033			--		0.033
Caprolactam			<0.067					0.067
Carbazole			<0.17					0.17
4-Chloro-3-methylphenol			<0.017					0.017
4-Chloroaniline			<0.067					0.067
Bis(2-chloroethoxy)methane			<0.017					0.017
Bis(2-chloroethyl) Ether			<0.017					0.017
Bis(2-chloroisopropyl) Ether			<0.017					0.017
2-Chloronaphthalene			<0.017					0.017
2-Chlorophenol			<0.017					0.017
4-Chlorophenyl Phenyl Ether			<0.017					0.017
Chrysene			<0.017			--		0.017
Dibenz(a,h)anthracene			<0.033					0.033
Dibenzofuran			<0.017					0.017
Di-n-butyl Phthalate			<0.067			--		0.067
3,3'-Dichlorobenzidine			<0.83					0.83
2,4-Dichlorophenol			<0.033					0.033
Diethyl Phthalate			<0.017			--		0.017
2,4-Dimethylphenol			<0.17					0.17
Dimethyl Phthalate			<0.017			--		0.017
4,6-Dinitro-2-methylphenol			<0.17			--		0.17

Continued on next page

**QUALITY CONTROL REPORT**
**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1302404 (Continued)** 3550C Sonication Extraction/USEPA-8270C

**Method Blank (Continued)**

Analyzed: 03/20/2013 By: JLB

Unit: mg/kg wet

Analytical Batch: 3C21061

2,4-Dinitrophenol			<0.17					0.17
2,6-Dinitrotoluene			<0.017					0.017
2,4-Dinitrotoluene			<0.033					0.033
Di-n-octyl Phthalate			<0.017			--		0.017
Bis(2-ethylhexyl) Phthalate			<0.033			--		0.033
Fluoranthene			<0.017			--		0.017
Fluorene			<0.033					0.033
Hexachlorobenzene			<0.017					0.017
Hexachlorobutadiene			<0.017					0.017
Hexachlorocyclopentadiene			<0.017					0.017
Hexachloroethane			<0.017			--		0.017
Indeno(1,2,3-cd)pyrene			<0.033					0.033
Isophorone			<0.017					0.017
2-Methylnaphthalene			<0.017					0.017
4-Methylphenol			<0.017			--		0.017
2-Methylphenol			<0.017					0.017
Naphthalene			<0.017					0.017
4-Nitroaniline			<0.033					0.033
2-Nitroaniline			<0.017					0.017
3-Nitroaniline			<0.033					0.033
Nitrobenzene			<0.017			--		0.017
2-Nitrophenol			<0.017					0.017
4-Nitrophenol			<0.67					0.67
N-Nitroso-diphenylamine			<0.017			--		0.017
N-Nitroso-di-n-propylamine			<0.017			--		0.017
Pentachlorophenol			<0.17					0.17
Phenanthrene			<0.017			--		0.017
Phenol			<0.17			--		0.17
Pyrene			<0.017					0.017
1,2,4,5-Tetrachlorobenzene			<0.033					0.033
2,3,4,6-Tetrachlorophenol			<0.033					0.033
1,2,4-Trichlorobenzene			<0.017					0.017
2,4,5-Trichlorophenol			<0.017					0.017
2,4,6-Trichlorophenol			<0.017					0.017

Continued on next page



**QUALITY CONTROL REPORT**

**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1302404 (Continued)** 3550C Sonication Extraction/USEPA-8270C

**Method Blank (Continued)**

Analyzed: 03/20/2013 By: JLB  
 Analytical Batch: 3C21061

Unit: mg/kg wet

**Surrogates:**

<i>2-Fluorophenol</i>		70		33-113
<i>Phenol-d6</i>		71		30-115
<i>Nitrobenzene-d5</i>		69		33-131
<i>2-Fluorobiphenyl</i>		79		46-122
<i>2,4,6-Tribromophenol</i>		61		12-124
<i>o-Terphenyl</i>		83		20-155

**Laboratory Control Sample**

Analyzed: 03/21/2013 By: JLB  
 Analytical Batch: 3C22004

Unit: mg/kg wet

Acenaphthene	0.331	<b>0.274</b>	83	55-113	--	20	0.017
4-Chloro-3-methylphenol	0.331	<b>0.186</b>	56	34-113	--	20	0.017
2-Chlorophenol	0.331	<b>0.229</b>	69	62-118	--	20	0.017
1,4-Dichlorobenzene	0.331	<b>0.240</b>	72	61-111	--	20	0.017
2,4-Dinitrotoluene	0.331	<b>0.273</b>	82	51-128	--	20	0.033
Naphthalene	0.331	<b>0.262</b>	79	52-128	--	20	0.017
4-Nitrophenol	0.331	<b>0.225</b>	68	36-131	--	20	0.67
N-Nitroso-di-n-propylamine	0.331	<b>0.236</b>	71	48-127	--	20	0.017
Pentachlorophenol	0.331	<b>0.170</b>	51	19-117	--	20	0.17
Phenol	0.331	<b>0.189</b>	57	53-120	--	20	0.17
Pyrene	0.331	<b>0.311</b>	94	60-132	--	20	0.017
1,2,4-Trichlorobenzene	0.331	<b>0.247</b>	75	51-110	--	20	0.017

**Surrogates:**

<i>2-Fluorophenol</i>		72		33-113
<i>Phenol-d6</i>		58		30-115
<i>Nitrobenzene-d5</i>		77		33-131

Continued on next page

**QUALITY CONTROL REPORT**
**Semivolatile Organic Compounds by EPA Method 8270C (Continued)**

Analyte	Sample Conc.	Spike Qty.	Result	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
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**QC Batch: 1302404 (Continued)** 3550C Sonication Extraction/USEPA-8270C

**Laboratory Control Sample (Continued)**

Unit: mg/kg wet

 Analyzed: 03/21/2013 By: JLB  
 Analytical Batch: 3C22004

**Surrogates (Continued):**

<i>2-Fluorobiphenyl</i>				82	46-122
<i>2,4,6-Tribromophenol</i>				77	12-124
<i>o-Terphenyl</i>				88	20-155

**Laboratory Control Sample Duplicate**

Unit: mg/kg wet

 Analyzed: 03/20/2013 By: JLB  
 Analytical Batch: 3C21061

Acenaphthene	0.325	<b>0.270</b>		83	55-113	2	20	0.017
4-Chloro-3-methylphenol	0.325	<b>0.184</b>		57	34-113	0.7	20	0.017
2-Chlorophenol	0.325	<b>0.228</b>		70	62-118	0.4	20	0.017
1,4-Dichlorobenzene	0.325	<b>0.240</b>		74	61-111	0.08	20	0.017
2,4-Dinitrotoluene	0.325	<b>0.274</b>		84	51-128	0.6	20	0.033
Naphthalene	0.325	<b>0.264</b>		81	52-128	0.8	20	0.017
4-Nitrophenol	0.325	<b>0.237</b>		73	36-131	5	20	0.67
N-Nitroso-di-n-propylamine	0.325	<b>0.228</b>		70	48-127	4	20	0.017
Pentachlorophenol	0.325	<b>0.182</b>		56	19-117	7	20	0.17
Phenol	0.325	<b>0.185</b>		57	53-120	2	20	0.17
Pyrene	0.325	<b>0.294</b>		91	60-132	6	20	0.017
1,2,4-Trichlorobenzene	0.325	<b>0.243</b>		75	51-110	2	20	0.017

**Surrogates:**

<i>2-Fluorophenol</i>				73	33-113
<i>Phenol-d6</i>				57	30-115
<i>Nitrobenzene-d5</i>				79	33-131
<i>2-Fluorobiphenyl</i>				84	46-122
<i>2,4,6-Tribromophenol</i>				74	12-124
<i>o-Terphenyl</i>				86	20-155

**QUALITY CONTROL REPORT**
**Physical/Chemical Parameters by EPA/APHA/ASTM Methods**

QC Type	Sample Conc.	Spike Qty.	Result	Unit	Spike % Rec.	Control Limits	RPD	RPD Limits	RL
<b>Analyte: Percent Solids/USEPA-3550C</b>									
QC Batch: 1302409 (Method Specific Preparation)						Analyzed: 03/19/2013		By: BAR	
Method Blank			<0.1	%					0.1



**STATEMENT OF DATA QUALIFICATIONS**

All analyses have been validated and comply with our Quality Control Program.  
No Qualification is required.



5560 Corporate Exchange Court SE  
 Grand Rapids, MI 49512  
 Phone (616) 975-4500 Fax (616) 942-7463  
 www.trimatrixlabs.com

### Chain of Custody Record

COC No. \_\_\_\_\_

#### Analyses Requested

Pg. \_\_\_ of \_\_\_

For Lab Use Only  
 GSI  
 VOA # 4712  
 Receipt #  
 Project #  
 Work Order No.  
 Scheduled  
 Method  
 Sample Number

Client Name: **FMC ESD**  
 Project Name: \_\_\_\_\_  
 Address: \_\_\_\_\_  
 Client Project No. / P.O. No.: \_\_\_\_\_  
 City, State Zip: \_\_\_\_\_  
 Invoice To:  Client  Other (comments)  
 Phone/Fax: \_\_\_\_\_  
 Contact/Report to: \_\_\_\_\_  
 Email: \_\_\_\_\_

1000ml Amber  
 125ml  
 125ml Amber bottle  
 VOC Phenols  
 (Method 8270)


- ← PRESERVATIVES
- A NONE pH=7
  - B HNO<sub>3</sub> pH<2
  - C H<sub>2</sub>SO<sub>4</sub> pH<2
  - D 1+1 HCl pH<2
  - E NaOH pH>12
  - F ZnAc/NaOH pH>9
  - G MeOH
  - H Other (note below)

Field Sample ID #	Container ID	Sample Date	Sample Time	C O M P	R A B	Matrix	Number of Containers Submitted	Label	Sample Comments
53183	Water	CONTROL					1		
53184	Water	10%					1		
53185	Water	20%					1		
53186	Carbon	CONTROL					1		
53187	Carbon	10%					1		
53188	Carbon	20%					1		

Sampled By (print): \_\_\_\_\_ Comments: \_\_\_\_\_  
 How Shipped? \_\_\_\_\_  
 Tracker's Signature \_\_\_\_\_ Tracking No. \_\_\_\_\_  
 Carrier: \_\_\_\_\_  
 Company: \_\_\_\_\_  
 1. Released By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 2. Received By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 3. Received For Lab By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_



### SAMPLE RECEIVING / LOG-IN CHECKLIST

 <b>TRIMATRIX</b> LABORATORIES		Client: <u>FMC Environmental Solutions</u> Receipt Record Page Line #: <u>47-2</u>		Work Order #: <u>1303154</u> New / Add To: _____ Project Chemist: _____ Sample #: _____																																																																			
Recorded by (initials/date): <u>SLR 3/9/13</u>		<input checked="" type="checkbox"/> Cooler Qty Received: <u>1</u> <input type="checkbox"/> Box <input type="checkbox"/> Other: _____		<input checked="" type="checkbox"/> IR Gun (#202) <input type="checkbox"/> Digital Thermometer (#54) <input type="checkbox"/> See Additional Cooler Information Form <input type="checkbox"/> Other (# _____)																																																																			
Cooler # <u>-</u> Time <u>0935</u> Custody Seals: <input checked="" type="checkbox"/> None <input type="checkbox"/> Present / Intact <input type="checkbox"/> Present / Not Intact Coolant Location: <u>Sides</u> Dispersed / Top / Middle / Bottom Coolant/Temperature Taken Via: <input type="checkbox"/> Loose Ice / Avg 2-3 containers <input checked="" type="checkbox"/> Bagged Ice / Avg 2-3 containers <input checked="" type="checkbox"/> Blue Ice / Avg 2-3 containers <input type="checkbox"/> None / Avg 2-3 containers Alternate Temperature Taken Via: <input type="checkbox"/> Temperature Blank (TB) <input type="checkbox"/> 1 Container Recorded °C: _____ Correction Factor °C: _____ Actual °C: _____ Temp Blank: <u>6.1</u> TB location: Representative / Not Representative <table border="1" style="width: 100%; border-collapse: collapse;"> <tr><td>1</td><td><u>5.7</u></td><td><u>-</u></td><td><u>5.7</u></td></tr> <tr><td>2</td><td><u>7.6</u></td><td><u>-</u></td><td><u>7.6</u></td></tr> <tr><td>3</td><td><u>7.4</u></td><td><u>-</u></td><td><u>7.4</u></td></tr> <tr><td colspan="4" style="text-align: center;">Average °C</td></tr> </table> <input type="checkbox"/> Cooler ID on COC? <u>69</u> <input type="checkbox"/> VOC Trip Blank received?		1	<u>5.7</u>	<u>-</u>	<u>5.7</u>	2	<u>7.6</u>	<u>-</u>	<u>7.6</u>	3	<u>7.4</u>	<u>-</u>	<u>7.4</u>	Average °C				Cooler # _____ Time _____ Custody Seals: <input type="checkbox"/> None <input type="checkbox"/> Present / Intact <input type="checkbox"/> Present / Not Intact Coolant Location: Dispersed / Top / Middle / Bottom Coolant/Temperature Taken Via: <input type="checkbox"/> Loose Ice / Avg 2-3 containers <input type="checkbox"/> Bagged Ice / Avg 2-3 containers <input type="checkbox"/> Blue Ice / Avg 2-3 containers <input type="checkbox"/> None / Avg 2-3 containers Alternate Temperature Taken Via: <input type="checkbox"/> Temperature Blank (TB) <input type="checkbox"/> 1 Container Recorded °C: _____ Correction Factor °C: _____ Actual °C: _____ Temp Blank: _____ TB location: Representative / Not Representative <table border="1" style="width: 100%; border-collapse: collapse;"> <tr><td>1</td><td></td><td></td><td></td></tr> <tr><td>2</td><td></td><td></td><td></td></tr> <tr><td>3</td><td></td><td></td><td></td></tr> <tr><td colspan="4" style="text-align: center;">Average °C</td></tr> </table> <input type="checkbox"/> Cooler ID on COC? <input type="checkbox"/> VOC Trip Blank received?		1				2				3				Average °C				Cooler # _____ Time _____ Custody Seals: <input type="checkbox"/> None <input type="checkbox"/> Present / Intact <input type="checkbox"/> Present / Not Intact Coolant Location: Dispersed / Top / Middle / Bottom Coolant/Temperature Taken Via: <input type="checkbox"/> Loose Ice / Avg 2-3 containers <input type="checkbox"/> Bagged Ice / Avg 2-3 containers <input type="checkbox"/> Blue Ice / Avg 2-3 containers <input type="checkbox"/> None / Avg 2-3 containers Alternate Temperature Taken Via: <input type="checkbox"/> Temperature Blank (TB) <input type="checkbox"/> 1 Container Recorded °C: _____ Correction Factor °C: _____ Actual °C: _____ Temp Blank: _____ TB location: Representative / Not Representative <table border="1" style="width: 100%; border-collapse: collapse;"> <tr><td>1</td><td></td><td></td><td></td></tr> <tr><td>2</td><td></td><td></td><td></td></tr> <tr><td>3</td><td></td><td></td><td></td></tr> <tr><td colspan="4" style="text-align: center;">Average °C</td></tr> </table> <input type="checkbox"/> Cooler ID on COC? <input type="checkbox"/> VOC Trip Blank received?		1				2				3				Average °C				Cooler # _____ Time _____ Custody Seals: <input type="checkbox"/> None <input type="checkbox"/> Present / Intact <input type="checkbox"/> Present / Not Intact Coolant Location: Dispersed / Top / Middle / Bottom Coolant/Temperature Taken Via: <input type="checkbox"/> Loose Ice / Avg 2-3 containers <input type="checkbox"/> Bagged Ice / Avg 2-3 containers <input type="checkbox"/> Blue Ice / Avg 2-3 containers <input type="checkbox"/> None / Avg 2-3 containers Alternate Temperature Taken Via: <input type="checkbox"/> Temperature Blank (TB) <input type="checkbox"/> 1 Container Recorded °C: _____ Correction Factor °C: _____ Actual °C: _____ Temp Blank: _____ TB location: Representative / Not Representative <table border="1" style="width: 100%; border-collapse: collapse;"> <tr><td>1</td><td></td><td></td><td></td></tr> <tr><td>2</td><td></td><td></td><td></td></tr> <tr><td>3</td><td></td><td></td><td></td></tr> <tr><td colspan="4" style="text-align: center;">Average °C</td></tr> </table> <input type="checkbox"/> Cooler ID on COC? <input type="checkbox"/> VOC Trip Blank received?		1				2				3				Average °C			
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<b>If any shaded areas checked, complete Sample Receiving Non-Conformance and/or Inventory Form</b>																																																																							
<b>Paperwork Received</b> Yes No <input checked="" type="checkbox"/> Chain of Custody record(s)? If No, Initiated By: _____ <input checked="" type="checkbox"/> Received for Lab Signed/Date/Time? <input type="checkbox"/> Shipping document? <input checked="" type="checkbox"/> Other: _____			<b>Check Sample Preservation</b> N/A Yes No <input checked="" type="checkbox"/> Average sample temperature ≤ 8° C? <input checked="" type="checkbox"/> Was thermal preservation required? If "No", Project Chemist Approval Initials: _____ <input checked="" type="checkbox"/> If "Yes" Completed Non Con Cooler - Cont Inventory Form? <input checked="" type="checkbox"/> Completed Sample Preservation Verification Form? <input checked="" type="checkbox"/> Samples chemically preserved correctly? If "No", added orange tag? <input checked="" type="checkbox"/> Received pre-preserved VOC soils? <input type="checkbox"/> MeOH <input type="checkbox"/> Na <sub>2</sub> SO <sub>4</sub>																																																																				
<b>COC Information</b> <input type="checkbox"/> TriMatrix COC <input checked="" type="checkbox"/> Other: _____ COC ID Numbers: _____			<b>Check for Short Hold-Time Prep/Analyses</b> <input type="checkbox"/> Bacteriological <input type="checkbox"/> Air Bags <input type="checkbox"/> EnCores / Methanol Pre-Preserved <input type="checkbox"/> Formaldehyde/Aldelyde <input type="checkbox"/> Green-tagged containers <input type="checkbox"/> Yellow/White-tagged 1L ampers (SV Prep-Lab)																																																																				
<b>Check COC for Accuracy</b> Yes No <input type="checkbox"/> Analysis Requested? <input checked="" type="checkbox"/> Sample ID matches COC? <input checked="" type="checkbox"/> Sample Date and Time matches COC? <input checked="" type="checkbox"/> Container type completed on COC? <input type="checkbox"/> All container types indicated are received?			<b>Check COC for Accuracy</b> Yes No <input type="checkbox"/> Analysis Requested? <input checked="" type="checkbox"/> Sample ID matches COC? <input checked="" type="checkbox"/> Sample Date and Time matches COC? <input checked="" type="checkbox"/> Container type completed on COC? <input type="checkbox"/> All container types indicated are received?																																																																				
<b>Sample Condition Summary</b> N/A Yes No <input checked="" type="checkbox"/> Broken containers/lids? <input checked="" type="checkbox"/> Missing or incomplete labels? <input checked="" type="checkbox"/> Illegible information on labels? <input checked="" type="checkbox"/> Low volume received? <input type="checkbox"/> Inappropriate or non-TriMatrix containers received? <input type="checkbox"/> VOC vials / TOX containers have headspace? <input type="checkbox"/> Extra sample locations / containers not listed on COC?			<b>Notes</b> <input type="checkbox"/> Trip Blank received <input type="checkbox"/> Trip Blank not listed on COC <table border="1" style="width: 100%; border-collapse: collapse;"> <tr> <td>Cooler Received (Date/Time): <u>3/9/13 0900</u></td> <td>Paperwork Delivered (Date/Time): <u>3/9/13 0955</u></td> <td>≤ 1 Hour Goal Met? <u>(Yes)</u> No</td> </tr> </table>			Cooler Received (Date/Time): <u>3/9/13 0900</u>	Paperwork Delivered (Date/Time): <u>3/9/13 0955</u>	≤ 1 Hour Goal Met? <u>(Yes)</u> No																																																															
Cooler Received (Date/Time): <u>3/9/13 0900</u>	Paperwork Delivered (Date/Time): <u>3/9/13 0955</u>	≤ 1 Hour Goal Met? <u>(Yes)</u> No																																																																					

47-2

Page 1 of 1 E-1303154

CHAIN OF CUSTODY RECORD 11352 Part 3

Project No.: FMC ESD 00439		Laboratory: TRIMATRIX LABORATORIES		Laboratory Contact: PHIL KOMAR		Adventus Remediation Technologies 1345 Fewster Drive Mississauga, Ontario Canada L4W 2A5 Tel: (905) 273-5374 Fax: (905) 273-4367		Analytical Laboratory to Complete Submission Number												
P.O. #		Adventus Remediation Tech. Contact (Name/Tel.): EVA JANZEN EXT. 232																		
Date	Time	Type		Matrix		No. of Containers	Adventus Remediation Tech. Sample Number				VOCs	SVOCs	Pesticides	Total Organic Carbon	Metals	TPH	CP	SVOC PHENOLS (METALS & DTC)	Remarks	Lab Sample #
		Composite	Grab	Soil	Water		ACTIVATED CARBON													
MAR 8/13		X		X		2	5	3	1	8	3							X	01	
		X		X		1	5	3	1	8	4							X	↓	
		X		X		1	5	3	1	8	5							X	↓	
		X		X		1	5	3	1	8	6							X	02	
		X		X		1	5	3	1	8	7							X	↓	
		X		X		1	5	3	1	8	8							X	↓	
Sampled by: (print name & initial) S. OWEN S.O + E. JANZEN 2P		Date Mar 8/13		Received by: (signature) <i>[Signature]</i>		Date Mar 8/13		Shipped by: <i>[Signature]</i>		Shipping Bill		NOTES Send analytical results to appropriate ART contact person.								
Relinquished by: (Signature) <i>[Signature]</i>		Date Mar 8/13		Received by Laboratory: <i>[Signature]</i>		Date/Time 3/9/13 0900														

Copies: White & Yellow-Laboratory, Pink-Sampler

