Summary Report for Chemical Pollution Control (CPC) Site Soil and Groundwater Investigation (1-52-015) Bay Shore, New York

Prepared for

New York State Department of Environmental Conservation
625 Broadway
Albany, New York 12233

Prepared by

EA Engineering, P.C. and Its Affiliate
EA Science and Technology
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(315) 431-4610

March 2008
Revision: FINAL
EA Project No.: 14368.12

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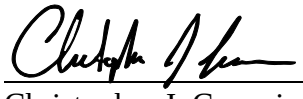
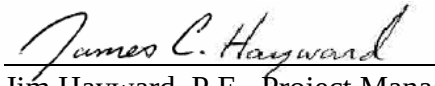


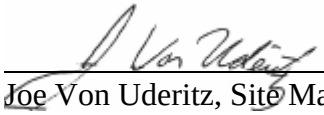
24 March 2008

Christopher J. Canonica, P.E., Program Manager — Date
EA Engineering, P.C.



24 March 2008

Jim Hayward, P.E., Project Manager — Date
EA Engineering, P.C.



24 March 2008

Joe Von Uderitz, Site Manager — Date
EA Science and Technology

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1. INTRODUCTION

1.1 PROJECT BACKGROUND

The New York State Department of Environmental Conservation (NYSDEC) tasked EA Engineering, P.C., and its affiliate EA Science and Technology (EA) to perform soil and groundwater sampling at the Chemical Pollution Control site (Site No. 152015). The site is located at 120 South 4th Street, in the City of Bay Shore, Suffolk County, New York (Figure 1).

This work assignment is being conducted under the NYSDEC State Superfund Standby Contract (Work Assignment No. D004438-12). The field activities for the sampling event were performed on 14 and 16 August 2007.

1.2 OBJECTIVES

The purpose of soil and groundwater sampling at Chemical Pollution Control (CPC) was to determine the appropriate classification listing of this site in the DEC's list of inactive hazardous waste disposal sites. The sampling program consisted of collecting sediment samples from four dry wells, soil samples from six borings and groundwater samples from nine monitoring wells throughout the target area (Figure 2).

This report was completed to discuss the field activities and summarize the soil and groundwater sample results.

1.3 REPORT ORGANIZATION

A summary of field activities completed in August 2007 is provided in Section 2. Analytical results are summarized in table format and presented in Section 3.

The following are provided as appendixes:

- Appendix A: Daily Field Reports
- Appendix B: Groundwater Sampling Forms
- Appendix C: Soil Boring Logs
- Appendix D: Data Usability Summary Report.

2. FIELD ACTIVITIES

This section presents the overall approach of the field investigation activities that were performed to meet the stated objectives of the groundwater and soil/sediment sampling. EA's approach for implementing the work assignment included field sampling activities designed to evaluate the presence or absence of contaminants of concern (COC) at the site and to summarize the concentrations of potential COC through laboratory analysis.

The field investigation program was performed during August 2007 and included the following activities:

- ***Groundwater Sampling—***Collection and analysis of groundwater samples from nine groundwater monitoring wells
- ***Soil Boring Sampling—***Installation, collection, and analysis of soil samples from six soil borings
- ***Dry Well Sampling—***Collection and analysis of sediment from four dry wells.

Copies of the daily field reports are provided in Appendix A. Site sampling locations are detailed in Figure 2.

2.1 GROUNDWATER SAMPLING

Nine groundwater samples were collected from the monitoring wells using a peristaltic pump and dedicated section of polyethylene tubing. One well (MW-02) was found to have been destroyed previously and therefore was unable to be sampled. Prior to sampling, water level measurements were taken from each monitoring well to prepare a groundwater contour map and evaluate groundwater flow direction (Figure 3). Depth to groundwater measurements ranged from a low of 6.75 ft below top of casing (TOC) in MW-01 to 8.89 feet below TOC in MW-04. The monitoring wells were purged until water quality parameters (pH, conductivity, oxygen reduction potential, temperature, dissolved oxygen, and turbidity) were stabilized. Once groundwater parameters were stabilized, samples were collected, placed in a cooler with ice, and delivered to Hampton-Clark Veritech located in West Fairfield, New Jersey.

Groundwater samples were analyzed for volatile organic compounds (VOCs) by U.S. Environmental Protection Agency (USEPA) Method 8260, semivolatile organic compounds (SVOCs) USEPA Method 8270, polychlorinated biphenyls (PCBs) Method 8082, pesticides Method 8081, and total metals (unfiltered) Method 6010. Groundwater sampling forms are provided in Appendix B.

2.2 SOIL SAMPLING

2.2.1 Soil Boring Sampling

A total of six soil borings were installed by Land Air Water Environmental Services, Inc. (Center Moriches, New York) using a geoprobe and direct-push technologies. Soil borings were installed to assess preferential seepage of contaminants due to historical and/or accidental spills. Soil samples were recovered using a 5-ft macro-core and a dedicated acetate sleeve. Soil samples were screened at 1 ft intervals using a photo-ionization detector (PID). If the PID indicated the presence of VOCs, the soil boring was extended to a depth where VOCs were no longer present, or groundwater was observed (which ever occurred first). Each soil sample was described and logged identifying its geologic characteristics, features, PID readings, and properties. The Unified Soil Classification System was used to characterize the soil samples. Soil boring logs are provided in Appendix C.

Soil samples were collected at 18 in. below ground surface (bgs) if no VOCs were detected. If VOCs were detected, soil samples were collected from a 2 ft interval starting at 18 in. bgs until the groundwater table interface.

A total of 11 soil samples were collected and sent for laboratory analysis. Soil samples were analyzed for VOCs by USEPA Method 8260, SVOCs Method 8270, PCBs Method 8082, pesticides Method 8081, and heavy metals Method 6010.

Upon completion of the sampling, each soil boring was backfilled with bentonite and the surface was restored to its original condition (asphalt or concrete).

2.2.2 Dry Well Sampling

Four dry wells were sampled using a combination of geoprobe and direct-push technologies. The dry well grates were removed, and the depth to water and depth to bottom were gauged prior to sampling. Borings were installed in each dry well to analyze any preferential seepage of contaminants in the event of any accidental spill. Dry well sediment samples were recovered using a 5-ft macro-core and a dedicated acetate sleeve. Dry well borings were monitored at 1 ft intervals using a PID. If the PID indicated the presence of VOCs, the boring was extended to a depth where VOCs were no longer observed. Each sediment sample was described and logged identifying its geologic characteristics, features, PID readings, and properties. The Unified Soil Classification System was used to characterize the sediment samples. Boring logs are provided in Appendix C.

Dry wells DW-01 through DW-03 were sampled at approximately 18 in. below the bottom of the dry well. In dry well DW-04, two samples were collected due to PID screening levels observed during the boring installation.

A total of five dry well sediment samples were collected and sent for laboratory analysis. Dry well sediment samples were analyzed for VOCs by USEPA Method 8260, SVOCs Method 8270, PCBs Method 8082, pesticides Method 8081, and heavy metals Method 6010.

3. FIELD SAMPLING RESULTS

The following section presents the analytical results of the groundwater and soil/sediment samples collected during the field investigation. Analytical data for the groundwater, soil boring and dry well samples are provided in Table 1, Table 2, and Table 3 respectively. All groundwater, soil, and sediment samples were analyzed by Hampton-Clarke Veritech (West Fairfield, New Jersey). Hampton-Clarke Veritech is a New York State Department of Health Environmental Laboratories Approval Program (ELAP)-certified laboratory for VOCs, SVOCs, PCBs, pesticides, and heavy metals analysis and in accordance with the NYSDEC Analytical Services Protocol (ASP).

Copies of the DUSR for the laboratory analytical data are provided in Appendix D. Laboratory analytical data and Form I's will be provided with the final report.

3.1 GROUNDWATER SAMPLING RESULTS

Three VOC analytes were detected above the Ambient Water Quality Standards (AWQS) class GA standards, including tetrachloroethene (PCE), trichloroethene (TCE), and cis-1,2-Dichlorethene. The AWQS for the analytes detected is 5 μg/L. The maximum concentration of TCE detected was in MW-04 at a concentration of 330 μg/L. The only detection of PCE was in MW-04 at a concentration of 14 ug/L; however, this is an estimated value. The maximum concentration of cis-1,2-Dichlorethene was 320 μg/L, which was also detected in MW-04. Toluene was detected in MW-06 at a concentration of 1.5 μg/L, which is below the AWQS standard of 5 μg/L.

The only SVOC detected was bis(2-Ethylhexyl)phthalate in MW-08 and MW-10, with both values below the standard of 5 μg/L. Several wells showed concentrations of iron and sodium above the corresponding AQWS standards of 300 and 20,000 μg/L respectively. The maximum detection of iron was 1,100 μg/L (MW-06) and the maximum detection of sodium was 27,000 μg/L (MW-09). A concentration of 330 μg/L of manganese was detected in MW-10, over the AQWS standard of 300 μg/L. No pesticides or PCBs were detected in any of the groundwater samples.

Validated analytical results for groundwater samples are included with the DUSR provided in Appendix D.

3.2 SOIL BORING SAMPLING RESULTS

Several VOC and SVOC analytes were detected above the laboratory detection limit, but below the respective standards.

Chromium was detected in five soil samples above the Part 375 standard of 1 mg/Kg, with the maximum concentration observed in SB-03 1.5-3.5' (180 mg/Kg). Silver was detected in one sample (SB-02 5-7') at a concentration of 3.4 mg/Kg, which was above the corresponding Part 375 standard of 2 mg/Kg.

One pesticide (chlordane) was detected above the reporting limit at soil boring location (SB-02 5-7') at a concentration of 0.021 mg/Kg. However, the concentration was below the corresponding standard of 0.094 mg/Kg.

No PCBs were detected in any of the soil boring samples.

Validated analytical results for soil boring samples are included with the DUSR provided in Appendix D.

3.3 DRY WELL SAMPLING RESULTS

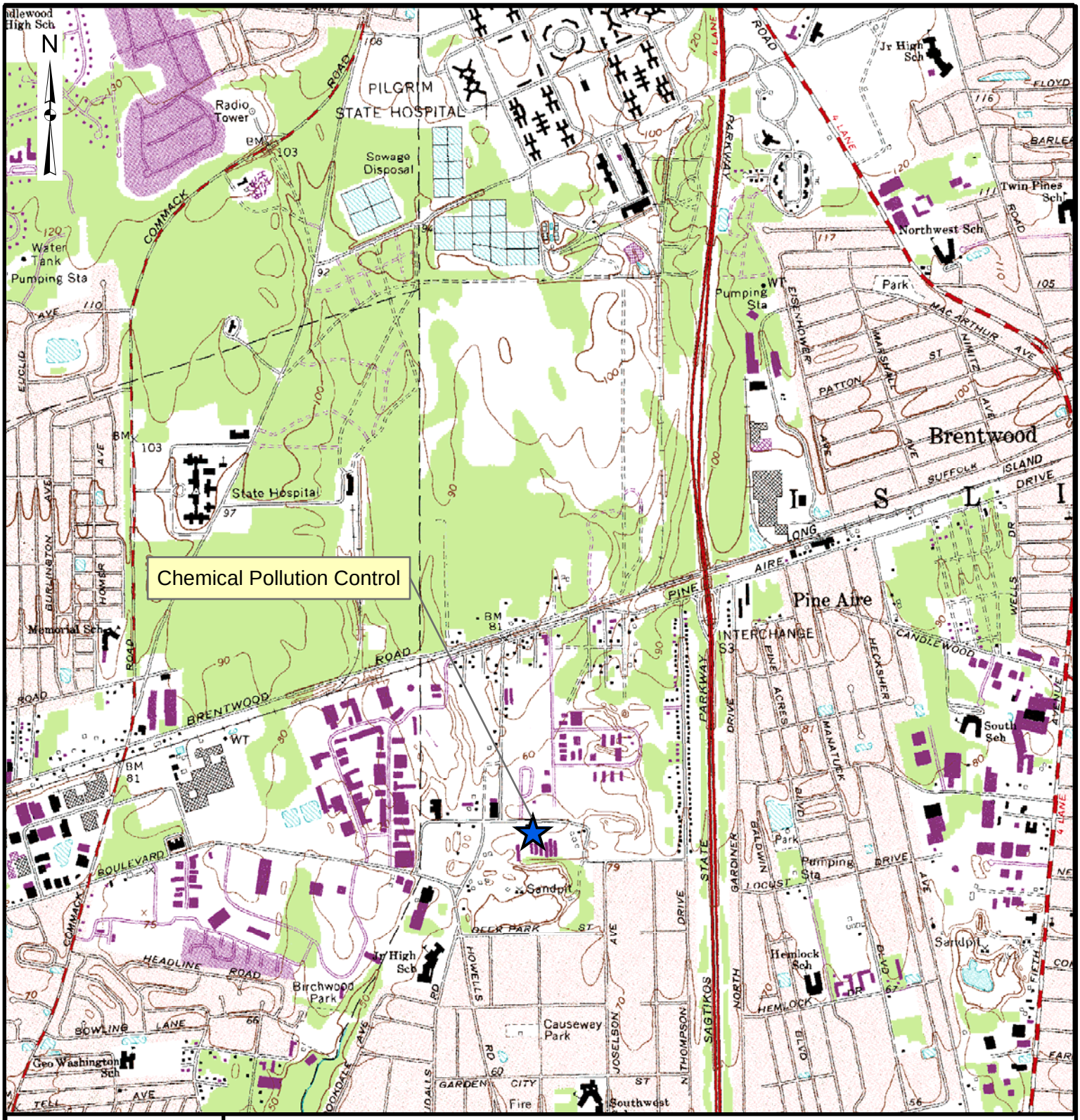
Several VOC and SVOC analytes were detected in dry well sediment samples, but below there respective 6 NYCRR Part 375 Environmental Remediation Programs-Unrestricted Use Standards.

Chromium was detected in DW-02 10-11' (6.7 mg/Kg) and DW-04 8-9' (28 mg/Kg) above the Part 375 standard of 1 mg/Kg. Lead (390 mg/Kg), silver (17 mg/kg), and zinc (140 mg/Kg) were all detected above the respective Part 375 standards in sample DW-04 8-9'. No other metals were detected above the respective standards.

Several pesticides were detected in DW-04 8-9' above the Part 375 standards, including p,p'-DDD (0.019 mg/Kg), p,p'-DDE (0.0036 mg/Kg) and p,p'-DDT (0.075 mg/Kg). Pesticides were not observed in any of the remaining dry well samples above the respective standards.

No PCBs were detected in any of the dry well sediment samples.

Validated analytical results for dry well sediment samples are included with the DUSR provided in Appendix D.

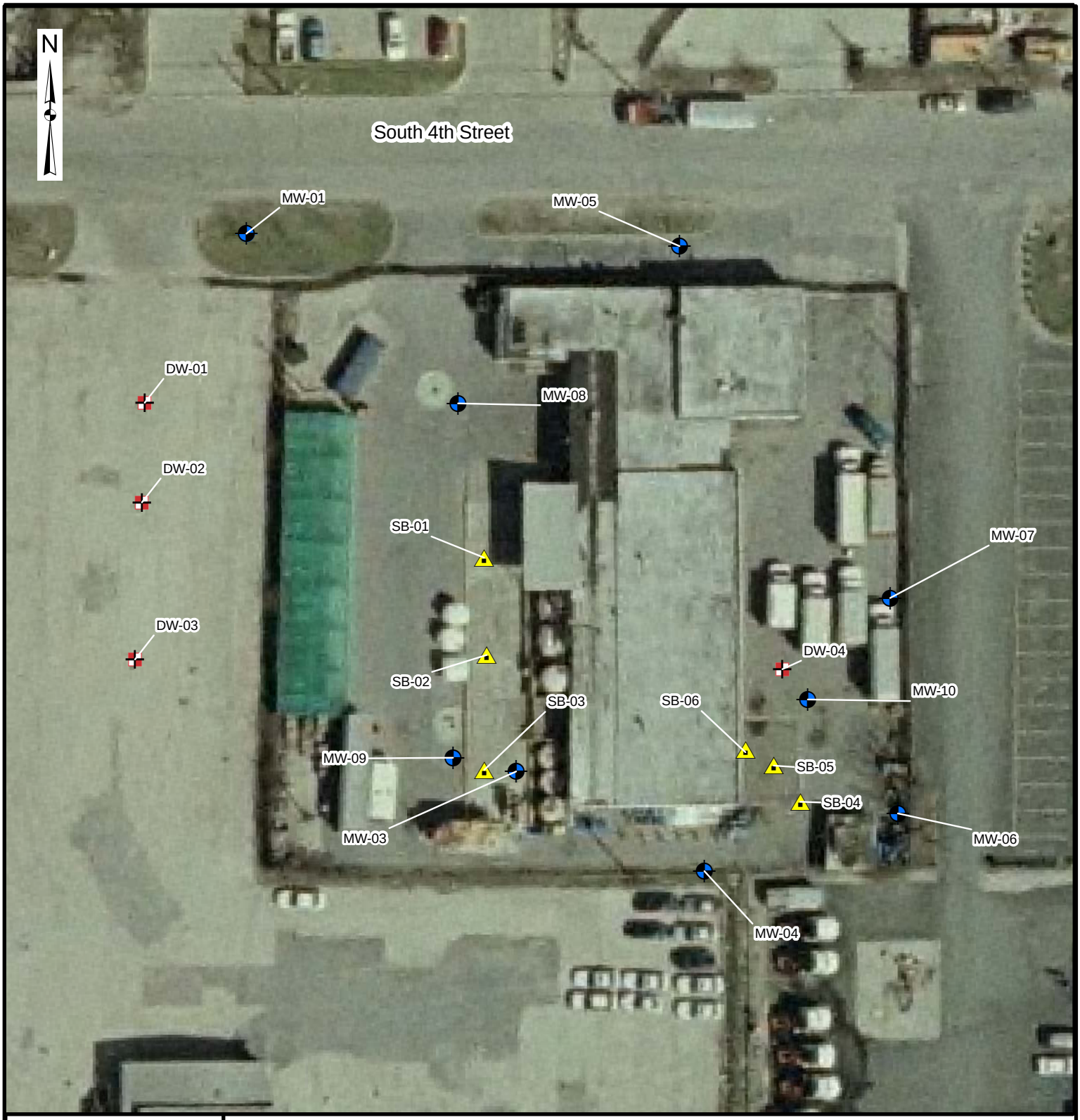


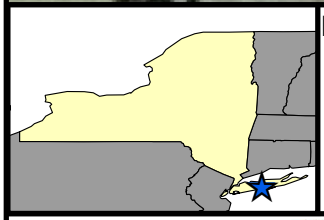
Source: NYSDEC, NYS GIS Clearinghouse 2005

CHEMICAL POLLUTION CONTROL (SITE #152015)
INVESTIGATION
SUMMARY REPORT
BAY SHORE, SUFFOLK COUNTY, NEW YORK

FIGURE 1
Site Location Map

PROJECT MGR:	DESIGNED BY:	CREATED BY:	CHECKED BY:	SCALE:	DATE:	PROJECT NO:	FILE NO:
JCH	CJS	CJS	JAV	AS SHOWN	SEPTEMBER 2007	14368.12	GIS/PROJECTS/ FIGURE1.MXD





Legend

- Drywell
- Monitoring Well
- Soil Boring

Feet

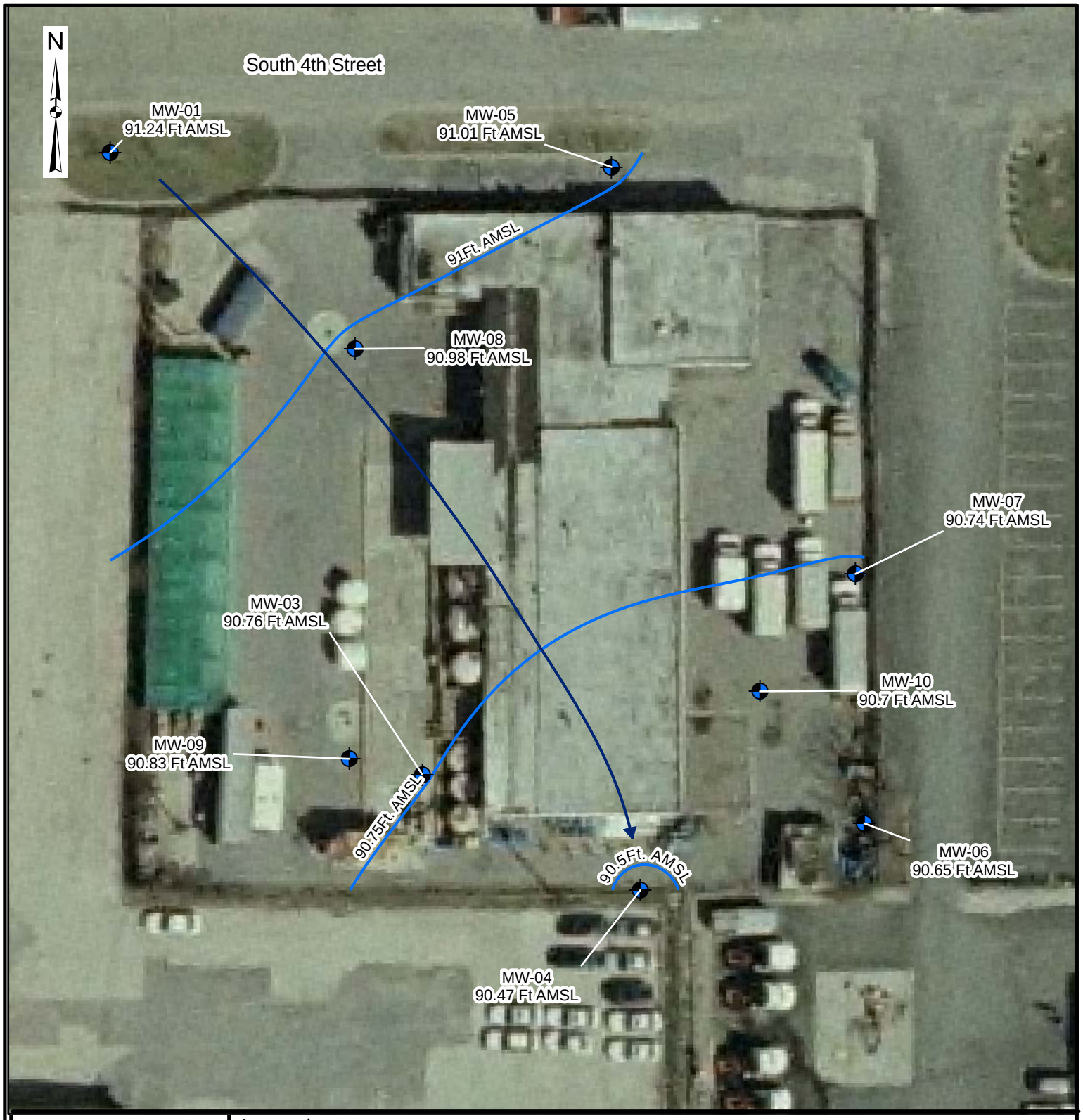
0 25 50 100

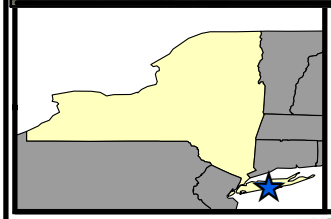
Source: NYSDEC, NYS GIS Clearinghouse 2005

CHEMICAL POLLUTION CONTROL (SITE #152015)
INVESTIGATION
SUMMARY REPORT
BAY SHORE, SUFFOLK COUNTY, NEW YORK

FIGURE 2
Site Sample Locations

PROJECT MGR:	DESIGNED BY:	CREATED BY:	CHECKED BY:	SCALE:	DATE:	PROJECT NO:	FILE NO:
JCH	CJS	CJS	JAV	AS SHOWN	SEPTEMBER 2007	14368.12	GIS/PROJECTS/ FIGURE2.MXD





Legend

- Monitoring Well
- Groundwater Contour
- Groundwater Flow Direction

Feet

0 25 50 100

Source: NYSDEC, NYS GIS Clearinghouse 2005

CHEMICAL POLLUTION CONTROL (SITE #152015)
INVESTIGATION
SUMMARY REPORT
BAY SHORE, SUFFOLK COUNTY, NEW YORK

FIGURE 3
Groundwater
Potentiometric Surface

PROJECT MGR:	DESIGNED BY:	CREATED BY:	CHECKED BY:	SCALE:	DATE:	PROJECT NO:	FILE NO:
JCH	CJS	CJS	JAV	AS SHOWN	SEPTEMBER 2007	14368.12	GIS/PROJECTS/ FIGURE3.MXD

TABLE 1 GROUNDWATER ANALYTICAL RESULTS 14 AUGUST 2007

Chemical Compound	MW-01	MW-03	MW-04	MW-05	MW-06	MW-07	MW-08	MW-09	MW-10	DUP (A)	AWQS (B)
VOLATILE ORGANIC COMPOUNDS (VOC) U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8260B (ug/L)											
cis-1,2-Dichloroethene	<5 U	1.4 J	**320**	<5 U	4.3 J	<5 U	**14**	<5 U	1.2 J	**20**	5
1,4-Dichlorobenzene	<5 U	<5 U	< 50 U	<5 U	<5 U	1.2 J	<5 U	<5 U	<5 U	<5 U	3
Tetrachloroethene	<5 U	4.4 J	**14 J**	<5 U	1.4 J	<5 U	1.1 J	1.7 J	1.1 J	1.8 J	5
Toluene	<1 U	<1.0 U	<10 U	<1 U	1.5	<1 U	<1.0 U	<1 U	<1 U	<1 U	5
Trichloroethene	<5 U	**6.6**	**330**	<5 U	**6.7**	<5 U	**15**	**6.0**	4.1 J	**19**	5
Total VOC	U	12.4	664	U	13.9	U	30.1	7.7	6.4	40.8	
TARGET ANALYTE LIST METALS U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 6010 (ug/L)											
Aluminum	<180 U	<180 U	<180 U	<180 U	<180 U	<180 U	250	<180 U	300	320	NL
Barium	<50 U	97	<50 U	<50 U	<50 U	<50 U	210	<50 U	58	210	1,000
Calcium	24,000	16,000	56,000	58,000	25,000	29,000	25,000	30,000	25,000	26,000	NL
Iron	<280 U	<280 U	<280 U	<280 U	**1,100**	<280 U	**530**	**340**	**710**	**660**	300
Lead	<4 U	<4 U	<4 U	<4 U	<4 U	<4 U	<4 U	<4 U	10	5.6	25
Magnesium	2,300	<2,000 U	4,600	3,100	2,200	2,900	2,500	2,400	<2,000 U	2,600	35,000*
Manganese	<40 U	74	<40 U	83	110	130	81	<40 U	330	83	300
Potassium	26,000	<5,000 U	11,000	26,000	8,700	13,000	13,000	16,000	17,000	13,000	NL
Sodium	**22,000**	<5,000 U	**25,000**	**22,000**	10,000	15,000	**20,000**	**27,000**	15,000	**20,000**	20,000

(A) The duplicate sample was collected from MW-08

(B) New York State Department of Environmental Conservation (NYSDEC). 1998. Division of Water Technical and Operational Guidance Series (1.1.1). Ambient Water Quality Standards (AWQS) and Guidance Values and Groundwater Effluent Limitations. Standards are compared to Class GA Standards (Source of drinking water-groundwater).

NOTE:
U = Analyzed but not reported at concentration above reporting limit. Sample quantitation limits are shown as (<__U)
J = Estimated Value, concentration below laboratory reporting limit
BOLD = Concentrations detected above the AWQS and Guidance Values and Groundwater Effluent Limitations.
NL = No Standard Listed
* = Guidance Value

SEMIVOLATILE ORGANIC COMPOUNDS U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8270B (ug/L)											
Bis (2-Ethylhexyl)phthalate	<10 U	<10 U	<10 U	<10 U	<10 U	1.2	1.5 J	<10 U	1.1 J	<10 U	5
PESTICIDES U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8081(ug/L)											
No Pesticides detected above laboratory detection limit											
POLYCHLORINATED BIPHENYLS (PCB) U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8082 (ug/L)											
No PCBs detected above laboratory detection limit											

TABLE 2 SOIL ANALYTICAL RESULTS 16 AUGUST 2007

Chemical Compound	SB-01 1.5-2.5'	SB-02 1-5'	SB-02 5-7'	SB-02 7-9'	SB-03 1.5-3.5'	SB-04 1.5-2.5'	SB-05 1-3'	SB-05 3-5'	SB-05 5-7'	SB-05 7-9'	SB-06 1.5-2.5'	DUP [A]	Table 375-6.8[B]
VOLATILE ORGANIC COMPOUNDS (VOC) U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8260B (mg/Kg)													
Acetone	<0.026 U	<0.027 U	<0.026 UJ	<0.026 UJ	<0.027 U	< 0.028 UJ	<0.027 U	<0.026 U	<0.026 U	<0.026 U	0.13 U	<0.027 U	0.05
Cis-1,2-Dichloroethene	0.0014 J	<0.0053 U	<0.0052 UJ	<0.0052 UJ	<0.0054 U	<0.0056 UJ	<0.0053 U	<0.0052 U	<0.0051 U	<0.0051 U	<0.027 UJ	0.0017 J	0.25
Ethylbenzene	<0.0011 U	<0.0011 U	<0.0010 UJ	<0.0010 UJ	<0.0011 U	< 0.0011 UJ	<0.0011 U	<0.0010 U	<0.0010 U	<0.0010 U	0.0069J	<0.0011 U	1
m&p-Xylene	<0.0021 U	<0.0021 U	<0.0021 UJ	<0.0021 UJ	0.0020 J	< 0.0022 UJ	< 0.0021UJ	<0.0021 U	<0.0020 U	<0.0020 U	0.021J	0.0015 J	NL
Methylene chloride	0.0080 U	0.0088 U	0.011 U	0.0073 U	0.0072 U	0.0068 U	0.0066 U	0.0069 U	0.0065 U	0.0066	0.027 U	0.0089 U	0.05
Tetrachloroethene	0.0093	0.0039	<0.0010 UJ	<0.0010 UJ	0.003	< 0.0011 UJ	<0.0011 U	<0.0010 U	<0.0010 U	<0.0010 U	0.33	0.0049	1.3
Tolune	0.0028	0.0026	<0.0010 UJ	<0.0010 UJ	<0.0011 U	< 0.0011 UJ	<0.0011 U	0.0012	<0.0010 U	<0.0010 U	<0.0054 UJ	0.0038	0.7
Trichloroethene	0.018	0.010	<0.0052 UJ	<0.0052 UJ	<0.0054 U	<0.0056 UJ	<0.0053 U	<0.0052 U	<0.0051 U	<0.0051 U	0.016 J	0.015	0.47
Total Volatile Organic Compounds	0.0395	0.0253	0.011	0.0073	0.0122	0.009	0.0087	0.0081	0.0065	0.0066	0.4919	0.0358	
SEMIVOLATILE ORGANIC COMPOUNDS U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8270B (mg/Kg)													
Benzo (a) anthracene	<0.35 U	<0.35 U	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	0.059 J	<0.34 U	<0.34 U	<0.36 U	<0.36 U	1
Benzo (a) pyrene	<0.35 U	<0.35 U	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	0.049 J	<0.34 U	<0.34 U	<0.36 U	<0.36 U	1
Benzo (b) fluoranthene	<0.35 U	<0.35 U	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	0.071 J	<0.34 U	<0.34 U	<0.36 U	<0.36 U	1
Benzo (g,h,i) perylene	<0.35 U	<0.35 U	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	0.041 J	<0.34 U	<0.34 U	<0.36 U	<0.36 U	100
Bis (2-Ethylhexyl)phthalate	<0.35 U	0.084 J	0.55	0.067 J	<0.36 U	0.17 J	0.11 J	0.14 J	0.060 J	0.051 J	0.13 J	0.085 J	NL
Chrysene	<0.35 U	<0.35 U	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	0.068 J	<0.34 U	<0.34 U	<0.36 U	<0.36 U	1
Di-n-butylphthalate	<0.35 U	0.038 J	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	<0.35 U	<0.34 U	<0.34 U	<0.36 U	<0.36 U	NL
Fluoranthene	<0.35 U	<0.35 U	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	0.12 J	<0.34 U	<0.34 U	<0.36 U	<0.36 U	100
Phenanthrene	<0.35 U	<0.35 U	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	0.070 J	<0.34 U	<0.34 U	<0.36 U	<0.36 U	100
Pyrene	<0.35 U	<0.35 U	<0.35 U	<0.34 U	<0.36 U	<0.37 U	<0.35 U	0.13 J	<0.34 U	<0.34 U	<0.36 U	<0.36 U	100

(A) The duplicate sample was collected from SB02-1-5'
(B) 6 NYCRR Part 375 Environmental Remediation Programs-Unrestricted Use Standards. 14 December 2006.

NOTE: U = Analyzed but not reported at concentrations above reporting limit. Sample quantitation limits are shown as <__U
J = Indicates an estimated value when a compound is detected at less than the specified detection limit
B = Indicates the analyte was found in the blank as well as in the sample
BOLD = Concentrations detected above the NYCRR Subpart 375-6 Remedial Program Soil Cleanup Objective
NL = Not Listed

Chemical Compound	SB-01 1.5-2.5'	SB-02 1-5'	SB-02 5-7'	SB-02 7-9'	SB-03 1.5-3.5'	SB-04 1.5-2.5'	SB-05 1-3'	SB-05 3-5'	SB-05 5-7'	SB-05 7-9'	SB-06 1.5-2.5'	DUP [A]	Table 375-6.8[B]
TARGET ANALYTE LIST METALS U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 6010 (mg/Kg)													
Aluminum	3,900	2,100	1,900	650	2,700	9,000	5,800	2,700	430	680	5,800	1,100	NL
Arsenic	< 2.1 U	< 2.1 U	< 2.1 U	< 2.1 U	< 2.2 U	3.2	2.5	< 2.1 U	< 2.0 U	< 2.0 U	<2.2	< 4.3 U	13
Barium	13	< 11 U	25	< 10 U	< 11 U	18	13	< 10 U	< 10 U	< 10 U	13	< 11 U	350
Calcium	< 1,100 U	< 1,100 U	24,000	< 1,000 U	< 1,100 U	< 1,100 U	< 1,100 U	< 1,000 U	< 1,000 U	< 1,000 U	< 1,100 U	130,000	NL
Chromium	< 5.3 U	< 5.3 U	**13**	< 5.2 U	**180 J**	**11**	**9.8**	< 5.2 U	< 5.1 U	< 5.1 U	**6.8**	< 5.4 U	1
Cobalt	< 2.6 U	< 2.7 U	< 2.6 U	< 2.6 U	< 2.7 U	4.4	2.9	< 2.6 U	< 2.6 U	< 2.6 U	< 2.7 U	< 11 U	NL
Copper	6	8.5	9.4	5.5	31	8.7	7.7	< 5.2 U	< 5.1 U	< 5.1 U	6.3	< 11 U	50
Iron	5,800	4,300	3,500	1,400	4,400	16,000	6,900	3,700	2,200	2,100	6,700	3,800	NL
Lead	< 5.3 U	< 5.3 U	< 5.2 U	< 5.2 U	< 5.4 U	5.7	5.6	< 5.2 U	< 5.1 U	< 5.1 U	< 5.4 U	< 11 U	63
Magnesium	770	< 530 U	7,800	< 520 U	< 540 U	1,400	930	< 520 U	< 510 U	< 510 U	850	78,000	NL
Manganese	120	64	70	53	70 J	98	75	53	52	75	83	120	1,600
Nickel	< 5.3 U	< 5.3 U	6.6	< 5.2 U	< 5.4 U	< 5.6 U	< 5.3 U	< 5.2 U	< 5.1 U	< 5.1 U	< 5.4 U	< 5.4 U	30
Silver	< 1.6 U	< 1.6 U	**3.4**	< 1.5 U	< 1.6 U	< 1.7 U	< 1.6 U	< 1.6 U	< 1.5 U	< 1.5 U	< 1.6 U	< 3.3 U	2
Vanadium	< 11 U	< 11 U	< 10 U	< 10 U	< 11 U	16	11	< 10 U	< 10 U	< 10 U	< 11 U	< 11 U	NL
Zinc	< 11 U	< 11 U	< 10 U	< 10 U	< 11 U	18	13	< 10 U	< 10 U	< 10 U	11	< 11 U	109
PESTICIDES U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8081(mg/Kg)													
Chlordane	<0.011 U	<0.011 U	0.021	<0.010 U	<0.011 U	<0.011 U	<0.011 U	<0.010 U	<0.010 U	<0.010 U	<0.011 U	<0.011 U	0.094
POLYCHLORINATED BIPHENYLS (PCBs) U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8082 (mg/Kg)													
No PCBs detected above laboratory detection limit													

TABLE 3 DRY WELL ANALYTICAL RESULTS 16 AUGUST 2007

Chemical Compound	DW-01 10-11'	DW-02 10-11'	DW-03 10-11'	DW-04 8-9'	DW-04 13-14'	Table 375-6.8(A)
VOLATILE ORGANIC COMPOUNDS U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8260B (mg/Kg)						
Acetone	<0.029 U	0.039 U	< 0.028 U	<0.029 U	<0.030 UJ	0.05
Benzene	0.0032	<0.0012 U	<0.0011 U	<0.0012 U	<0.0012 UJ	0.06
Carbon disulfide	<0.0058 UJ	0.0028 J	<0.0056 UJ	<0.0059 UJ	<0.0061 UJ	NL
Ethylbenzene	0.0065	<0.0012 U	<0.0011 U	0.0012	<0.0012 UJ	1
m&p-Xylene	0.027	<0.0024 U	<0.0022 U	0.0032	<0.0024 UJ	NL
Methylene chloride	0.010 U	0.0096 U	0.015 U	0.0060 U	0.014 U	0.05
o-Xylene	0.008	<0.0012 U	<0.0011 U	0.0033	<0.0012 UJ	NL
Toluene	0.029	0.0021	0.0016	<0.0012 U	<0.0012 UJ	0.7
Total Volatile Organic Compound	0.0837	0.0535	0.0166	0.0137	0.014	

(A) 6 NYCRR Part 375 Environmental Remediation Programs-Unrestricted Use Standards. 14 December 2006.

NOTE: U = Analyzed but not reported at concentrations above reporting limit. Sample quantitation limits are shown as <__U
J = Indicates an estimated value when a compound is detected at less than the specified detection limit
B = Indicates the analyte was found in the blank as well as the sample
NL = Not Listed
BOLD = Concentrations detected above the NYCRR Subpart 375-6 Remedial Program Soil Cleanup Objective.

Chemical Compound	DW-01 10-11'	DW-02 10-11'	DW-03 10-11'	DW-04 8-9'	DW-04 13-14'	Table 375-6.8[(A)]
SEMI-VOLATILE ORGANIC COMPOUNDS U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8270B (mg/Kg)						
2-Methylnaphthalene	<0.39 U	<0.40 U	<0.037 U	0.97	<0.41 U	NL
Benzo (a) anthracene	<0.39 U	<0.40 U	<0.037 U	0.20 J	<0.41 U	1
Benzo (a) pyrene	<0.39 U	0.050 J	<0.037 U	0.18 J	<0.41 U	1
Benzo (b) fluoranthene	<0.39 U	0.072 J	<0.037 U	0.43 J	<0.41 U	1
Benzo (g,h,i) perylene	<0.39 U	0.061 J	<0.037 U	0.19 J	<0.41 U	100
Benzo (k) fluoranthene	<0.39 U	<0.40 U	<0.037 U	0.10 J	<0.41 U	0.8
Bis (2-Ethylhexyl)phthalate	0.049 J	0.42	0.10 J	7.1	0.053 J	NL
Butylbenzylphthalate	<0.39 U	<0.40 U	<0.037 U	0.63 J	<0.41 U	NL
Chrysene	<0.39 U	0.052 J	<0.037 U	0.39 J	<0.41 U	1
Di-n-butylphthalate	<0.39 U	<0.40 U	0.070 J	0.23 J	<0.41 U	NL
Fluoranthene	<0.39 U	0.071 J	<0.037 U	0.61 J	<0.41 U	100
Fluorene	<0.39 U	<0.40 U	<0.037 U	0.12 J	<0.41 U	30
Indeno (1,2,3-cd) pyrene	<0.39 U	0.042 J	<0.037 U	0.16 J	<0.41 U	0.5
Naphthalene	<0.39 U	<0.40 U	<0.037 U	0.47 J	<0.41 U	12
Phenanthrene	<0.39 U	<0.40 U	<0.037 U	0.48 J	<0.41 U	100
Pyrene	<0.39 U	0.088 J	<0.037 U	0.61 J	<0.41 U	100

Chemical Compound	DW-01 10-11'	DW-02 10-11'	DW-03 10-11'	DW-04 8-9'	DW-04 13-14'	Table 375-6.8[(A)]
TARGET ANALYTE LIST METALS U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 6010 (mg/Kg)						
Aluminum	530	1,600	530	1,800	2,300	NL
Arsenic	<2.3 U	<2.4 U	<2.2 U	2.7	<2.4 U	13
Barium	<12 U	24	<11 U	51	< 12	350
Cadmium	<0.70 U	<0.72 U	<0.67 U	1.7	<0.73 U	2.5
Calcium	<1200 U	2,400	<1100 U	1,700	<1200 U	NL
Chromium	<5.8 U	**6.7 J**	<5.6 U	**28**	<6.1 U	1
Copper	<5.8 U	13	<5.6 U	42	<6.1 U	50
Iron	1,400	1,900	1,700	5,600	4,400	NL
Lead	<5.8 U	16	<5.6 U	**390**	<6.1 U	63
Magnesium	<580 U	930	<560 U	1,200	1,100	NL
Manganese	30	85	20	28	36	1,600
Nickel	<5.8 U	<6.0 U	<5.6 U	6.2	<6.1 U	30
Potassium	<580 U	<600 U	<560 U	<590 U	810	NL
Silver	<1.7 U	<1.8 U	<1.7 U	**17**	<1.8 U	2
Vanadium	<12 U	17	<11 U	<120 U	<12	NL
Zinc	<12 U	68	<11 U	**140**	17	109
PESTICIDES U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8081(mg/Kg)						
Chlordane	<0.012 U	0.021	<0.011 U	0.05	<0.012 U	0.094
p,p'-DDD	<0.0029 U	<0.0030 U	<0.0028 U	**0.019**	<0.0030 U	0.0033
p,p'-DDE	<0.0029 U	<0.0030 U	<0.0028 U	**0.0036 D**	<0.0030 U	0.0033
p,p'-DDT	<0.0029 UJ	<0.0030 UJ	<0.0028 UJ	**0.075J**	<0.0030 UJ	0.0033
POLYCHLORINATED BIPHENYLS (PCBS) U.S. ENVIRONMENTAL PROTECTION AGENCY METHOD 8082 (mg/Kg)						
No PCBs detected above laboratory detection limit						

Appendix A

Daily Field Reports

DAILY OBSERVATION REPORT **Day: 1** **Date: August 14, 2007**

NYSDEC

Temperature: (F) 65 (am) 80 (pm)

Wind Direction: 2 W (am) 2 W (pm)

Weather: (am) Sunny
(pm) Sunny

Project Name
Chemical Pollution Control
NYSDEC Site # 152015

Contract #

Location: Bay Shore, New York

Arrive at site 0630 (am)

Leave site: 1500 (pm)

HEALTH & SAFETY:

Are there any changes to the Health & Safety Plan? Yes ( ) No (x)
(If yes, list the deviation under items for concern)

Are monitoring results at acceptable levels?

Soil	Yes (x)	n/a ( )	* No ( )
Waters	Yes (x)	n/a ( )	* No ( )
Air	Yes (x)	n/a ( )	* No ( )

- *If No, provide comments*

OTHER ITEMS:

Site Sketch Attached: Yes ( ) No (X)
Photos Taken: Yes ( ) No (X)

DESCRIPTION OF DAILY WORK PERFORMED:

Arrived onsite, and walked site. All wells are accessible, except MW02 which is destroyed. Gauged all wells in order to obtain groundwater elevations for groundwater contour map. Purged all monitoring wells with a peristaltic pump and dedicated polyethelyne tubing. Collected all samples and put in a cooler on Ice. Once all samples were collected called and set up a sample pick up time with Hampton-Clarke-Veritech laboratories. Wells MW01 and MW05 were locked, had to get key from CPC. Locked wells after sampling.

SAMPLING (Soil/Water/Air) Collected

Contractor Sample ID:	DEC ID:	Description:
CPC-MW01	NA	GROUNDWATER SAMPLE FROM MW01
CPC-MW03	NA	GROUNDWATER SAMPLE FROM MW03
CPC-MW04	NA	GROUNDWATER SAMPLE FROM MW04
CPC-MW05	NA	GROUNDWATER SAMPLE FROM MW05
CPC-MW-06	NA	GROUNDWATER SAMPLE FROM MW06
CPC-MW07	NA	GROUNDWATER SAMPLE FROM MW07
CPC-MW08	NA	GROUNDWATER SAMPLE FROM MW08
CPC-MW09	NA	GROUNDWATER SAMPLE FROM MW09
CPC-MW10	NA	GROUNDWATER SAMPLE FROM MW09
CPC-MW10 (MS/MSD)	NA	MATRIX SPIKE / MATRIX SPIKE DUPLICATE MW10
CPC-DUPLICATE	NA	DUPLICATE SAMPLE COLLECTED FROM MW08

DAILY OBSERVATION REPORT **Day: 1** **Date: August 14, 2007**

_____

CONTRACTOR/SUBCONTRACTOR EQUIPMENT AND PERSONNEL ON SITE:

(Name of contractor) personnel: NA

(Name of Subcontractor) personnel:

(Name of contractor) equipment:

(*Indicates active equipment)

Other Subcontractors:

VISITORS TO SITE:

1. None

PROJECT SCHEDULE ISSUES:

None.

PROJECT BUDGET ISSUES:

None.

ITEMS OF CONCERN:

MW02 was destroyed and unable to be sampled. MW06 was gauged with a total depth of 9.80-ft below ground surface, the well log shows a total depth of 24-ft below ground surface. Approximately 14-ft of material (soft, sandy) has entered the well. When sampling, noticed the well sections were not put together tight, this may be where some the material has entered the well.

COMMENTS:

ATTACHMENT(S) TO THIS REPORT:

SITE REPRESENTATIVE: **Joseph Von Uderitz-Geologist EA Engineering, Science, Tech.**

Name: (*signature*)
cc:

DAILY PHOTOLOG

DAILY OBSERVATION REPORT **Day: 2** **Date: August 16, 2007**

NYSDEC

Temperature: (F) 65 (am) 80 (pm)

Wind Direction: 2 W (am) 2 W (pm)

Weather: (am) Sunny
(pm) Sunny

Project Name
Chemical Pollution Control
NYSDEC Site # 152015

Contract #

Location: Bay Shore, New York

Arrive at site 0645 (am)

Leave site: 1300 (pm)

HEALTH & SAFETY:

Are there any changes to the Health & Safety Plan? Yes ( ) No (x)
(If yes, list the deviation under items for concern)

Are monitoring results at acceptable levels?

Soil	Yes (x)	n/a ( )	* No ( )
Waters	Yes (x)	n/a ( )	* No ( )
Air	Yes (x)	n/a ( )	* No ( )

- *If No, provide comments*

OTHER ITEMS:

Site Sketch Attached: Yes ( ) No (X)
Photos Taken: Yes ( ) No (X)

DESCRIPTION OF DAILY WORK PERFORMED:

Arrived onsite, and walked site to make sure all drilling locations were accessible. Started sampling three Dry Wells (DW01-DW3) on offsite property and collected samples. Once sampling was completed moved to CPC and installed six soil borings, collected samples as appropriate. Collected two samples from DW04. Once all sampling was complete, did a site survey with a GPS unit. Surveyed in all monitoring well locations, soil boring locations and dry well locations. Hand delivered samples to Hampton-Clarke-Veritech Laboratories in West Fairfield, NJ.

SAMPLING (Soil/Water/Air) Collected

Contractor Sample ID:	DEC ID:	Description:
CPC-SB01-1.5-2.5’	NA	SOIL SAMPLE FROM SB01
CPC-SB02-1-5’	NA	SOIL SAMPLE FROM SB02
CPC-SB02-5-7’	NA	SOIL SAMPLE FROM SB02
CPC-SB02-7-9’	NA	SOIL SAMPLE FROM SB02
CPC-SB03-1.5-3.5’	NA	SOIL SAMPLE FROM SB03
CPC-SB04-1.5-2.5	NA	SOIL SAMPLE FROM SB04
CPC-SB05-1-3’	NA	SOIL SAMPLE FROM SB05
CPC-SB05-3-5’	NA	SOIL SAMPLE FROM SB05
CPC-SB05-5-7’	NA	SOIL SAMPLE FROM SB05
CPC-SB05-7-9’	NA	SOIL SAMPLE FROM SB05
CPC-SB06-1.5-2.5’	NA	SOIL SAMPLE FROM SB06

CPC-DW01-10-11'	NA	SOIL SAMPLE FROM DW01
CPC-DW02-10-11'	NA	SOIL SAMPLE FROM DW02
CPC-DW03-10-11'	NA	SOIL SAMPLE FROM DW03
CPC-DW04-8-9'	NA	SOIL SAMPLE FROM DW04
CPC-DW04-13-14'	NA	SOIL SAMPLE FROM DW05
CPC-SOIL DUPLICATE	NA	SOIL SAMPLE DUPLICATE FROM SB02
CPC-SB03-1.5-3.5' (MS/MSD)	NA	MATRIX SPIKE / MATRIX SPIKE DUPLICATE SB03
CPC-RINSATE	NA	EQUIPMENT BLANK

CONTRACTOR/SUBCONTRACTOR EQUIPMENT AND PERSONNEL ON SITE:

(Name of contractor) personnel: Land Air and Water Environmental

(Name of Subcontractor) personnel: Nesto Santiago and Eric Bedell

(Name of contractor) equipment:

(*Indicates active equipment)

Other Subcontractors: None

VISITORS TO SITE:

1. NYSDEC Personnel: Chek Ng, Bob Corcoran
2. NYSDOH: Renata Ockerby

PROJECT SCHEDULE ISSUES:

None.

PROJECT BUDGET ISSUES:

None.

ITEMS OF CONCERN:

None

COMMENTS:

None

ATTACHMENT(S) TO THIS REPORT:

SITE REPRESENTATIVE: **Joseph Von Uderitz-Geologist EA Engineering, Science, Tech.**

Name: (*signature*)
cc:

DAILY PHOTOLOG

Appendix B

Groundwater Sampling Forms

EA Engineering PC and its Affliate,
EA Science and Technology

GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW-01	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: Western most well on South St.	**Well Condition:** Good	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 4-inch

Purge Date: 8/14/2007	**Purge Time:** 1329 to 1353 (24 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 17.07	**D. Well Volume (ft):** 0.6532	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 6.75	**E. Well Volume (gal) C*D):** 6.741024	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 10.32	**F. Three Well Volumes (gal) (E3):** 20.223072	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
1332	6.75	0.3	0.1	6.41	65	19.51	0.439	1.01	32
1335	6.75	0.6	0.1	6.61	60	19.57	0.682	0.79	28
1338	6.75	0.9	0.1	6.61	61	19.40	0.726	0.56	29
1341	6.75	1.2	0.1	6.61	62	19.29	0.806	0.52	45
1344	6.75	1.5	0.1	6.64	61	19.26	0.810	0.53	45
1347	6.75	1.8	0.1	6.64	62	19.36	0.795	0.59	50
1350	6.75	2.1	0.1	6.65	63	19.46	0.802	0.56	52
1353	6.75	2.4	0.1	6.65	62	19.51	0.806	0.51	50

Total Quantity of Water Removed (gal):	2.4	**Sampling Time:**	1355
Samplers:	Joe Von Uderitz	**Split Sample With:**	NA
Sampling Date:	8/14/2007	**Sample Type:**	Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors. Well in good condition.

EA Engineering PC and its Affliate,
EA Science and Technology

GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW-03	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: Southern well on loading dock	**Well Condition:** Good	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 4-inch

Purge Date: 8/14/2007	**Purge Time:** 1005 to 1029 (24 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 17.72	**D. Well Volume (ft):** 0.6532	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 8.42	**E. Well Volume (gal) C*D):** 6.07476	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 9.3	**F. Three Well Volumes (gal) (E3):** 18.22428	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
1008	8.42	0.3	0.1	6.69	57	21.66	1.580	0.05	1
1011	8.42	0.6	0.1	6.67	51	21.64	0.473	0	0
1014	8.42	0.9	0.1	6.66	49	21.68	0.431	0	0
1017	8.42	1.2	0.1	6.67	47	21.69	0.388	0	0
1020	8.42	1.5	0.1	6.67	42	21.73	0.345	0	0
1023	8.42	1.8	0.1	6.67	33	21.71	0.340	0	0
1026	8.42	2.1	0.1	6.68	28	21.70	0.321	0	0
1029	8.42	2.4	0.1	6.68	26	21.71	0.306	0	0

Total Quantity of Water Removed (gal): 2.4
Samplers: Joe Von Uderitz
Sampling Date: 8/14/2007
Sampling Time: 1030
Split Sample With: NA
Sample Type: Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors. Well in good condition.

EA Engineering PC and its Affliate,
EA Science and Technology

GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW-04	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: On fence at southern property line	**Well Condition:** Good	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 4-inch

Purge Date: 8/14/2007	**Purge Time:** 713 to 740 (27 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 17.59	**D. Well Volume (ft):** 0.6532	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 8.89	**E. Well Volume (gal) C*D):** 5.68284	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 8.7	**F. Three Well Volumes (gal) (E3):** 17.04852	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
716	8.91	0.3	0.1	6.71	87	20.48	9.16	5.36	2.3
713	8.92	0.6	0.1	6.53	95	20.28	5.99	4.92	2
722	8.91	0.9	0.1	6.47	102	20.20	5.10	4.60	2
725	8.91	1.2	0.1	6.47	106	20.10	4.61	4.38	1
728	8.91	1.5	0.1	6.48	107	20.06	4.40	4.27	0
731	8.91	1.8	0.1	6.50	109	20.00	4.25	4.19	0
734	8.91	2.1	0.1	6.51	111	19.93	4.14	4.16	0
737	8.91	2.4	0.1	6.52	112	19.87	4.06	4.10	0
740	8.91	2.7	0.1	6.52	112	19.81	3.99	4.11	0

Total Quantity of Water Removed (gal): 2.7
Samplers: Joe Von Uderitz
Sampling Date: 8/14/2007
Sampling Time: 740
Split Sample With: NA
Sample Type: Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors. Well in good condition.

EA Engineering PC and its Affliate,
EA Science and Technology

GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW-05	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: Front of shop on South St	**Well Condition:** Good	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 4-inch

Purge Date: 8/14/2007	**Purge Time:** 1415 to 1439 (24 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 18.34	**D. Well Volume (ft):** 0.6532	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 8.2	**E. Well Volume (gal) C*D):** 6.623448	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 10.14	**F. Three Well Volumes (gal) (E3):** 19.870344	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
1418	8.20	0.3	0.1	6.78	51	20.16	3.36	0.85	2.3
1421	8.20	0.6	0.1	6.86	49	19.85	2.45	0.34	6
1424	8.20	0.9	0.1	6.88	46	19.89	2.14	0.27	5
1427	8.20	1.2	0.1	6.89	43	19.87	1.90	0.11	4
1430	8.20	1.5	0.1	6.90	42	19.75	1.85	0.08	8
1433	8.20	1.8	0.1	6.90	41	19.83	1.92	0.03	9
1436	8.20	2.1	0.1	6.91	40	20.00	1.85	0.00	12
1439	8.20	2.4	0.1	6.91	41	20.09	1.88	0.00	13

Total Quantity of Water Removed (gal):	2.4	**Sampling Time:**	1440
Samplers:	Joe Von Uderitz	**Split Sample With:**	NA
Sampling Date:	8/14/2007	**Sample Type:**	Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors. Well in good condition.

EA Engineering PC and its Affliate,
EA Science and Technology

GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW-06	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: South east corner of property	**Well Condition:** Good	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 2-inch

Purge Date: 8/14/2007	**Purge Time:** 1415 to 1439 (24 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 9.8	**D. Well Volume (ft):** 0.1632	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 8.87	**E. Well Volume (gal) C*D):** 0.151776	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 0.93	**F. Three Well Volumes (gal) (E3):** 0.455328	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
1106	8.89	0.3	0.1	6.10	-88	20.58	3.37	0.54	35
1109	8.87	0.6	0.1	6.07	-91	20.42	2.98	0.40	27
1112	8.87	0.9	0.1	6.07	-92	20.25	2.04	0.31	9
11156	8.87	1.2	0.1	6.08	-90	20.13	1.71	0.27	1
1118	8.87	1.5	0.1	6.09	-90	20.11	1.68	0.28	1
1121	8.87	1.8	0.1	6.09	-89	20.11	1.66	0.30	1
1124	8.87	2.1	0.1	6.10	-88	20.02	1.63	0.30	1
1127	8.87	2.4	0.1	6.11	-87	20.01	1.77	0.34	5
1130	8.87	2.7	0.1	6.11	-88	20.06	1.73	0.37	3

Total Quantity of Water Removed (gal):	2.7	**Sampling Time:**	1130
Samplers:	Joe Von Uderitz	**Split Sample With:**	NA
Sampling Date:	8/14/2007	**Sample Type:**	Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors. Well in good condition.

EA Engineering PC and its Affliate,
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GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW-07	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: Eastern property line along fence	**Well Condition:** Good	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 2-inch

Purge Date: 8/14/2007	**Purge Time:** 1233 to 1300 (27 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 21.94	**D. Well Volume (ft):** 0.1632	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 8.81	**E. Well Volume (gal) C*D):** 2.142816	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 13.13	**F. Three Well Volumes (gal) (E3):** 6.428448	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
1236	8.81	0.3	0.1	6.24	58	30.16	2.65	0.17	1
1239	8.81	0.6	0.1	6.23	62	20.05	1.85	0.00	7
1242	8.81	0.9	0.1	6.25	65	19.94	1.34	0.00	3
1245	8.81	1.2	0.1	6.27	66	19.86	1.13	0.00	9
1248	8.81	1.5	0.1	6.29	66	19.80	1.12	0.00	6
1251	8.81	1.8	0.1	6.30	66	19.75	1.03	0.00	6
1254	8.81	2.1	0.1	6.31	66	19.74	1.01	0.00	5
1257	8.81	2.4	0.1	6.31	65	19.72	0.98	0.00	5
1300	8.81	2.7	0.1	6.30	66	19.71	0.94	0.00	4

Total Quantity of Water Removed (gal):	2.7	**Sampling Time:**	1300
Samplers:	Joe Von Uderitz	**Split Sample With:**	NA
Sampling Date:	8/14/2007	**Sample Type:**	Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors. Well in good condition. TD on well log is 24-ft bgs. Well had some fill material in it. Can tell that the well is not put together tight, was getting tubing stuck at a joint in the well casing.

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GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW-08	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: Just west of Office	**Well Condition:** Poor-see notes	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 2-inch

Purge Date: 8/14/2007	**Purge Time:** 829 to 850 (21 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 30	**D. Well Volume (ft):** 0.1632	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 7.87	**E. Well Volume (gal) C*D):** 3.611616	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 22.13	**F. Three Well Volumes (gal) (E3):** 10.834848	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
832	7.94	0.3	0.1	6.74	-137				
835	7.91	0.6	0.1	6.72	-138				
838	7.90	0.9	0.1	6.69	-135				
841	7.90	1.2	0.1	6.67	-127				
844	7.90	1.5	0.1	6.63	-116				
847	7.91	1.8	0.1	6.62	-110				
850	7.91	2.1	0.1	6.61	-105				

Total Quantity of Water Removed (gal):	2.7	**Sampling Time:**	1130
Samplers:	Joe Von Uderitz	**Split Sample With:**	Duplicate sample
Sampling Date:	8/14/2007	**Sample Type:**	Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors.
Compression plug is wedged down the hole, protective casing cover is sitting on top of the well. A new pad and cover should be installed. Inside of pvc has some build-up from run-off getting into the well.

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GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW-09	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: West of tanker loading dock	**Well Condition:** Good	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 2-inch

Purge Date: 8/14/2007	**Purge Time:** 922 to 949 (27 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 29.2	**D. Well Volume (ft):** 0.1632	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 8.08	**E. Well Volume (gal) C*D):** 3.446784	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 21.12	**F. Three Well Volumes (gal) (E3):** 10.340352	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
925	8.10	0.3	0.1	6.64	27	21.24	3.00	1.42	25
928	8.10	0.6	0.1	6.62	33	20.90	2.17	1.44	20
931	8.11	0.9	0.1	6.63	27	20.83	2.09	1.70	27
934	8.10	1.2	0.1	6.64	40	20.73	1.98	1.57	29
937	8.10	1.5	0.1	6.64	44	20.66	1.78	1.67	14
940	8.10	1.8	0.1	6.65	45	20.60	1.63	1.45	16
943	8.10	2.1	0.1	6.64	47	20.46	1.39	0.80	15
946	8.10	2.4	0.1	6.64	48	20.45	1.27	0.71	14
949	8.10	2.7	0.1	6.64	47	20.41	1.21	0.64	17

Total Quantity of Water Removed (gal):	2.7	**Sampling Time:**	950
Samplers:	Joe Von Uderitz	**Split Sample With:**	NA
Sampling Date:	8/14/2007	**Sample Type:**	Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors. Well in good condition.

EA Engineering PC and its Affliate,
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GROUNDWATER SAMPLING
PURGE FORM

Well I.D.: MW10	**EA Personnel:** Joe Von Uderitz	**Client:** New York State Department of Environmental Conservation (NYSDEC)
Location: West of tanker loading dock	**Well Condition:** Good	**Weather:** Sunny 80 F
Sounding Method: Interface probe	**Gauge Date:** 8/14/2007	**Measurement Ref:** PVC
Stick Up/Down (ft): Flushmount	**Gauge Time:**	**Well Diameter (in):** 2-inch

Purge Date: 8/14/2007	**Purge Time:** 1143 to 1207 (24 minutes)
Purge Method: Low-Flow Peristalic Pump	**Field Technician:**

Well Volume

A. Well Depth (ft): 29.91	**D. Well Volume (ft):** 0.1632	**Depth/Height of Top of PVC:**
B. Depth to Water (ft): 8.27	**E. Well Volume (gal) C*D):** 3.531648	**Pump Type:** Peristalic Pump
C. Liquid Depth (ft) (A-B): 21.64	**F. Three Well Volumes (gal) (E3):** 10.594944	**Pump Designation:** Peristalic Pump number 2

Water Quality Parameters

Time (hrs)	DTW (ft btoc)	Volume (gallons)	Rate (Gpm)	pH (pH units)	ORP (mV)	Temperature (oC)	Conductivity (mS/cm)	DO (mg/L)	Turbidity (ntu)
1146	8.28	0.3	0.1	6.37	-19	19.50	0.66	0.66	2
1149	8.27	0.6	0.1	6.44	-14	19.48	0.74	0.25	43
1152	8.27	0.9	0.1	6.46	-12	19.44	0.62	0.16	2
1155	8.27	1.2	0.1	6.47	-7	19.32	0.53	0.11	5
1158	8.27	1.5	0.1	6.48	-4	19.25	0.53	0.10	3
1201	8.27	1.8	0.1	6.48	-1	19.20	0.53	0.80	4
1204	8.27	2.1	0.1	6.48	0	19.20	0.49	0.80	5
1207	8.27	2.4	0.1	6.48	3	19.18	0.47	0.90	5

Total Quantity of Water Removed (gal): 2.4
Samplers: Joe Von Uderitz
Sampling Date: 8/14/2007
Sampling Time: 1210
Split Sample With: MS/MSD collected
Sample Type: Groundwater

COMMENTS AND OBSERVATIONS: Purge water was clear, no visible solids, no noticeable odors. Well in good condition.

Appendix C

Soil Boring Logs

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: DW-01
Easting 644305.028
Northing 4513787.659

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:		Geoprobe	Well Number: DW-01
Sampling Method:		Macro-Core	Sheet 1 of 1

					Drilling Start	Drilling Finish
Water Lev.						
Time					8/16/07 730	8/16/07 750
Date						
Reference						

Surface Conditions: Steel Grate
Weather: Cloudy
Temperature: 75 F

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet	USCS Log	
				0		
				1		
				2		
				3		
				4		
				5		
				6		
				7		
			0	8		Water at 8 ft below ground surface
	2.5 / 2.5					8.5-9 Black ORGANICS, Very fine black SAND
			0	9		9-11 Tan Coarse SAND with many Quartz Cobbles.
			0	10		
				11		
				12		End of boring at 11 ft bgs
				13		
				14		
				15		
				16		
				17		
				18		
				19		
				20		
				21		
				22		
				23		
				24		
				25		

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate MS/MSD

Sample Depth: 10-11 feet Sample Time: 750 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: DW-02
Easting 644304.727
Northing 4513777.28

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:	Geoprobe		Well Number: DW-02
Sampling Method:	Macro-Core		Sheet 1 of 1

					Drilling Start	Drilling Finish
Water Lev.						
Time					8/16/07 800	8/16/07 830
Date						
Reference						

Surface Conditions: Steel Grate
Weather: Cloudy
Temperature: 75 F

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet	USCS Log	
				0		
				1		
				2		
				3		
				4		
				5		
				6		
				7		
			0	8		Water at 8 ft below ground surface
	2.5 / 3		0	9		8.5-11 Black coarse SAND with trace very fine SAND. Many quartz cobbles.
			0	10		
				11		
				12		End of boring at 11 ft bgs
				13		
				14		
				15		
				16		
				17		
				18		
				19		
				20		
				21		
				22		
				23		
				24		
				25		

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate MS/MSD

Sample Depth: 10-11 feet Sample Time: 830 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: DW-03
Easting 644304.011
Northing 4513760.801

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:		Geoprobe	Well Number: DW-03
Sampling Method:		Macro-Core	Sheet 1 of 1

					Drilling Start	Drilling Finish
Water Lev.						
Time					8/16/07 830	8/16/07 845
Date						
Reference						

Surface Conditions: Steel Grade
Weather: Cloudy
Temperature: 75 F

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet	USCS Log	
				0		
				1		
				2		
				3		Water at 3 ft bgs
				4		
				5		
				6		
				7		
	3 / 3		0	8	8-9	Black fine SAND, some SILT, some coarse SAND and ORGANICS
			0	9	9-11	Light brown medium to coarse SAND with trace quartz cobbles.
			0	10		
				11		
				12		End of boring at 11 ft bgs
				13		
				14		
				15		
				16		
				17		
				18		
				19		
				20		
				21		
				22		
				23		
				24		
				25		

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate MS/MSD

Sample Depth: 10-11 feet Sample Time: 845 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: DW-04
Easting 644371.868
Northing 4513759.775

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:	Geoprobe		Well Number: DW-04
Sampling Method:	Macro-Core		Sheet 1 of 1

					Drilling Start	Drilling Finish
Water Lev.						
Time					8/16/07 1130	8/16/07 1145
Date						
Reference						

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet		USCS Log		
				0			Surface Conditions:	Steel Grate
				1			Weather:	Cloudy
				2			Temperature:	75 F
				3				
				4				
				5				
								Water at 5.5 ft bgs
				6				
				7				
	3 / 3		4.2	8			8-9	Black ORGANICS with some quartz cobbles
			2.3	9			9-10	Light gray fine to medium SAND with trace SILT, loose, non-cohesive
			0	10			10-11	Tan medium SAND with trace SILT.
			0	11				
			0	12				
	5 / 5		0	13			13-14	Light gray fine SAND
			0	14			14-17	Light gray coarse SAND with quartz cobbles
			0	15				
			0	16				
			0	17			17-18	Brown medium SAND with trace quartz
				18				
				19				
				20				
				21				
				22				
				23				
				24				
				25				

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate MS/MSD

Sample Depth: 8-9 feet Sample Time: 1130 Sample Date: 16 August 2007
Sample Depth: 13-14 feet Sample Time: 1145 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: SB-01
Easting 644340.601
Northing 4513771.454

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:		Geoprobe	Well Number: SB-01
Sampling Method:		Macro-Core	Sheet 1 of 1

					Drilling Start	Drilling Finish
Water Lev.						
Time					8/16/07 1100	8/16/07 1115
Date						
Reference						

Surface Conditions: Concrete
Weather: Cloudy
Temperature: 75 F

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet	USCS Log		
	5/5		0	0	0-1	CONCRETE and Subbase	
			0	1	1-4	Brown medium SAND with trace quartz pebbles, tight, non-cohesive.	
			0	2			
			0	3			
			0	4	4-4.5	Brown very fine SAND and SILT.	
					4.5-5	Brown medium SAND with some quartz	
				5			
				6			
				7			
				8			
				9			
				10			
				11			
				12			
				13			
				14			
				15			
				16			
				17			
				18			
				19			
				20			
				21			
				22			
				23			
				24			
				25			

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate MS/MSD

Sample Depth: 1.5-2.5 feet Sample Time: 1115 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: SB-02
Easting 644340.874
Northing 4513761.262

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:	Geoprobe		Well Number: SB-02
Sampling Method:	Macro-Core		Sheet 1 of 1

					Drilling Start	Drilling Finish
Water Lev.						
Time					8/16/07 1035	8/16/07 1055
Date						
Reference						

Surface Conditions: Concrete
Weather: Cloudy
Temperature: 75 F

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet		USCS Log	
				0	0-1		CONCRETE and Subbase
			9.1	1	1-4		Brown medium SAND with some quartz
	5/5		8.9	2			
			6.7	3			
			3.7	4	4-5		Tan medium SAND with some quartz
			4.4	5	5-6		Brown fine SAND with trace SILT. Some quartz
			4.1	6	6-9.5		Tan medium SAND with trace quartz pebbles
	4/5		4.6	7			
			3.9	8			
							Wet at approximatley 9 ft below ground surface
			4.2	9			
					9.5-10		Gray medium SAND.
			3.1	10			
				11			
				12			
				13			
				14			
				15			
				16			
				17			
				18			
				19			
				20			
				21			
				22			
				23			
				24			
				25			

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** **Duplicate 1-5 ft** MS/MSD

Sample Depth: 1-5 feet Sample Time: 1045 Sample Date: 16 August 2007
Sample Depth: 5-7 feet Sample Time: 1050 Sample Date: 16 August 2007
Sample Depth: 7-9 feet Sample Time: 1055 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: SB-03
Easting 644341.186
Northing 4513769.773

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:		Geoprobe	Well Number: SB-03
Sampling Method:		Macro-Core	Sheet 1 of 1

					Drilling	
Water Lev.					Start	Finish
Time					8/16/07 1015	8/16/07 1030
Date						
Reference						

Surface Conditions: Concrete
Weather: Cloudy
Temperature: 75 F

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet	USCS Log	
				0	0-1	CONCRETE and Subbase
			0	1	1-3.5	Brown medium SAND.
	5 / 5		0	2		
			0	3	3.5-4	Brown tight SITLY SAND. Moist.
			0	4	4-5	Tan medium SAND
			0	5		
				6		
				7		
				8		
				9		
				10		
				11		
				12		
				13		
				14		
				15		
				16		
				17		
				18		
				19		
				20		
				21		
				22		
				23		
				24		
				25		

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate **MS/MSD**

Sample Depth: 1.5-2.5 feet Sample Time: 1030 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: SB-04
Easting 644373.758
Northing 4513745.907

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:		Geoprobe	Well Number: SB-04
Sampling Method:		Macro-Core	Sheet 1 of 1

					Drilling	
Water Lev.					Start	Finish
Time					8/16/07 900	8/16/07 915
Date						
Reference						

Surface Conditions: Asphalt and Subbase
Weather: Cloudy
Temperature: 75 F

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet	USCS Log	
				0		0-0.5 Aspalt and Subbase
						0.5-3 Dark brown tight SANDY SILT, with some quartz cobbles.
			0	1		
	4 / 5		0	2		
			0	3		3-4 Tan fine to medium SAND with quartz pebbles.
			0	4		
			0	5		
				6		
				7		
				8		
				9		
				10		
				11		
				12		
				13		
				14		
				15		
				16		
				17		
				18		
				19		
				20		
				21		
				22		
				23		
				24		
				25		

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate MS/MSD

Sample Depth: 1.5-2.5 feet Sample Time: 915 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: SB-05
Easting 644370.974
Northing 4513749.694

Job. No.	Client	Location
	New York State Departement of Environmental Conservation	Chemical Pollution Control (CPC)
Drilling Method:	Geoprobe	Well Number: SB-05
Sampling Method:	Macro-Core	Sheet 1 of 1

					Drilling	
Water Lev.					Start	Finish
Time					8/16/07 900	8/16/07 915
Date						
Reference						

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet		USCS Log	Surface Conditions: Asphalt and Subbase Weather: Cloudy Temperature: 75 F
				0			0-0.5 Aspalt and Subbase
							0.5-1 Tan fine to medium SAND.
			50	1			1-2 Black medium SILTY SAND with quartz pebbles.
							*PID readings may be assoicated with newly layed asphalt.
	4 / 5		6	2			2-4 Brown tan medium SAND with trace quartz.
			5	3			
			3	4			
				5			5-6 Brown medium SAND with trace SILT, loose, non-cohesive.
			11	6			6-10 Tan medium SAND with trace quartz pebbles, wet at 9-ft
	5 / 5		1.5	7			
			0	8			
			0.3	9			
			0.1	10			
				11			
				12			
				13			
				14			
				15			
				16			
				17			
				18			
				19			
				20			
				21			
				22			
				23			
				24			
				25			

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO

Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate MS/MSD

Sample Depth: 1-3 feet Sample Time: 930 Sample Date: 16 August 2007
Sample Depth: 3-5 feet Sample Time: 935 Sample Date: 16 August 2007
Sample Depth: 5-7 feet Sample Time: 940 Sample Date: 16 August 2007
Sample Depth: 7-9 feet Sample Time: 945 Sample Date: 16 August 2007

EA Engineering, P.C.
EA Science and Technology

LOG OF SOIL BORING

Coordinates Description: SB-06
Easting 644367.142
Northing 4513751.203

Job. No.	Client	New York State Departement of Environmental Conservation	Location: Chemical Pollution Control (CPC)
Drilling Method:		Geoprobe	Well Number: SB-06
Sampling Method:		Macro-Core	Sheet 1 of 1

					Drilling Start	Drilling Finish
Water Lev.						
Time					8/16/07 900	8/16/07 915
Date						
Reference						

Blow Counts (140-lb)	Feet Drvn/Ft. Recvrd	Well Diagram	PID (ppm) HNu	Depth in Feet	USCS Log	Surface Conditions: Asphalt and Subbase Weather: Cloudy Temperature: 75 F
	4 / 5		0.3	0		0-0.5 Aspalt and Subbase
						0.5-1 Brown medium SAND
			1.2	1		1-2 Black medium SAND with trace SILT. Tight, non-cohesive. Dry
			0.6	2		2-3 Orange brown medium SAND with trace SILT.
			0.7	3		3-4 Tan medium SAND.
			0.2	4		
				5		
				6		
				7		
				8		
				9		
				10		
				11		
				12		
				13		
				14		
				15		
				16		
				17		
				18		
				19		
				20		
				21		
				22		
				23		
				24		
				25		

Logged by: Joe Von Uderitz Date: 16 August 2007

Drilling Contractor: L.A.W.E. Driller: Nesto / Erik

SOIL SAMPLE COLLECTED **YES** / NO
Samples Collected for **VOC SVOC PCB Pesticides Metals** Duplicate MS/MSD

Sample Depth: 1.5-2.5 feet Sample Time: 1000 Sample Date: 16 August 2007

Appendix D

Data Usability Summary Report

CW CHEMWORLD ENVIRONMENTAL, INC.
Environmental Consulting Services

March 5, 2008

Mr. Joe Von Uderitz
Geologist
EA Engineering , P.C.
6712 Brooklawn Parkway
Suite 104
Syracuse, New York 13211-2158

RE: Data Usability Summary Report (DUSR) #1
CPC Project
Hampton-Clarke, Inc. Veritech Laboratories, Fairfield, NJ
Lab Project Nos. 7081532 and 7081622
Water and Soil Samples
Analyses for Volatile Organics, Semi-Volatiles (Base/Neutral and Acid Extractable Organics), Pesticides, Polychlorinated Biphenyls (PCB's), Inorganics (Metals) and Cyanide

Dear Mr. Von Uderitz:

Data Usability Summary Report (DUSR) technical services were performed by ChemWorld Environmental, Inc. for the CPC Project for the water and soil sampling events of August 14 - 16, 2007. The DUSR review was performed in accordance with United States Environmental Protection Agency (USEPA) Region II data validation guidelines and New York State Department of Environmental Conservation (NYSDEC) Analytical Service Protocols (ASP) requirements, where applicable.

The analytical data from the Lab Project Nos. referenced above was reviewed (screened) for the parameters noted. The data screening consisted of a review of the Quality Control (QC) Summary Forms and a brief review of various chromatograms and quantitation reports. The QC Forms were reviewed to determine whether any data required qualification based upon QC deviations noted on the Forms. The associated Analytical Data Result Forms are included as Attachment A. These Forms include data qualifiers as described within this letter report. Unless otherwise noted, all results included on the Forms are considered usable, based upon the DUSR review items noted below. Attachment B includes copies of the associated Case Narratives and the Chain-of-Custody forms.

The DUSR review items include the following, as method appropriate:

- Completeness of Data Package
- Chain-of-Custody Review
- Holding Times from Verified Time of Sample Receipt (VTSR)
- Surrogate Recovery
- GC/MS Instrument Performance Check
- Initial and Continuing Calibration
- Matrix Spike / Matrix Spike Duplicates (MS/MSD)
- Matrix Spike Blanks (MSB)
- Internal Standards
- Method and Field Blanks
- CRDL Standards for ICP
- Laboratory Duplicate Samples
- Laboratory Control Samples (LCS)
- ICP Interference Check
- ICP Serial Dilutions

14 Orchard Way North, Rockville, Maryland 20854, Tel. 301-294-6144, Fax 301-~~309-6640~~ 294-6144

The QC Summary Forms included various deviations based upon the acceptable limits for quality control. The following should be noted regarding qualification of the data set for the review items above.

Volatiles - Water, Lab Project No. 7081532

Surrogate Recovery: Sample CPC-MW06 generated high surrogate recovery for Dibromofluoromethane at 146% (Limit 78-135). This sample was qualified as 'J', estimated, for the positive results, only, for Volatiles.

MS/MSD: The site-specific MS and MSD for CPC-MW10 generated low recovery for Trichloroethene at 64% and 63%, respectively (Limit 72-127). Sample CPC-MW10 was qualified as 'J', estimated, for the positive result for Trichloroethene. Additional qualification was not required.

Continuing Calibration: The continuing calibration analyzed on 8/20/07 at 04:08 generated Percent Difference (%D's) of greater than the 25% limit for various compounds. These compounds include: Trichlorofluoromethane, Carbon Tetrachloride, 1,1,2-Trichloro-1,2,2-Trifluoroethane, Methylene Chloride, 4-Methyl-2-pentanone and o-Xylene. The %D's ranged from 26.6% to 38.7%. The associated samples were qualified as 'UJ', estimated, for the non-detectable results for the compounds noted. Positive results were not detected for the compounds affected.

Volatiles - Soil, Lab Project No. 7081622

Holding Times: Six samples were analyzed 1 day beyond the acceptable NYSDEC holding time of 7 days from Verified Time of Sample Receipt (VTSR). These samples include: CPC-SB-04 1.5-2.5, CPC-SB-06 1.5-2.5, CPC-SB-05 7-9, CPC-SB-02 5-7, CPC-SB-02 7-9 and CPC-DW-04-13-14. These samples were qualified as 'J', estimated, for the positive results and 'UJ', estimated, for the non-detectable results.

Continuing Calibration: The continuing calibrations analyzed on 8/20/07 at 04:08, 8/23/07 at 07:12 and 8/24/07 at 09:38 generated %D's greater than the 25% limit for various compounds. These compounds include: Trichlorofluoromethane, Carbon Tetrachloride, Carbon Disulfide, 1,1,2-Trichloro-1,2,2-Trifluoroethane, Methylene Chloride, 4-Methyl-2-pentanone, o-Xylene, trans-1,2-Dichloroethene, Methylcyclohexane, Acetone, 2-Butanone and Vinyl Acetate. The %D's ranged from 26.6% to 161.5%. The associated samples were qualified as 'J', estimated, for the positive results and 'UJ', estimated, for the non-detectable results for the compounds noted.

Method Blanks: Four method blank were analyzed for the associated soil samples. Volatile compounds found in the method blanks included Methylene Chloride and Acetone, as follows: The Method Blank analyzed on 8/23/07 was found to contain Methylene Chloride at 0.0044 mg/Kg and Methylene Chloride was also detected in the Method Blank on 8/24/07 at 0.0080 mg/Kg. A second Method Blank analyzed on 8/24/07 was found to contain Methylene Chloride at 0.0065 mg/Kg and Acetone at 0.014 mg/Kg. Limits of ten times the highest Methylene Chloride and Acetone results were used for review and qualification of the associated soil samples. Sample results for Methylene Chloride and Acetone found to be less than the respective method blank limit and reported at greater than the Contract Required Quantitation Limit (CRQL), were qualified as 'U', not detected.

It should be noted that the Laboratory re-evaluated and re-issued Method Blank results analyzed on 8/24/07 at 12:44 to include Acetone at 0.014 mg/Kg. This subsequent review and update was due to a request by the Client during evaluation of Acetone results detected for the project. All Acetone results detected in the project samples were qualified as 'U', not detected, through the associated Method Blank analyzed on 8/24/07.

Semi-Volatiles – Water and Soil, Lab Project Nos. 7081532 and 7081622

Qualification of the data sets for Semi-Volatiles was not required. The associated quality control information was found to be generated within acceptable limits.

Pesticides - Water, Lab Project No. 7081532

Continuing Calibration: The continuing calibration analyzed on 8/17/07 at 13:01 generated %D's greater than the 15% limit for various compounds. These compounds include: delta-BHC, p,p-DDE, Endrin Aldehyde, Endrin Ketone, p,p-DDD, Endosulfan Sulfate, Methoxychlor and Endrin Ketone. The %D's ranged from 17.8% to 31.4%. The associated samples were qualified as 'UJ', estimated, for the non-detectable results for the compounds noted. Positive results were not detected for the compounds affected.

Pesticides - Soil, Lab Project No. 7081622

Continuing Calibration: The continuing calibrations analyzed on 8/20/07 at 05:41, 8/20/07 at 08:13, 8/20/07 at 11:22, 8/21/07 at 05:27 and 8/21/07 at 10:27 generated %D's greater than the 15% limit for various compounds. These compounds include: delta-BHC, Endrin Ketone, p,p-DDD, Endosulfan Sulfate, Methoxychlor, Endosulfan I, gamma-BHC, Heptachlor, delta-BHC, p,p-DDT and alpha-BHC. The %D's ranged from 16.6% to 50.2%. The associated samples were qualified as 'J', estimated, for the positive results and 'UJ', estimated, for the non-detectable results for the compounds noted.

PCB's – Water and Soil, Lab Project Nos. 7081532 and 7081622

Qualification of the data sets for PCB's was not required. The associated quality control information was found to be generated within acceptable limits.

Inorganics - Water, Lab Project No. 7081532

MS/MSD: The site-specific MS and MSD for CPC-MW10 generated low recovery for Manganese at 59% and 60%, respectively (Limit 75-125). Sample CPC-MW10 was qualified as 'J', estimated, for the positive result for Manganese. Additional qualification was not required.

Inorganics and Cyanide - Soil, Lab Project No. 7081622

MS/MSD: The site-specific MS and MSD for CPC-SB03 1.5-3.5 generated low and high recovery for Chromium at 158% and 64%, respectively and 132% and 128% for Manganese (Limit 75-125). Sample CPC-SB03 1.5-3.5was qualified as 'J', estimated, for the positive results for Chromium and Manganese. Additional qualification was not required.

Please contact me by telephone or Fax at 301-294-6144, should you require additional information or clarification regarding this Letter Report.

Sincerely,

Andrea P. Schuessler

Andrea P. Schuessler, CHMM
ChemWorld Environmental, Inc.

c: EA-2007.2 file

ORGANIC DATA QUALIFIERS

U - Indicates that the compound was analyzed for, but not detected at or above the Contract Required Quantitation Limit (CRQL), or the compound is not detected due to qualification through the method or field blank.

J - The associated numerical value is an estimated quantity.

JN - Tentatively identified with approximated concentrations (Volatile and Semi-Volatile Organics). Presumptively present at an approximated quantity (Pesticides/PCBs).

UJ - The compound was analyzed for, but not detected. The sample quantitation limit is an estimated quantity due to variance from quality control limits.

C - Applies to Pesticide results where the identification has been confirmed by GC/MS.

E - Reported value is estimated due to quantitation above the calibration range.

D - Reported result taken from diluted sample analysis.

A - Aldol condensation product.

R - Reported value is unusable and rejected due to variance from quality control limits.

NA - Not Analyzed.

INORGANIC DATA QUALIFIERS

U - Indicates analyte not detected at or above the Contract Required Detection Limit (CRDL), or the compound is not detected due to qualification through the method or field blank.

B - Indicates analyte result is between Instrument Detection Limit (IDL) and CRDL.

J - The reported value is estimated due to variance from quality control limits.

UJ - The element was analyzed for, but not detected. The sample quantitation limit is an estimate due to variance from quality control limits.

E - Reported value is estimated because of the presence of interference.

R - Reported value is unusable and rejected due to variance from quality control limits.

NA - Not analyzed.

ATTACHMENT A

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-001
Client Id: CPC-MW01
Data File: 8M17915.D
Analysis Date: 08/20/07 22:57
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	156-59-2	cis-1,2-Dichloroethene	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	U J
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	127-18-4	Tetrachloroethene	5.0	U
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	79-01-6	Trichloroethene	5.0	U
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-002
Client Id: CPC-MW08
Data File: 8M17964.D
Analysis Date: 08/22/07 16:50
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	**156-59-2**	**cis-1,2-Dichloroethene**	**5.0**	**14**
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	U J
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	**127-18-4**	**Tetrachloroethene**	**5.0**	**1.1 J**
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	**79-01-6**	**Trichloroethene**	**5.0**	**15**
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 30.1

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-003
Client Id: CPC-MW03
Data File: 8M17917.D
Analysis Date: 08/20/07 23:48
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	**156-59-2**	**cis-1,2-Dichloroethene**	**5.0**	**1.4 J**
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	U J
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	**127-18-4**	**Tetrachloroethene**	**5.0**	**4.4 J**
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	**79-01-6**	**Trichloroethene**	**5.0**	**6.6**
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 12.4

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-004(10X)
Client Id: CPC-MW04
Data File: 8M17967.D
Analysis Date: 08/22/07 18:08
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 10
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	50	U	75-00-3	Chloroethane	50	U
79-34-5	1,1,2,2-Tetrachloroethane	50	U	67-66-3	Chloroform	50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	50	U J	74-87-3	Chloromethane	50	U
79-00-5	1,1,2-Trichloroethane	50	U	**156-59-2**	**cis-1,2-Dichloroethene**	**50**	**320**
75-34-3	1,1-Dichloroethane	50	U	10061-01-5	cis-1,3-Dichloropropene	50	U
75-35-4	1,1-Dichloroethene	50	U	110-82-7	Cyclohexane	50	U
120-82-1	1,2,4-Trichlorobenzene	50	U	124-48-1	Dibromochloromethane	50	U
96-12-8	1,2-Dibromo-3-Chloropropa	50	U	75-71-8	Dichlorodifluoromethane	50	U
106-93-4	1,2-Dibromoethane	50	U	100-41-4	Ethylbenzene	10	U
95-50-1	1,2-Dichlorobenzene	50	U	98-82-8	Isopropylbenzene	10	U
107-06-2	1,2-Dichloroethane	50	U	1330-20-7	m&p-Xylenes	20	U
78-87-5	1,2-Dichloropropane	50	U	79-20-9	Methyl Acetate	50	U
541-73-1	1,3-Dichlorobenzene	50	U	108-87-2	Methylcyclohexane	50	U
106-46-7	1,4-Dichlorobenzene	50	U	75-09-2	Methylene Chloride	50	U J
78-93-3	2-Butanone	50	U	1634-04-4	Methyl-t-butyl ether	10	U
591-78-6	2-Hexanone	50	U	95-47-6	o-Xylene	10	U J
108-10-1	4-Methyl-2-Pentanone	50	U J	100-42-5	Styrene	50	U
67-64-1	Acetone	250	U	**127-18-4**	**Tetrachloroethene**	**50**	**14 J**
71-43-2	Benzene	10	U	108-88-3	Toluene	10	U
75-27-4	Bromodichloromethane	50	U	156-60-5	trans-1,2-Dichloroethene	50	U
75-25-2	Bromoform	50	U	10061-02-6	trans-1,3-Dichloropropene	50	U
74-83-9	Bromomethane	50	U	**79-01-6**	**Trichloroethene**	**50**	**330**
75-15-0	Carbon Disulfide	50	U	75-69-4	Trichlorofluoromethane	50	U J
56-23-5	Carbon Tetrachloride	50	U J	75-01-4	Vinyl Chloride	50	U
108-90-7	Chlorobenzene	50	U				

Worksheet #: 57389

Total Target Concentration 664

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-005
Client Id: CPC-MW05
Data File: 8M17919.D
Analysis Date: 08/21/07 00:38
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	UJ	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	156-59-2	cis-1,2-Dichloroethene	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	UJ
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	UJ
108-10-1	4-Methyl-2-Pentanone	5.0	UJ	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	127-18-4	Tetrachloroethene	5.0	U
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	79-01-6	Trichloroethene	5.0	U
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	UJ
56-23-5	Carbon Tetrachloride	5.0	UJ	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration use

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-006
Client Id: CPC-MW06
Data File: 8M17920.D
Analysis Date: 08/21/07 01:03
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	**156-59-2**	**cis-1,2-Dichloroethene**	**5.0**	**4.3 J**
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	U J
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	**127-18-4**	**Tetrachloroethene**	**5.0**	**1.4 J**
71-43-2	Benzene	1.0	U	**108-88-3**	**Toluene**	**1.0**	**1.5 J**
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	**79-01-6**	**Trichloroethene**	**5.0**	**6.7 J**
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 13.9

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-006
Client Id: CPC-MW06
Data File: 8M17965.D
Analysis Date: 08/22/07 17:15
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-triflüor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	**156-59-2**	**cis-1,2-Dichloroethene**	**5.0**	**2.5 J**
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	U J
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	**127-18-4**	**Tetrachloroethene**	**5.0**	**1.1 J**
71-43-2	Benzene	1.0	U	**108-88-3**	**Toluene**	**1.0**	**1.2 J**
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	**79-01-6**	**Trichloroethene**	**5.0**	**4.2 J**
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-007
Client Id: CPC-MW07
Data File: 8M17966.D
Analysis Date: 08/22/07 17:40
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	156-59-2	cis-1,2-Dichloroethene	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	**1,4-Dichlorobenzene**	**5.0**	**1.2 J**	75-09-2	Methylene Chloride	5.0	U J
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	127-18-4	Tetrachloroethene	5.0	U
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	79-01-6	Trichloroethene	5.0	U
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 1.2

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-008
Client Id: CPC-MW09
Data File: 8M17925.D
Analysis Date: 08/21/07 03:09
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	156-59-2	cis-1,2-Dichloroethene	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	U J
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	**127-18-4**	**Tetrachloroethene**	**5.0**	**1.7 J**
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	**79-01-6**	**Trichloroethene**	**5.0**	**6.0**
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 7.7

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-009
Client Id: CPC-MW10
Data File: 8M17922.D
Analysis Date: 08/21/07 01:54
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	UJ	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	**156-59-2**	**cis-1,2-Dichloroethene**	**5.0**	**1.2 J**
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	UJ
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	UJ
108-10-1	4-Methyl-2-Pentanone	5.0	UJ	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	**127-18-4**	**Tetrachloroethene**	**5.0**	**1.1 J**
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	**79-01-6**	**Trichloroethene**	**5.0**	**4.1 J**
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	UJ
56-23-5	Carbon Tetrachloride	5.0	UJ	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 6.4

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-012
Client Id: CPC-DUPLICATE
Data File: 8M17926.D
Analysis Date: 08/21/07 03:34
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	**156-59-2**	**cis-1,2-Dichloroethene**	**5.0**	**20**
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	U J
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	**127-18-4**	**Tetrachloroethene**	**5.0**	**1.8 J**
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	**79-01-6**	**Trichloroethene**	**5.0**	**19**
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 40.8

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32351-013
Client Id: TB
Data File: 8M17903.D
Analysis Date: 08/20/07 17:55
Date Rec/Extracted: 08/15/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	156-59-2	cis-1,2-Dichloroethene	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	**75-09-2**	**Methylene Chloride**	**5.0**	**2.4 J**
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	127-18-4	Tetrachloroethene	5.0	U
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	79-01-6	Trichloroethene	5.0	U
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57389

Total Target Concentration 2.4

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-001
Client Id: CPC-DW01-10-11
Data File: 1M24248.D
Analysis Date: 08/23/07 09:05
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 86

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0058	U	56-23-5	Carbon Tetrachloride	0.0058	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0058	U	108-90-7	Chlorobenzene	0.0058	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0058	U	75-00-3	Chloroethane	0.0058	U
79-00-5	1,1,2-Trichloroethane	0.0058	U	67-66-3	Chloroform	0.0058	U
75-34-3	1,1-Dichloroethane	0.0058	U	74-87-3	Chloromethane	0.0058	U
75-35-4	1,1-Dichloroethene	0.0058	U	156-59-2	cis-1,2-Dichloroethene	0.0058	U
10061-01-5	cis-1,3-Dichloropropene	0.0058	U	120-82-1	1,2,4-Trichlorobenzene	0.0058	U
110-82-7	Cyclohexane	0.0058	U	124-48-1	Dibromochloromethane	0.0058	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0058	U	75-71-8	Dichlorodifluoromethane	0.0058	U
106-93-4	1,2-Dibromoethane	0.0058	U	**100-41-4**	**Ethylbenzene**	**0.0012**	**0.0065**
95-50-1	1,2-Dichlorobenzene	0.0058	U	98-82-8	Isopropylbenzene	0.0012	U
107-06-2	1,2-Dichloroethane	0.0058	U	**1330-20-7**	**m&p-Xylenes**	**0.0023**	**0.027**
78-87-5	1,2-Dichloropropane	0.0058	U	79-20-9	Methyl Acetate	0.0058	U
108-87-2	Methylcyclohexane	0.0058	U J	541-73-1	1,3-Dichlorobenzene	0.0058	U
75-09-2	**Methylene Chloride**	**0.0058**	U 0.010 B J	1634-04-4	Methyl-t-butyl ether	0.0012	U
106-46-7	1,4-Dichlorobenzene	0.0058	U	**95-47-6**	**o-Xylene**	**0.0012**	**0.0080**
78-93-3	2-Butanone	0.0058	U	100-42-5	Styrene	0.0058	U
591-78-6	2-Hexanone	0.0058	U	127-18-4	Tetrachloroethene	0.0012	U
108-10-1	4-Methyl-2-Pentanone	0.0058	U	**108-88-3**	**Toluene**	**0.0012**	**0.029**
67-64-1	Acetone	0.029	U	156-60-5	trans-1,2-Dichloroethene	0.0058	U J
71-43-2	**Benzene**	**0.0012**	**0.0032**	10061-02-6	trans-1,3-Dichloropropene	0.0058	U
75-27-4	Bromodichloromethane	0.0058	U	79-01-6	Trichloroethene	0.0058	U
75-25-2	Bromoform	0.0058	U	75-69-4	Trichlorofluoromethane	0.0058	U J
74-83-9	Bromomethane	0.0058	U	75-01-4	Vinyl Chloride	0.0058	U
75-15-0	Carbon Disulfide	0.0058	U J				

Worksheet #: 57391

Total Target Concentration 0.0837

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration use

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-002
Client Id: CPC-DW02-10-11
Data File: 1M24249.D
Analysis Date: 08/23/07 09:31
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 83

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0060	U	56-23-5	Carbon Tetrachloride	0.0060	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0060	U	108-90-7	Chlorobenzene	0.0060	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0060	U	75-00-3	Chloroethane	0.0060	U
79-00-5	1,1,2-Trichloroethane	0.0060	U	67-66-3	Chloroform	0.0060	U
75-34-3	1,1-Dichloroethane	0.0060	U	74-87-3	Chloromethane	0.0060	U
75-35-4	1,1-Dichloroethene	0.0060	U	156-59-2	cis-1,2-Dichloroethene	0.0060	U
10061-01-5	cis-1,3-Dichloropropene	0.0060	U	120-82-1	1,2,4-Trichlorobenzene	0.0060	U
110-82-7	Cyclohexane	0.0060	U	124-48-1	Dibromochloromethane	0.0060	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0060	U	75-71-8	Dichlorodifluoromethane	0.0060	U
106-93-4	1,2-Dibromoethane	0.0060	U	100-41-4	Ethylbenzene	0.0012	U
95-50-1	1,2-Dichlorobenzene	0.0060	U	98-82-8	Isopropylbenzene	0.0012	U
107-06-2	1,2-Dichloroethane	0.0060	U	1330-20-7	m&p-Xylenes	0.0024	U
78-87-5	1,2-Dichloropropane	0.0060	U	79-20-9	Methyl Acetate	0.0060	U
108-87-2	Methylcyclohexane	0.0060	U J	541-73-1	1,3-Dichlorobenzene	0.0060	U
75-09-2	**Methylene Chloride**	**0.0060**	U 0.0096 B J	1634-04-4	Methyl-t-butyl ether	0.0012	U
106-46-7	1,4-Dichlorobenzene	0.0060	U	95-47-6	o-Xylene	0.0012	U
78-93-3	2-Butanone	0.0060	U	100-42-5	Styrene	0.0060	U
591-78-6	2-Hexanone	0.0060	U	127-18-4	Tetrachloroethene	0.0012	U
108-10-1	4-Methyl-2-Pentanone	0.0060	U	**108-88-3**	**Toluene**	**0.0012**	**0.0021**
67-64-1	**Acetone**	**0.030**	0.039 J	156-60-5	trans-1,2-Dichloroethene	0.0060	U J
71-43-2	Benzene	0.0012	U	10061-02-6	trans-1,3-Dichloropropene	0.0060	U
75-27-4	Bromodichloromethane	0.0060	U	79-01-6	Trichloroethene	0.0060	U
75-25-2	Bromoform	0.0060	U	75-69-4	Trichlorofluoromethane	0.0060	U J
74-83-9	Bromomethane	0.0060	U	75-01-4	Vinyl Chloride	0.0060	U
75-15-0	**Carbon Disulfide**	**0.0060**	**0.0028 J**				

Worksheet #: 57391

Total Target Concentration 0.0535

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-003
Client Id: CPC-DW03-10-11
Data File: 1M24250.D
Analysis Date: 08/23/07 09:56
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 90

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0056	U	56-23-5	Carbon Tetrachloride	0.0056	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0056	U	108-90-7	Chlorobenzene	0.0056	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0056	U	75-00-3	Chloroethane	0.0056	U
79-00-5	1,1,2-Trichloroethane	0.0056	U	67-66-3	Chloroform	0.0056	U
75-34-3	1,1-Dichloroethane	0.0056	U	74-87-3	Chloromethane	0.0056	U
75-35-4	1,1-Dichloroethene	0.0056	U	156-59-2	cis-1,2-Dichloroethene	0.0056	U
10061-01-5	cis-1,3-Dichloropropene	0.0056	U	120-82-1	1,2,4-Trichlorobenzene	0.0056	U
110-82-7	Cyclohexane	0.0056	U	124-48-1	Dibromochloromethane	0.0056	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0056	U	75-71-8	Dichlorodifluoromethane	0.0056	U
106-93-4	1,2-Dibromoethane	0.0056	U	100-41-4	Ethylbenzene	0.0011	U
95-50-1	1,2-Dichlorobenzene	0.0056	U	98-82-8	Isopropylbenzene	0.0011	U
107-06-2	1,2-Dichloroethane	0.0056	U	1330-20-7	m&p-Xylenes	0.0022	U
78-87-5	1,2-Dichloropropane	0.0056	U	79-20-9	Methyl Acetate	0.0056	U
108-87-2	Methylcyclohexane	0.0056	U J	541-73-1	1,3-Dichlorobenzene	0.0056	U
75-09-2	**Methylene Chloride**	**0.0056**	U 0.015 B J	1634-04-4	Methyl-t-butyl ether	0.0011	U
106-46-7	1,4-Dichlorobenzene	0.0056	U	95-47-6	o-Xylene	0.0011	U
78-93-3	2-Butanone	0.0056	U	100-42-5	Styrene	0.0056	U
591-78-6	2-Hexanone	0.0056	U	127-18-4	Tetrachloroethene	0.0011	U
108-10-1	4-Methyl-2-Pentanone	0.0056	U	**108-88-3**	**Toluene**	**0.0011**	**0.0016**
67-64-1	Acetone	0.028	U	156-60-5	trans-1,2-Dichloroethene	0.0056	U J
71-43-2	Benzene	0.0011	U	10061-02-6	trans-1,3-Dichloropropene	0.0056	U
75-27-4	Bromodichloromethane	0.0056	U	79-01-6	Trichloroethene	0.0056	U
75-25-2	Bromoform	0.0056	U	75-69-4	Trichlorofluoromethane	0.0056	U J
74-83-9	Bromomethane	0.0056	U	75-01-4	Vinyl Chloride	0.0056	U
75-15-0	Carbon Disulfide	0.0056	U J				

Worksheet #: 57391

Total Target Concentration 0.0166

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-004
Client Id: CPC-DW04-8-9
Data File: 1M24251.D
Analysis Date: 08/23/07 10:22
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 85

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0059	U	56-23-5	Carbon Tetrachloride	0.0059	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0059	U	108-90-7	Chlorobenzene	0.0059	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0059	U	75-00-3	Chloroethane	0.0059	U
79-00-5	1,1,2-Trichloroethane	0.0059	U	67-66-3	Chloroform	0.0059	U
75-34-3	1,1-Dichloroethane	0.0059	U	74-87-3	Chloromethane	0.0059	U
75-35-4	1,1-Dichloroethene	0.0059	U	156-59-2	cis-1,2-Dichloroethene	0.0059	U
10061-01-5	cis-1,3-Dichloropropene	0.0059	U	110-82-7	Cyclohexane	0.0059	U
124-48-1	Dibromochloromethane	0.0059	U	96-12-8	1,2-Dibromo-3-Chloropropa	0.0059	U
75-71-8	Dichlorodifluoromethane	0.0059	U	106-93-4	1,2-Dibromoethane	0.0059	U
100-41-4	**Ethylbenzene**	**0.0012**	**0.0012**	98-82-8	Isopropylbenzene	0.0012	U
107-06-2	1,2-Dichloroethane	0.0059	U	**1330-20-7**	**m&p-Xylenes**	**0.0024**	**0.0032**
78-87-5	1,2-Dichloropropane	0.0059	U	79-20-9	Methyl Acetate	0.0059	U
108-87-2	Methylcyclohexane	0.0059	U J	**75-09-2**	**Methylene Chloride**	**0.0059**	U 0.0060 B J
1634-04-4	Methyl-t-butyl ether	0.0012	U	**95-47-6**	**o-Xylene**	**0.0012**	**0.0033**
78-93-3	2-Butanone	0.0059	U	100-42-5	Styrene	0.0059	U
591-78-6	2-Hexanone	0.0059	U	127-18-4	Tetrachloroethene	0.0012	U
108-10-1	4-Methyl-2-Pentanone	0.0059	U	108-88-3	Toluene	0.0012	U
67-64-1	Acetone	0.029	U	156-60-5	trans-1,2-Dichloroethene	0.0059	U J
71-43-2	Benzene	0.0012	U	10061-02-6	trans-1,3-Dichloropropene	0.0059	U
75-27-4	Bromodichloromethane	0.0059	U	79-01-6	Trichloroethene	0.0059	U
75-25-2	Bromoform	0.0059	U	75-69-4	Trichlorofluoromethane	0.0059	U J
74-83-9	Bromomethane	0.0059	U	75-01-4	Vinyl Chloride	0.0059	U
75-15-0	Carbon Disulfide	0.0059	U J				

Worksheet #: 57391

Total Target Concentration 0.0137

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-005
Client Id: CPC-SOIL-Duplicate
Data File: 1M24259.D
Analysis Date: 08/23/07 13:53
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 92

Units: mg/Kg

Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0054	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0054	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0054	U
79-00-5	1,1,2-Trichloroethane	0.0054	U
75-34-3	1,1-Dichloroethane	0.0054	U
75-35-4	1,1-Dichloroethene	0.0054	U
10061-01-5	cis-1,3-Dichloropropene	0.0054	U
110-82-7	Cyclohexane	0.0054	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0054	U
106-93-4	1,2-Dibromoethane	0.0054	U
95-50-1	1,2-Dichlorobenzene	0.0054	U
107-06-2	1,2-Dichloroethane	0.0054	U
78-87-5	1,2-Dichloropropane	0.0054	U
108-87-2	Methylcyclohexane	0.0054	U J
75-09-2	**Methylene Chloride**	**0.0054**	U 0.0089 B J
106-46-7	1,4-Dichlorobenzene	0.0054	U
78-93-3	2-Butanone	0.0054	U
591-78-6	2-Hexanone	0.0054	U
108-10-1	4-Methyl-2-Pentanone	0.0054	U
67-64-1	Acetone	0.027	U
71-43-2	Benzene	0.0011	U
75-27-4	Bromodichloromethane	0.0054	U
75-25-2	Bromoform	0.0054	U
74-83-9	Bromomethane	0.0054	U
75-15-0	Carbon Disulfide	0.0054	U J
56-23-5	Carbon Tetrachloride	0.0054	U
108-90-7	Chlorobenzene	0.0054	U
75-00-3	Chloroethane	0.0054	U
67-66-3	Chloroform	0.0054	U
74-87-3	Chloromethane	0.0054	U
156-59-2	**cis-1,2-Dichloroethene**	**0.0054**	**0.0017 J**
120-82-1	1,2,4-Trichlorobenzene	0.0054	U
124-48-1	Dibromochloromethane	0.0054	U
75-71-8	Dichlorodifluoromethane	0.0054	U
100-41-4	Ethylbenzene	0.0011	U
98-82-8	Isopropylbenzene	0.0011	U
1330-20-7	**m&p-Xylenes**	**0.0022**	**0.0015 J**
79-20-9	Methyl Acetate	0.0054	U
541-73-1	1,3-Dichlorobenzene	0.0054	U
1634-04-4	Methyl-t-butyl ether	0.0011	U
95-47-6	o-Xylene	0.0011	U
100-42-5	Styrene	0.0054	U
127-18-4	**Tetrachloroethene**	**0.0011**	**0.0049**
108-88-3	**Toluene**	**0.0011**	**0.0038**
156-60-5	trans-1,2-Dichloroethene	0.0054	U J
10061-02-6	trans-1,3-Dichloropropene	0.0054	U
79-01-6	**Trichloroethene**	**0.0054**	**0.015**
75-69-4	Trichlorofluoromethane	0.0054	U J
75-01-4	Vinyl Chloride	0.0054	U

Worksheet #: 57391

Total Target Concentration 0.0358

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-006
Client Id: CPC-SB01-1.5-2.5
Data File: 1M24260.D
Analysis Date: 08/23/07 14:19
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 95

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0053	U	56-23-5	Carbon Tetrachloride	0.0053	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0053	U	108-90-7	Chlorobenzene	0.0053	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0053	U	75-00-3	Chloroethane	0.0053	U
79-00-5	1,1,2-Trichloroethane	0.0053	U	67-66-3	Chloroform	0.0053	U
75-34-3	1,1-Dichloroethane	0.0053	U	74-87-3	Chloromethane	0.0053	U
75-35-4	1,1-Dichloroethene	0.0053	U	**156-59-2**	**cis-1,2-Dichloroethene**	**0.0053**	**0.0014 J**
10061-01-5	cis-1,3-Dichloropropene	0.0053	U	120-82-1	1,2,4-Trichlorobenzene	0.0053	U
110-82-7	Cyclohexane	0.0053	U	124-48-1	Dibromochloromethane	0.0053	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0053	U	75-71-8	Dichlorodifluoromethane	0.0053	U
106-93-4	1,2-Dibromoethane	0.0053	U	100-41-4	Ethylbenzene	0.0011	U
95-50-1	1,2-Dichlorobenzene	0.0053	U	98-82-8	Isopropylbenzene	0.0011	U
107-06-2	1,2-Dichloroethane	0.0053	U	1330-20-7	m&p-Xylenes	0.0021	U
78-87-5	1,2-Dichloropropane	0.0053	U	79-20-9	Methyl Acetate	0.0053	U
108-87-2	Methylcyclohexane	0.0053	U J	541-73-1	1,3-Dichlorobenzene	0.0053	U
75-09-2	**Methylene Chloride**	**0.0053**	U 0.0080 B J	1634-04-4	Methyl-t-butyl ether	0.0011	U
106-46-7	1,4-Dichlorobenzene	0.0053	U	95-47-6	o-Xylene	0.0011	U
78-93-3	2-Butanone	0.0053	U	100-42-5	Styrene	0.0053	U
591-78-6	2-Hexanone	0.0053	U	**127-18-4**	**Tetrachloroethene**	**0.0011**	**0.0093**
108-10-1	4-Methyl-2-Pentanone	0.0053	U	**108-88-3**	**Toluene**	**0.0011**	**0.0028**
67-64-1	Acetone	0.026	U	156-60-5	trans-1,2-Dichloroethene	0.0053	U J
71-43-2	Benzene	0.0011	U	10061-02-6	trans-1,3-Dichloropropene	0.0053	U
75-27-4	Bromodichloromethane	0.0053	U	**79-01-6**	**Trichloroethene**	**0.0053**	**0.018**
75-25-2	Bromoform	0.0053	U	75-69-4	Trichlorofluoromethane	0.0053	U J
74-83-9	Bromomethane	0.0053	U	75-01-4	Vinyl Chloride	0.0053	U
75-15-0	Carbon Disulfide	0.0053	U J				

Worksheet #: 57391

Total Target Concentration 0.0395

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-007
Client Id: CPC-SB02-1-5
Data File: 1M24261.D
Analysis Date: 08/23/07 14:44
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 94

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0053	U	56-23-5	Carbon Tetrachloride	0.0053	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0053	U	108-90-7	Chlorobenzene	0.0053	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0053	U	75-00-3	Chloroethane	0.0053	U
79-00-5	1,1,2-Trichloroethane	0.0053	U	67-66-3	Chloroform	0.0053	U
75-34-3	1,1-Dichloroethane	0.0053	U	74-87-3	Chloromethane	0.0053	U
75-35-4	1,1-Dichloroethene	0.0053	U	156-59-2	cis-1,2-Dichloroethene	0.0053	U
10061-01-5	cis-1,3-Dichloropropene	0.0053	U	120-82-1	1,2,4-Trichlorobenzene	0.0053	U
110-82-7	Cyclohexane	0.0053	U	124-48-1	Dibromochloromethane	0.0053	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0053	U	75-71-8	Dichlorodifluoromethane	0.0053	U
106-93-4	1,2-Dibromoethane	0.0053	U	100-41-4	Ethylbenzene	0.0011	U
95-50-1	1,2-Dichlorobenzene	0.0053	U	98-82-8	Isopropylbenzene	0.0011	U
107-06-2	1,2-Dichloroethane	0.0053	U	1330-20-7	m&p-Xylenes	0.0021	U
78-87-5	1,2-Dichloropropane	0.0053	U	79-20-9	Methyl Acetate	0.0053	U
108-87-2	Methylcyclohexane	0.0053	U J	541-73-1	1,3-Dichlorobenzene	0.0053	U
75-09-2	**Methylene Chloride**	**0.0053**	U 0.0088 B J	1634-04-4	Methyl-t-butyl ether	0.0011	U
106-46-7	1,4-Dichlorobenzene	0.0053	U	95-47-6	o-Xylene	0.0011	U
78-93-3	2-Butanone	0.0053	U	100-42-5	Styrene	0.0053	U
591-78-6	2-Hexanone	0.0053	U	**127-18-4**	**Tetrachloroethene**	**0.0011**	**0.0039**
108-10-1	4-Methyl-2-Pentanone	0.0053	U	**108-88-3**	**Toluene**	**0.0011**	**0.0026**
67-64-1	Acetone	0.027	U	156-60-5	trans-1,2-Dichloroethene	0.0053	U J
71-43-2	Benzene	0.0011	U	10061-02-6	trans-1,3-Dichloropropene	0.0053	U
75-27-4	Bromodichloromethane	0.0053	U	**79-01-6**	**Trichloroethene**	**0.0053**	**0.010**
75-25-2	Bromoform	0.0053	U	75-69-4	Trichlorofluoromethane	0.0053	U J
74-83-9	Bromomethane	0.0053	U	75-01-4	Vinyl Chloride	0.0053	U
75-15-0	Carbon Disulfide	0.0053	U J				

Worksheet #: 57391

Total Target Concentration 0.0253

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-008
Client Id: CPC-SB03-1.5-3.5
Data File: 1M24264.D
Analysis Date: 08/23/07 16:02
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 93

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0054	U	56-23-5	Carbon Tetrachloride	0.0054	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0054	U	108-90-7	Chlorobenzene	0.0054	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0054	U	75-00-3	Chloroethane	0.0054	U
79-00-5	1,1,2-Trichloroethane	0.0054	U	67-66-3	Chloroform	0.0054	U
75-34-3	1,1-Dichloroethane	0.0054	U	74-87-3	Chloromethane	0.0054	U
75-35-4	1,1-Dichloroethene	0.0054	U	156-59-2	cis-1,2-Dichloroethene	0.0054	U
10061-01-5	cis-1,3-Dichloropropene	0.0054	U	110-82-7	Cyclohexane	0.0054	U
95-63-6	1,2,4-Trimethylbenzene	0.0011	U	124-48-1	Dibromochloromethane	0.0054	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0054	U	75-71-8	Dichlorodifluoromethane	0.0054	U
106-93-4	1,2-Dibromoethane	0.0054	U	100-41-4	Ethylbenzene	0.0011	U
95-50-1	1,2-Dichlorobenzene	0.0054	U	98-82-8	Isopropylbenzene	0.0011	U
107-06-2	1,2-Dichloroethane	0.0054	U	**1330-20-7**	**m&p-Xylenes**	**0.0022**	**0.0020 J**
78-87-5	1,2-Dichloropropane	0.0054	U	79-20-9	Methyl Acetate	0.0054	U
108-87-2	Methylcyclohexane	0.0054	U J	541-73-1	1,3-Dichlorobenzene	0.0054	U
75-09-2	**Methylene Chloride**	**0.0054**	U 0.0072 B J	1634-04-4	Methyl-t-butyl ether	0.0011	U
106-46-7	1,4-Dichlorobenzene	0.0054	U	95-47-6	o-Xylene	0.0011	U
78-93-3	2-Butanone	0.0054	U	100-42-5	Styrene	0.0054	U
591-78-6	2-Hexanone	0.0054	U	**127-18-4**	**Tetrachloroethene**	**0.0011**	**0.0030**
108-10-1	4-Methyl-2-Pentanone	0.0054	U	108-88-3	Toluene	0.0011	U
67-64-1	Acetone	0.027	U	156-60-5	trans-1,2-Dichloroethene	0.0054	U J
71-43-2	Benzene	0.0011	U	10061-02-6	trans-1,3-Dichloropropene	0.0054	U
75-27-4	Bromodichloromethane	0.0054	U	79-01-6	Trichloroethene	0.0054	U
75-25-2	Bromoform	0.0054	U	75-69-4	Trichlorofluoromethane	0.0054	U J
74-83-9	Bromomethane	0.0054	U	75-01-4	Vinyl Chloride	0.0054	U
75-15-0	Carbon Disulfide	0.0054	U J				

Worksheet #: 57391

Total Target Concentration 0.0122

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-011
Client Id: CPC-SB04-1.5-2.5
Data File: 1M24285.D
Analysis Date: 08/24/07 11:06
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 90

Units: mg/Kg

Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0056	U J
79-34-5	1,1,2,2-Tetrachloroethane	0.0056	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0056	U
79-00-5	1,1,2-Trichloroethane	0.0056	U
75-34-3	1,1-Dichloroethane	0.0056	U
75-35-4	1,1-Dichloroethene	0.0056	U
120-82-1	1,2,4-Trichlorobenzene	0.0056	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0056	U
106-93-4	1,2-Dibromoethane	0.0056	U
95-50-1	1,2-Dichlorobenzene	0.0056	U
107-06-2	1,2-Dichloroethane	0.0056	U
78-87-5	1,2-Dichloropropane	0.0056	U
541-73-1	1,3-Dichlorobenzene	0.0056	U
106-46-7	1,4-Dichlorobenzene	0.0056	U
78-93-3	2-Butanone	0.0056	U
591-78-6	2-Hexanone	0.0056	U
108-10-1	4-Methyl-2-Pentanone	0.0056	U
67-64-1	Acetone	0.028	U
71-43-2	Benzene	0.0011	U
75-27-4	Bromodichloromethane	0.0056	U
75-25-2	Bromoform	0.0056	U
74-83-9	Bromomethane	0.0056	U
75-15-0	Carbon Disulfide	0.0056	U
56-23-5	Carbon Tetrachloride	0.0056	U
108-90-7	Chlorobenzene	0.0056	U
75-00-3	Chloroethane	0.0056	U J
67-66-3	Chloroform	0.0056	U
74-87-3	Chloromethane	0.0056	U
156-59-2	cis-1,2-Dichloroethene	0.0056	U
10061-01-5	cis-1,3-Dichloropropene	0.0056	U
110-82-7	Cyclohexane	0.0056	U
124-48-1	Dibromochloromethane	0.0056	U
75-71-8	Dichlorodifluoromethane	0.0056	U
100-41-4	Ethylbenzene	0.0011	U
98-82-8	Isopropylbenzene	0.0011	U
1330-20-7	m&p-Xylenes	0.0022	U
79-20-9	Methyl Acetate	0.0056	U
108-87-2	Methylcyclohexane	0.0056	U
75-09-2	**Methylene Chloride**	**0.0056**	U 0.0068 B J
1634-04-4	Methyl-t-butyl ether	0.0011	U J
95-47-6	o-Xylene	0.0011	U
100-42-5	Styrene	0.0056	U
127-18-4	Tetrachloroethene	0.0011	U
108-88-3	Toluene	0.0011	U
156-60-5	trans-1,2-Dichloroethene	0.0056	U
10061-02-6	trans-1,3-Dichloropropene	0.0056	U
79-01-6	Trichloroethene	0.0056	U
75-69-4	Trichlorofluoromethane	0.0056	U
75-01-4	Vinyl Chloride	0.0056	U

Worksheet #: 57392

Total Target Concentration 0.0068

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration use

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-012(5X)
Client Id: CPC-SB06-1.5-2.5
Data File: 2M23126.D
Analysis Date: 08/24/07 20:53
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 5
Solids: 93

Method: EPA 8260B

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.027	U	75-00-3	Chloroethane	0.027	U
79-34-5	1,1,2,2-Tetrachloroethane	0.027	U	67-66-3	Chloroform	0.027	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.027	U	74-87-3	Chloromethane	0.027	U
79-00-5	1,1,2-Trichloroethane	0.027	U	156-59-2	cis-1,2-Dichloroethene	0.027	U
75-34-3	1,1-Dichloroethane	0.027	U	10061-01-5	cis-1,3-Dichloropropene	0.027	U
75-35-4	1,1-Dichloroethene	0.027	U	110-82-7	Cyclohexane	0.027	U
120-82-1	1,2,4-Trichlorobenzene	0.027	U	124-48-1	Dibromochloromethane	0.027	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.027	U	75-71-8	Dichlorodifluoromethane	0.027	U
106-93-4	1,2-Dibromoethane	0.027	U	100-41-4	Ethylbenzene	0.0054	0.0069
95-50-1	1,2-Dichlorobenzene	0.027	U	98-82-8	Isopropylbenzene	0.0054	U
107-06-2	1,2-Dichloroethane	0.027	U	1330-20-7	m&p-Xylenes	0.011	0.021
78-87-5	1,2-Dichloropropane	0.027	U	79-20-9	Methyl Acetate	0.027	U
541-73-1	1,3-Dichlorobenzene	0.027	U	108-87-2	Methylcyclohexane	0.027	U
106-46-7	1,4-Dichlorobenzene	0.027	U	75-09-2	Methylene Chloride	0.027 U	~~0.027 B~~
78-93-3	2-Butanone	0.027	U	1634-04-4	Methyl-t-butyl ether	0.0054	U
591-78-6	2-Hexanone	0.027	U	95-47-6	o-Xylene	0.0054	U
108-10-1	4-Methyl-2-Pentanone	0.027	U	100-42-5	Styrene	0.027	U
67-64-1	Acetone	0.13 U	~~0.091 JB~~	127-18-4	Tetrachloroethene	0.027	0.33
71-43-2	Benzene	0.0054	U	108-88-3	Toluene	0.0054	U
75-27-4	Bromodichloromethane	0.027	U	156-60-5	trans-1,2-Dichloroethene	0.027	U
75-25-2	Bromoform	0.027	U	10061-02-6	trans-1,3-Dichloropropene	0.027	U
74-83-9	Bromomethane	0.027	U	79-01-6	Trichloroethene	0.027	0.016 J
75-15-0	Carbon Disulfide	0.027	U	75-69-4	Trichlorofluoromethane	0.027	U
56-23-5	Carbon Tetrachloride	0.027	U	75-01-4	Vinyl Chloride	0.027	U
108-90-7	Chlorobenzene	0.027	U				

Worksheet #: 73611

Total Target Concentration 0.4919

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Reported %D/D>40% between columns due to coelution. Lower concentration used

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-013
Client Id: CPC-SB05-1-3
Data File: 1M24267.D
Analysis Date: 08/23/07 17:19
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 94

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0053	U	75-00-3	Chloroethane	0.0053	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0053	U	67-66-3	Chloroform	0.0053	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0053	U	74-87-3	Chloromethane	0.0053	U
79-00-5	1,1,2-Trichloroethane	0.0053	U	156-59-2	cis-1,2-Dichloroethene	0.0053	U
75-34-3	1,1-Dichloroethane	0.0053	U	10061-01-5	cis-1,3-Dichloropropene	0.0053	U
75-35-4	1,1-Dichloroethene	0.0053	U	110-82-7	Cyclohexane	0.0053	U
120-82-1	1,2,4-Trichlorobenzene	0.0053	U	124-48-1	Dibromochloromethane	0.0053	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0053	U	75-71-8	Dichlorodifluoromethane	0.0053	U
106-93-4	1,2-Dibromoethane	0.0053	U	100-41-4	Ethylbenzene	0.0011	U
95-50-1	1,2-Dichlorobenzene	0.0053	U	98-82-8	Isopropylbenzene	0.0011	U
107-06-2	1,2-Dichloroethane	0.0053	U	1330-20-7	m&p-Xylenes	0.0021	U
78-87-5	1,2-Dichloropropane	0.0053	U	79-20-9	Methyl Acetate	0.0053	U
541-73-1	1,3-Dichlorobenzene	0.0053	U	108-87-2	Methylcyclohexane	0.0053	U J
106-46-7	1,4-Dichlorobenzene	0.0053	U	**75-09-2**	**Methylene Chloride**	**0.0053**	U 0.0066 B J
78-93-3	2-Butanone	0.0053	U	1634-04-4	Methyl-t-butyl ether	0.0011	U
591-78-6	2-Hexanone	0.0053	U	95-47-6	o-Xylene	0.0011	U
108-10-1	4-Methyl-2-Pentanone	0.0053	U	100-42-5	Styrene	0.0053	U
67-64-1	Acetone	0.027	U	127-18-4	Tetrachloroethene	0.0011	U
71-43-2	Benzene	0.0011	U	108-88-3	Toluene	0.0011	U
75-27-4	Bromodichloromethane	0.0053	U	156-60-5	trans-1,2-Dichloroethene	0.0053	U J
75-25-2	Bromoform	0.0053	U	10061-02-6	trans-1,3-Dichloropropene	0.0053	U
74-83-9	Bromomethane	0.0053	U	79-01-6	Trichloroethene	0.0053	U
75-15-0	Carbon Disulfide	0.0053	U J	75-69-4	Trichlorofluoromethane	0.0053	U J
56-23-5	Carbon Tetrachloride	0.0053	U	75-01-4	Vinyl Chloride	0.0053	U
108-90-7	Chlorobenzene	0.0053	U				

Worksheet #: 57392

Total Target Concentration 0.0066

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-014
Client Id: CPC-SB05-3-5
Data File: 1M24268.D
Analysis Date: 08/23/07 17:45
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 96

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0052	U	75-00-3	Chloroethane	0.0052	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0052	U	67-66-3	Chloroform	0.0052	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0052	U	74-87-3	Chloromethane	0.0052	U
79-00-5	1,1,2-Trichloroethane	0.0052	U	156-59-2	cis-1,2-Dichloroethene	0.0052	U
75-34-3	1,1-Dichloroethane	0.0052	U	10061-01-5	cis-1,3-Dichloropropene	0.0052	U
75-35-4	1,1-Dichloroethene	0.0052	U	110-82-7	Cyclohexane	0.0052	U
120-82-1	1,2,4-Trichlorobenzene	0.0052	U	124-48-1	Dibromochloromethane	0.0052	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0052	U	75-71-8	Dichlorodifluoromethane	0.0052	U
106-93-4	1,2-Dibromoethane	0.0052	U	100-41-4	Ethylbenzene	0.0010	U
95-50-1	1,2-Dichlorobenzene	0.0052	U	98-82-8	Isopropylbenzene	0.0010	U
107-06-2	1,2-Dichloroethane	0.0052	U	1330-20-7	m&p-Xylenes	0.0021	U
78-87-5	1,2-Dichloropropane	0.0052	U	79-20-9	Methyl Acetate	0.0052	U
541-73-1	1,3-Dichlorobenzene	0.0052	U	108-87-2	Methylcyclohexane	0.0052	U J
106-46-7	1,4-Dichlorobenzene	0.0052	U	**75-09-2**	**Methylene Chloride**	**0.0052**	U **0.0069** B
78-93-3	2-Butanone	0.0052	U	1634-04-4	Methyl-t-butyl ether	0.0010	U
591-78-6	2-Hexanone	0.0052	U	95-47-6	o-Xylene	0.0010	U
108-10-1	4-Methyl-2-Pentanone	0.0052	U	100-42-5	Styrene	0.0052	U
67-64-1	Acetone	0.026	U	127-18-4	Tetrachloroethene	0.0010	U
71-43-2	Benzene	0.0010	U	**108-88-3**	**Toluene**	**0.0010**	**0.0012**
75-27-4	Bromodichloromethane	0.0052	U	156-60-5	trans-1,2-Dichloroethene	0.0052	U J
75-25-2	Bromoform	0.0052	U	10061-02-6	trans-1,3-Dichloropropene	0.0052	U
74-83-9	Bromomethane	0.0052	U	79-01-6	Trichloroethene	0.0052	U
75-15-0	Carbon Disulfide	0.0052	U J	75-69-4	Trichlorofluoromethane	0.0052	U J
56-23-5	Carbon Tetrachloride	0.0052	U	75-01-4	Vinyl Chloride	0.0052	U
108-90-7	Chlorobenzene	0.0052	U				

Worksheet #: 57392

Total Target Concentration 0.0081

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-015
Client Id: CPC-SB05-5-7
Data File: 1M24269.D
Analysis Date: 08/23/07 18:11
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 98

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0051	U	75-00-3	Chloroethane	0.0051	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0051	U	67-66-3	Chloroform	0.0051	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0051	U	74-87-3	Chloromethane	0.0051	U
79-00-5	1,1,2-Trichloroethane	0.0051	U	156-59-2	cis-1,2-Dichloroethene	0.0051	U
75-34-3	1,1-Dichloroethane	0.0051	U	10061-01-5	cis-1,3-Dichloropropene	0.0051	U
75-35-4	1,1-Dichloroethene	0.0051	U	110-82-7	Cyclohexane	0.0051	U
120-82-1	1,2,4-Trichlorobenzene	0.0051	U	124-48-1	Dibromochloromethane	0.0051	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0051	U	75-71-8	Dichlorodifluoromethane	0.0051	U
106-93-4	1,2-Dibromoethane	0.0051	U	100-41-4	Ethylbenzene	0.0010	U
95-50-1	1,2-Dichlorobenzene	0.0051	U	98-82-8	Isopropylbenzene	0.0010	U
107-06-2	1,2-Dichloroethane	0.0051	U	1330-20-7	m&p-Xylenes	0.0020	U
78-87-5	1,2-Dichloropropane	0.0051	U	79-20-9	Methyl Acetate	0.0051	U
541-73-1	1,3-Dichlorobenzene	0.0051	U	108-87-2	Methylcyclohexane	0.0051	U J
106-46-7	1,4-Dichlorobenzene	0.0051	U	**75-09-2**	**Methylene Chloride**	**0.0051**	U 0.0065 B J
78-93-3	2-Butanone	0.0051	U	1634-04-4	Methyl-t-butyl ether	0.0010	U
591-78-6	2-Hexanone	0.0051	U	95-47-6	o-Xylene	0.0010	U
108-10-1	4-Methyl-2-Pentanone	0.0051	U	100-42-5	Styrene	0.0051	U
67-64-1	Acetone	0.026	U	127-18-4	Tetrachloroethene	0.0010	U
71-43-2	Benzene	0.0010	U	108-88-3	Toluene	0.0010	U
75-27-4	Bromodichloromethane	0.0051	U	156-60-5	trans-1,2-Dichloroethene	0.0051	U J
75-25-2	Bromoform	0.0051	U	10061-02-6	trans-1,3-Dichloropropene	0.0051	U
74-83-9	Bromomethane	0.0051	U	79-01-6	Trichloroethene	0.0051	U
75-15-0	Carbon Disulfide	0.0051	U J	75-69-4	Trichlorofluoromethane	0.0051	U J
56-23-5	Carbon Tetrachloride	0.0051	U	75-01-4	Vinyl Chloride	0.0051	U
108-90-7	Chlorobenzene	0.0051	U				

Worksheet #: 57392

Total Target Concentration 0.0065

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-016
Client Id: CPC-SB05-7-9
Data File: 1M24297.D
Analysis Date: 08/24/07 16:17
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 98

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0051	U J	75-00-3	Chloroethane	0.0051	U J
79-34-5	1,1,2,2-Tetrachloroethane	0.0051	U	67-66-3	Chloroform	0.0051	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0051	U	74-87-3	Chloromethane	0.0051	U
79-00-5	1,1,2-Trichloroethane	0.0051	U	156-59-2	cis-1,2-Dichloroethene	0.0051	U
75-34-3	1,1-Dichloroethane	0.0051	U	10061-01-5	cis-1,3-Dichloropropene	0.0051	U
75-35-4	1,1-Dichloroethene	0.0051	U	110-82-7	Cyclohexane	0.0051	U
120-82-1	1,2,4-Trichlorobenzene	0.0051	U	124-48-1	Dibromochloromethane	0.0051	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0051	U	75-71-8	Dichlorodifluoromethane	0.0051	U
106-93-4	1,2-Dibromoethane	0.0051	U	100-41-4	Ethylbenzene	0.0010	U
95-50-1	1,2-Dichlorobenzene	0.0051	U	98-82-8	Isopropylbenzene	0.0010	U
107-06-2	1,2-Dichloroethane	0.0051	U	1330-20-7	m&p-Xylenes	0.0020	U
78-87-5	1,2-Dichloropropane	0.0051	U	79-20-9	Methyl Acetate	0.0051	U
541-73-1	1,3-Dichlorobenzene	0.0051	U	108-87-2	Methylcyclohexane	0.0051	U
106-46-7	1,4-Dichlorobenzene	0.0051	U	**75-09-2**	**Methylene Chloride**	**0.0051**	U 0.0066 B J
78-93-3	2-Butanone	0.0051	U	1634-04-4	Methyl-t-butyl ether	0.0010	U J
591-78-6	2-Hexanone	0.0051	U	95-47-6	o-Xylene	0.0010	U
108-10-1	4-Methyl-2-Pentanone	0.0051	U	100-42-5	Styrene	0.0051	U
67-64-1	Acetone	0.026	U	127-18-4	Tetrachloroethene	0.0010	U
71-43-2	Benzene	0.0010	U	108-88-3	Toluene	0.0010	U
75-27-4	Bromodichloromethane	0.0051	U	156-60-5	trans-1,2-Dichloroethene	0.0051	U
75-25-2	Bromoform	0.0051	U	10061-02-6	trans-1,3-Dichloropropene	0.0051	U
74-83-9	Bromomethane	0.0051	U	79-01-6	Trichloroethene	0.0051	U
75-15-0	Carbon Disulfide	0.0051	U	75-69-4	Trichlorofluoromethane	0.0051	U
56-23-5	Carbon Tetrachloride	0.0051	U	75-01-4	Vinyl Chloride	0.0051	U
108-90-7	Chlorobenzene	0.0051	U				

Worksheet #: 57392

Total Target Concentration 0.0066

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-017
Client Id: CPC-SB02-5-7
Data File: 1M24292.D
Analysis Date: 08/24/07 14:07
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 96

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0052	U J	75-00-3	Chloroethane	0.0052	U J
79-34-5	1,1,2,2-Tetrachloroethane	0.0052	U	67-66-3	Chloroform	0.0052	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0052	U	74-87-3	Chloromethane	0.0052	U
79-00-5	1,1,2-Trichloroethane	0.0052	U	156-59-2	cis-1,2-Dichloroethene	0.0052	U
75-34-3	1,1-Dichloroethane	0.0052	U	10061-01-5	cis-1,3-Dichloropropene	0.0052	U
75-35-4	1,1-Dichloroethene	0.0052	U	110-82-7	Cyclohexane	0.0052	U
120-82-1	1,2,4-Trichlorobenzene	0.0052	U	124-48-1	Dibromochloromethane	0.0052	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0052	U	75-71-8	Dichlorodifluoromethane	0.0052	U
106-93-4	1,2-Dibromoethane	0.0052	U	100-41-4	Ethylbenzene	0.0010	U
95-50-1	1,2-Dichlorobenzene	0.0052	U	98-82-8	Isopropylbenzene	0.0010	U
107-06-2	1,2-Dichloroethane	0.0052	U	1330-20-7	m&p-Xylenes	0.0021	U
78-87-5	1,2-Dichloropropane	0.0052	U	79-20-9	Methyl Acetate	0.0052	U
541-73-1	1,3-Dichlorobenzene	0.0052	U	108-87-2	Methylcyclohexane	0.0052	U
106-46-7	1,4-Dichlorobenzene	0.0052	U	**75-09-2**	**Methylene Chloride**	**0.0052**	U 0.011 B J
78-93-3	2-Butanone	0.0052	U	1634-04-4	Methyl-t-butyl ether	0.0010	U J
591-78-6	2-Hexanone	0.0052	U	95-47-6	o-Xylene	0.0010	U
108-10-1	4-Methyl-2-Pentanone	0.0052	U	100-42-5	Styrene	0.0052	U
67-64-1	Acetone	0.026	U	127-18-4	Tetrachloroethene	0.0010	U
71-43-2	Benzene	0.0010	U	108-88-3	Toluene	0.0010	U
75-27-4	Bromodichloromethane	0.0052	U	156-60-5	trans-1,2-Dichloroethene	0.0052	U
75-25-2	Bromoform	0.0052	U	10061-02-6	trans-1,3-Dichloropropene	0.0052	U
74-83-9	Bromomethane	0.0052	U	79-01-6	Trichloroethene	0.0052	U
75-15-0	Carbon Disulfide	0.0052	U	75-69-4	Trichlorofluoromethane	0.0052	U
56-23-5	Carbon Tetrachloride	0.0052	U	75-01-4	Vinyl Chloride	0.0052	U
108-90-7	Chlorobenzene	0.0052	U				

Worksheet #: 57392

Total Target Concentration 0.011

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-018
Client Id: CPC-SB02-7-9
Data File: 1M24293.D
Analysis Date: 08/24/07 14:33
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 97

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0052	U J	75-00-3	Chloroethane	0.0052	U J
79-34-5	1,1,2,2-Tetrachloroethane	0.0052	U	67-66-3	Chloroform	0.0052	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0052	U	74-87-3	Chloromethane	0.0052	U
79-00-5	1,1,2-Trichloroethane	0.0052	U	156-59-2	cis-1,2-Dichloroethene	0.0052	U
75-34-3	1,1-Dichloroethane	0.0052	U	10061-01-5	cis-1,3-Dichloropropene	0.0052	U
75-35-4	1,1-Dichloroethene	0.0052	U	110-82-7	Cyclohexane	0.0052	U
120-82-1	1,2,4-Trichlorobenzene	0.0052	U	124-48-1	Dibromochloromethane	0.0052	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0052	U	75-71-8	Dichlorodifluoromethane	0.0052	U
106-93-4	1,2-Dibromoethane	0.0052	U	100-41-4	Ethylbenzene	0.0010	U
95-50-1	1,2-Dichlorobenzene	0.0052	U	98-82-8	Isopropylbenzene	0.0010	U
107-06-2	1,2-Dichloroethane	0.0052	U	1330-20-7	m&p-Xylenes	0.0021	U
78-87-5	1,2-Dichloropropane	0.0052	U	79-20-9	Methyl Acetate	0.0052	U
541-73-1	1,3-Dichlorobenzene	0.0052	U	108-87-2	Methylcyclohexane	0.0052	U
106-46-7	1,4-Dichlorobenzene	0.0052	U	**75-09-2**	**Methylene Chloride**	**0.0052**	U 0.0073 B
78-93-3	2-Butanone	0.0052	U	1634-04-4	Methyl-t-butyl ether	0.0010	U J
591-78-6	2-Hexanone	0.0052	U	95-47-6	o-Xylene	0.0010	U
108-10-1	4-Methyl-2-Pentanone	0.0052	U	100-42-5	Styrene	0.0052	U
67-64-1	Acetone	0.026	U	127-18-4	Tetrachloroethene	0.0010	U
71-43-2	Benzene	0.0010	U	108-88-3	Toluene	0.0010	U
75-27-4	Bromodichloromethane	0.0052	U	156-60-5	trans-1,2-Dichloroethene	0.0052	U
75-25-2	Bromoform	0.0052	U	10061-02-6	trans-1,3-Dichloropropene	0.0052	U
74-83-9	Bromomethane	0.0052	U	79-01-6	Trichloroethene	0.0052	U
75-15-0	Carbon Disulfide	0.0052	U	75-69-4	Trichlorofluoromethane	0.0052	U
56-23-5	Carbon Tetrachloride	0.0052	U	75-01-4	Vinyl Chloride	0.0052	U
108-90-7	Chlorobenzene	0.0052	U				

Worksheet #: 57392

Total Target Concentration 0.0073

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-019
Client Id: CPC-DW04-13-14
Data File: 1M24294.D
Analysis Date: 08/24/07 14:59
Date Rec/Extracted: 08/16/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 82

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0061	U J	75-00-3	Chloroethane	0.0061	U J
79-34-5	1,1,2,2-Tetrachloroethane	0.0061	U	67-66-3	Chloroform	0.0061	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	0.0061	U	74-87-3	Chloromethane	0.0061	U
79-00-5	1,1,2-Trichloroethane	0.0061	U	156-59-2	cis-1,2-Dichloroethene	0.0061	U
75-34-3	1,1-Dichloroethane	0.0061	U	10061-01-5	cis-1,3-Dichloropropene	0.0061	U
75-35-4	1,1-Dichloroethene	0.0061	U	110-82-7	Cyclohexane	0.0061	U
120-82-1	1,2,4-Trichlorobenzene	0.0061	U	124-48-1	Dibromochloromethane	0.0061	U
96-12-8	1,2-Dibromo-3-Chloropropa	0.0061	U	75-71-8	Dichlorodifluoromethane	0.0061	U
106-93-4	1,2-Dibromoethane	0.0061	U	100-41-4	Ethylbenzene	0.0012	U
95-50-1	1,2-Dichlorobenzene	0.0061	U	98-82-8	Isopropylbenzene	0.0012	U
107-06-2	1,2-Dichloroethane	0.0061	U	1330-20-7	m&p-Xylenes	0.0024	U
78-87-5	1,2-Dichloropropane	0.0061	U	79-20-9	Methyl Acetate	0.0061	U
541-73-1	1,3-Dichlorobenzene	0.0061	U	108-87-2	Methylcyclohexane	0.0061	U
106-46-7	1,4-Dichlorobenzene	0.0061	U	**75-09-2**	**Methylene Chloride**	**0.0061**	U 0.014 B J
78-93-3	2-Butanone	0.0061	U	1634-04-4	Methyl-t-butyl ether	0.0012	U J
591-78-6	2-Hexanone	0.0061	U	95-47-6	o-Xylene	0.0012	U
108-10-1	4-Methyl-2-Pentanone	0.0061	U	100-42-5	Styrene	0.0061	U
67-64-1	Acetone	0.030	U	127-18-4	Tetrachloroethene	0.0012	U
71-43-2	Benzene	0.0012	U	108-88-3	Toluene	0.0012	U
75-27-4	Bromodichloromethane	0.0061	U	156-60-5	trans-1,2-Dichloroethene	0.0061	U
75-25-2	Bromoform	0.0061	U	10061-02-6	trans-1,3-Dichloropropene	0.0061	U
74-83-9	Bromomethane	0.0061	U	79-01-6	Trichloroethene	0.0061	U
75-15-0	Carbon Disulfide	0.0061	U	75-69-4	Trichlorofluoromethane	0.0061	U
56-23-5	Carbon Tetrachloride	0.0061	U	75-01-4	Vinyl Chloride	0.0061	U
108-90-7	Chlorobenzene	0.0061	U				

Worksheet #: 57392

Total Target Concentration 0.014

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC32378-020
Client Id: CPC-Rinsate
Data File: 8M17901.D
Analysis Date: 08/20/07 17:04
Date Rec/Extracted: 08/16/07-NA

Matrix: Aqueous
Initial Vol: 5ml
Final Vol: NA
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	67-66-3	Chloroform	5.0	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluor	5.0	U J	74-87-3	Chloromethane	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	156-59-2	cis-1,2-Dichloroethene	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	110-82-7	Cyclohexane	5.0	U
120-82-1	1,2,4-Trichlorobenzene	5.0	U	124-48-1	Dibromochloromethane	5.0	U
96-12-8	1,2-Dibromo-3-Chloropropa	5.0	U	75-71-8	Dichlorodifluoromethane	5.0	U
106-93-4	1,2-Dibromoethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	98-82-8	Isopropylbenzene	1.0	U
107-06-2	1,2-Dichloroethane	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-87-5	1,2-Dichloropropane	5.0	U	79-20-9	Methyl Acetate	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	108-87-2	Methylcyclohexane	5.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	75-09-2	Methylene Chloride	5.0	U
78-93-3	2-Butanone	5.0	U	1634-04-4	Methyl-t-butyl ether	1.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U J
108-10-1	4-Methyl-2-Pentanone	5.0	U J	100-42-5	Styrene	5.0	U
67-64-1	Acetone	25	U	127-18-4	Tetrachloroethene	5.0	U
71-43-2	Benzene	1.0	U	108-88-3	Toluene	1.0	U
75-27-4	Bromodichloromethane	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
75-25-2	Bromoform	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
74-83-9	Bromomethane	5.0	U	79-01-6	Trichloroethene	5.0	U
75-15-0	Carbon Disulfide	5.0	U	75-69-4	Trichlorofluoromethane	5.0	U J
56-23-5	Carbon Tetrachloride	5.0	U J	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 57392

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-001
Client Id: CPC-MW01
Data File: 5M34146.D
Analysis Date: 08/21/07 21:39
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	25	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-002
Client Id: CPC-MW08
Data File: 5M34147.D
Analysis Date: 08/21/07 22:01
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**10**	**1.5 J**
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	25	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 1.5

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-003
Client Id: CPC-MW03
Data File: 5M34149.D
Analysis Date: 08/21/07 22:45
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 980ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	26	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-004
Client Id: CPC-MW04
Data File: 5M34133.D
Analysis Date: 08/21/07 16:52
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 980ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	26	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-005
Client Id: CPC-MW05
Data File: 5M34134.D
Analysis Date: 08/21/07 17:15
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 980ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	26	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-006
Client Id: CPC-MW06
Data File: 5M34135.D
Analysis Date: 08/21/07 17:37
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 980ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	26	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-007
Client Id: CPC-MW07
Data File: 5M34136.D
Analysis Date: 08/21/07 17:59
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 980ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	26	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-008
Client Id: CPC-MW09
Data File: 5M34145.D
Analysis Date: 08/21/07 21:17
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	25	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-009
Client Id: CPC-MW10
Data File: 5M33991.D
Analysis Date: 08/17/07 19:45
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]peryiene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	25	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**10**	**1.1 J**
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 1.1

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32351-012
Client Id: CPC-DUPLICATE
Data File: 5M34148.D
Analysis Date: 08/21/07 22:23
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
105-67-9	2,4-Dimethylphenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
121-14-2	2,4-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	105-60-2	Caprolactam	10	U
91-58-7	2-Chloronaphthalene	10	U	86-74-8	Carbazole	10	U
95-57-8	2-Chlorophenol	10	U	218-01-9	Chrysene	10	U
91-57-6	2-Methylnaphthalene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
95-48-7	2-Methylphenol	10	U	132-64-9	Dibenzofuran	10	U
88-74-4	2-Nitroaniline	10	U	84-66-2	Diethylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	131-11-3	Dimethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	117-84-0	Di-n-octylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	206-44-0	Fluoranthene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	86-73-7	Fluorene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	118-74-1	Hexachlorobenzene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	87-68-3	Hexachlorobutadiene	10	U
106-47-8	4-Chloroaniline	10	U	77-47-4	Hexachlorocyclopentadiene	25	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
100-01-6	4-Nitroaniline	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
100-02-7	4-Nitrophenol	10	U	78-59-1	Isophorone	10	U
83-32-9	Acenaphthene	10	U	91-20-3	Naphthalene	10	U
208-96-8	Acenaphthylene	10	U	98-95-3	Nitrobenzene	10	U
98-86-2	Acetophenone	10	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
120-12-7	Anthracene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
1912-24-9	Atrazine	10	U	87-86-5	Pentachlorophenol	10	U
100-52-7	Benzaldehyde	10	U	85-01-8	Phenanthrene	10	U
56-55-3	Benzo[a]anthracene	10	U	108-95-2	Phenol	10	U
50-32-8	Benzo[a]pyrene	10	U	129-00-0	Pyrene	10	U
205-99-2	Benzo[b]fluoranthene	10	U				

Worksheet #: 57381

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-001
Client Id: CPC-DW01-10-11
Data File: 5M34186.D
Analysis Date: 08/22/07 21:36
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 86

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.39	U	205-99-2	Benzo[b]fluoranthene	0.39	U
95-95-4	2,4,5-Trichlorophenol	0.39	U	191-24-2	Benzo[g,h,i]perylene	0.39	U
88-06-2	2,4,6-Trichlorophenol	0.39	U	207-08-9	Benzo[k]fluoranthene	0.39	U
120-83-2	2,4-Dichlorophenol	0.39	U	111-91-1	bis(2-Chloroethoxy)methan	0.39	U
105-67-9	2,4-Dimethylphenol	0.39	U	111-44-4	bis(2-Chloroethyl)ether	0.39	U
51-28-5	2,4-Dinitrophenol	0.39	U	39638-32-9	bis(2-chloroisopropyl)ether	0.39	U
121-14-2	2,4-Dinitrotoluene	0.39	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.39**	**0.049 J**
606-20-2	2,6-Dinitrotoluene	0.39	U	85-68-7	Butylbenzylphthalate	0.39	U
91-58-7	2-Chloronaphthalene	0.39	U	105-60-2	Caprolactam	0.39	U
95-57-8	2-Chlorophenol	0.39	U	86-74-8	Carbazole	0.39	U
91-57-6	2-Methylnaphthalene	0.39	U	218-01-9	Chrysene	0.39	U
95-48-7	2-Methylphenol	0.39	U	53-70-3	Dibenzo[a,h]anthracene	0.39	U
88-74-4	2-Nitroaniline	0.39	U	132-64-9	Dibenzofuran	0.39	U
88-75-5	2-Nitrophenol	0.39	U	84-66-2	Diethylphthalate	0.39	U
106-44-5	3&4-Methylphenol	0.39	U	131-11-3	Dimethylphthalate	0.39	U
91-94-1	3,3'-Dichlorobenzidine	0.39	U	84-74-2	Di-n-butylphthalate	0.39	U
99-09-2	3-Nitroaniline	0.39	U	117-84-0	Di-n-octylphthalate	0.39	U
534-52-1	4,6-Dinitro-2-methylphenol	0.39	U	206-44-0	Fluoranthene	0.39	U
101-55-3	4-Bromophenyl-phenylether	0.39	U	86-73-7	Fluorene	0.39	U
59-50-7	4-Chloro-3-methylphenol	0.39	U	118-74-1	Hexachlorobenzene	0.39	U
106-47-8	4-Chloroaniline	0.39	U	87-68-3	Hexachlorobutadiene	0.39	U
7005-72-3	4-Chlorophenyl-phenylether	0.39	U	77-47-4	Hexachlorocyclopentadiene	0.97	U
100-01-6	4-Nitroaniline	0.39	U	67-72-1	Hexachloroethane	0.39	U
100-02-7	4-Nitrophenol	0.39	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.39	U
83-32-9	Acenaphthene	0.39	U	78-59-1	Isophorone	0.39	U
208-96-8	Acenaphthylene	0.39	U	91-20-3	Naphthalene	0.39	U
98-86-2	Acetophenone	0.39	U	98-95-3	Nitrobenzene	0.39	U
621-64-7	N-Nitroso-di-n-propylamine	0.39	U	120-12-7	Anthracene	0.39	U
86-30-6	n-Nitrosodiphenylamine	0.39	U	1912-24-9	Atrazine	0.39	U
87-86-5	Pentachlorophenol	0.39	U	100-52-7	Benzaldehyde	0.39	U
85-01-8	Phenanthrene	0.39	U	56-55-3	Benzo[a]anthracene	0.39	U
108-95-2	Phenol	0.39	U	50-32-8	Benzo[a]pyrene	0.39	U
129-00-0	Pyrene	0.39	U				

Worksheet #: 57379

Total Target Concentration 0.049

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-002
Client Id: CPC-DW02-10-11
Data File: 9M05059.D
Analysis Date: 08/23/07 04:19
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 83

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.40	U	**205-99-2**	**Benzo[b]fluoranthene**	**0.40**	**0.072 J**
95-95-4	2,4,5-Trichlorophenol	0.40	U	**191-24-2**	**Benzo[g,h,i]perylene**	**0.40**	**0.061 J**
88-06-2	2,4,6-Trichlorophenol	0.40	U	207-08-9	Benzo[k]fluoranthene	0.40	U
120-83-2	2,4-Dichlorophenol	0.40	U	111-91-1	bis(2-Chloroethoxy)methan	0.40	U
105-67-9	2,4-Dimethylphenol	0.40	U	111-44-4	bis(2-Chloroethyl)ether	0.40	U
51-28-5	2,4-Dinitrophenol	0.40	U	39638-32-9	bis(2-chloroisopropyl)ether	0.40	U
121-14-2	2,4-Dinitrotoluene	0.40	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.40**	**0.42**
606-20-2	2,6-Dinitrotoluene	0.40	U	85-68-7	Butylbenzylphthalate	0.40	U
91-58-7	2-Chloronaphthalene	0.40	U	105-60-2	Caprolactam	0.40	U
95-57-8	2-Chlorophenol	0.40	U	86-74-8	Carbazole	0.40	U
91-57-6	2-Methylnaphthalene	0.40	U	**218-01-9**	**Chrysene**	**0.40**	**0.052 J**
95-48-7	2-Methylphenol	0.40	U	53-70-3	Dibenzo[a,h]anthracene	0.40	U
88-74-4	2-Nitroaniline	0.40	U	132-64-9	Dibenzofuran	0.40	U
88-75-5	2-Nitrophenol	0.40	U	84-66-2	Diethylphthalate	0.40	U
106-44-5	3&4-Methylphenol	0.40	U	131-11-3	Dimethylphthalate	0.40	U
91-94-1	3,3'-Dichlorobenzidine	0.40	U	84-74-2	Di-n-butylphthalate	0.40	U
99-09-2	3-Nitroaniline	0.40	U	117-84-0	Di-n-octylphthalate	0.40	U
534-52-1	4,6-Dinitro-2-methylphenol	0.40	U	**206-44-0**	**Fluoranthene**	**0.40**	**0.071 J**
101-55-3	4-Bromophenyl-phenylether	0.40	U	86-73-7	Fluorene	0.40	U
59-50-7	4-Chloro-3-methylphenol	0.40	U	118-74-1	Hexachlorobenzene	0.40	U
106-47-8	4-Chloroaniline	0.40	U	87-68-3	Hexachlorobutadiene	0.40	U
7005-72-3	4-Chlorophenyl-phenylether	0.40	U	77-47-4	Hexachlorocyclopentadiene	0.40	U
100-01-6	4-Nitroaniline	0.40	U	67-72-1	Hexachloroethane	0.40	U
100-02-7	4-Nitrophenol	0.40	U	**193-39-5**	**Indeno[1,2,3-cd]pyrene**	**0.40**	**0.042 J**
83-32-9	Acenaphthene	0.40	U	78-59-1	Isophorone	0.40	U
208-96-8	Acenaphthylene	0.40	U	91-20-3	Naphthalene	0.40	U
98-86-2	Acetophenone	0.40	U	98-95-3	Nitrobenzene	0.40	U
621-64-7	N-Nitroso-di-n-propylamine	0.40	U	120-12-7	Anthracene	0.40	U
86-30-6	n-Nitrosodiphenylamine	0.40	U	1912-24-9	Atrazine	0.40	U
87-86-5	Pentachlorophenol	0.40	U	100-52-7	Benzaldehyde	0.40	U
85-01-8	Phenanthrene	0.40	U	56-55-3	Benzo[a]anthracene	0.40	U
108-95-2	Phenol	0.40	U	**50-32-8**	**Benzo[a]pyrene**	**0.40**	**0.050 J**
129-00-0	**Pyrene**	**0.40**	**0.088 J**				

Worksheet #: 57379

Total Target Concentration 0.856

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-003
Client Id: CPC-DW03-10-11
Data File: 9M05048.D
Analysis Date: 08/22/07 23:57
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 90

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.37	U	205-99-2	Benzo[b]fluoranthene	0.37	U
95-95-4	2,4,5-Trichlorophenol	0.37	U	191-24-2	Benzo[g,h,i]perylene	0.37	U
88-06-2	2,4,6-Trichlorophenol	0.37	U	207-08-9	Benzo[k]fluoranthene	0.37	U
120-83-2	2,4-Dichlorophenol	0.37	U	111-91-1	bis(2-Chloroethoxy)methan	0.37	U
105-67-9	2,4-Dimethylphenol	0.37	U	111-44-4	bis(2-Chloroethyl)ether	0.37	U
51-28-5	2,4-Dinitrophenol	0.37	U	39638-32-9	bis(2-chloroisopropyl)ether	0.37	U
121-14-2	2,4-Dinitrotoluene	0.37	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.37**	**0.10 J**
606-20-2	2,6-Dinitrotoluene	0.37	U	85-68-7	Butylbenzylphthalate	0.37	U
91-58-7	2-Chloronaphthalene	0.37	U	105-60-2	Caprolactam	0.37	U
95-57-8	2-Chlorophenol	0.37	U	86-74-8	Carbazole	0.37	U
91-57-6	2-Methylnaphthalene	0.37	U	218-01-9	Chrysene	0.37	U
95-48-7	2-Methylphenol	0.37	U	53-70-3	Dibenzo[a,h]anthracene	0.37	U
88-74-4	2-Nitroaniline	0.37	U	132-64-9	Dibenzofuran	0.37	U
88-75-5	2-Nitrophenol	0.37	U	84-66-2	Diethylphthalate	0.37	U
106-44-5	3&4-Methylphenol	0.37	U	131-11-3	Dimethylphthalate	0.37	U
91-94-1	3,3'-Dichlorobenzidine	0.37	U	**84-74-2**	**Di-n-butylphthalate**	**0.37**	**0.070 J**
99-09-2	3-Nitroaniline	0.37	U	117-84-0	Di-n-octylphthalate	0.37	U
534-52-1	4,6-Dinitro-2-methylphenol	0.37	U	206-44-0	Fluoranthene	0.37	U
101-55-3	4-Bromophenyl-phenylether	0.37	U	86-73-7	Fluorene	0.37	U
59-50-7	4-Chloro-3-methylphenol	0.37	U	118-74-1	Hexachlorobenzene	0.37	U
106-47-8	4-Chloroaniline	0.37	U	87-68-3	Hexachlorobutadiene	0.37	U
7005-72-3	4-Chlorophenyl-phenylether	0.37	U	77-47-4	Hexachlorocyclopentadiene	0.37	U
100-01-6	4-Nitroaniline	0.37	U	67-72-1	Hexachloroethane	0.37	U
100-02-7	4-Nitrophenol	0.37	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.37	U
83-32-9	Acenaphthene	0.37	U	78-59-1	Isophorone	0.37	U
208-96-8	Acenaphthylene	0.37	U	91-20-3	Naphthalene	0.37	U
98-86-2	Acetophenone	0.37	U	98-95-3	Nitrobenzene	0.37	U
621-64-7	N-Nitroso-di-n-propylamine	0.37	U	120-12-7	Anthracene	0.37	U
86-30-6	n-Nitrosodiphenylamine	0.37	U	1912-24-9	Atrazine	0.37	U
87-86-5	Pentachlorophenol	0.37	U	100-52-7	Benzaldehyde	0.37	U
85-01-8	Phenanthrene	0.37	U	56-55-3	Benzo[a]anthracene	0.37	U
108-95-2	Phenol	0.37	U	50-32-8	Benzo[a]pyrene	0.37	U
129-00-0	Pyrene	0.37	U				

Worksheet #: 57379

Total Target Concentration 0.17

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-004(2X)
Client Id: CPC-DW04-8-9
Data File: 9M05092.D
Analysis Date: 08/23/07 23:28
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 2
Solids: 85

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.78	U	**205-99-2**	**Benzo[b]fluoranthene**	**0.78**	**0.43 J**
95-95-4	2,4,5-Trichlorophenol	0.78	U	**191-24-2**	**Benzo[g,h,i]perylene**	**0.78**	**0.19 J**
88-06-2	2,4,6-Trichlorophenol	0.78	U	**207-08-9**	**Benzo[k]fluoranthene**	**0.78**	**0.10 J**
120-83-2	2,4-Dichlorophenol	0.78	U	111-91-1	bis(2-Chloroethoxy)methan	0.78	U
105-67-9	2,4-Dimethylphenol	0.78	U	111-44-4	bis(2-Chloroethyl)ether	0.78	U
51-28-5	2,4-Dinitrophenol	0.78	U	39638-32-9	bis(2-chloroisopropyl)ether	0.78	U
121-14-2	2,4-Dinitrotoluene	0.78	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.78**	**7.1**
606-20-2	2,6-Dinitrotoluene	0.78	U	**85-68-7**	**Butylbenzylphthalate**	**0.78**	**0.63 J**
91-58-7	2-Chloronaphthalene	0.78	U	105-60-2	Caprolactam	0.78	U
95-57-8	2-Chlorophenol	0.78	U	86-74-8	Carbazole	0.78	U
91-57-6	**2-Methylnaphthalene**	**0.78**	**0.97**	**218-01-9**	**Chrysene**	**0.78**	**0.39 J**
95-48-7	2-Methylphenol	0.78	U	53-70-3	Dibenzo[a,h]anthracene	0.78	U
88-74-4	2-Nitroaniline	0.78	U	132-64-9	Dibenzofuran	0.78	U
88-75-5	2-Nitrophenol	0.78	U	84-66-2	Diethylphthalate	0.78	U
106-44-5	3&4-Methylphenol	0.78	U	131-11-3	Dimethylphthalate	0.78	U
91-94-1	3,3'-Dichlorobenzidine	0.78	U	**84-74-2**	**Di-n-butylphthalate**	**0.78**	**0.23 J**
99-09-2	3-Nitroaniline	0.78	U	117-84-0	Di-n-octylphthalate	0.78	U
534-52-1	4,6-Dinitro-2-methylphenol	0.78	U	**206-44-0**	**Fluoranthene**	**0.78**	**0.61 J**
101-55-3	4-Bromophenyl-phenylether	0.78	U	**86-73-7**	**Fluorene**	**0.78**	**0.12 J**
59-50-7	4-Chloro-3-methylphenol	0.78	U	118-74-1	Hexachlorobenzene	0.78	U
106-47-8	4-Chloroaniline	0.78	U	87-68-3	Hexachlorobutadiene	0.78	U
7005-72-3	4-Chlorophenyl-phenylether	0.78	U	77-47-4	Hexachlorocyclopentadiene	0.78	U
100-01-6	4-Nitroaniline	0.78	U	67-72-1	Hexachloroethane	0.78	U
100-02-7	4-Nitrophenol	0.78	U	**193-39-5**	**Indeno[1,2,3-cd]pyrene**	**0.78**	**0.16 J**
83-32-9	Acenaphthene	0.78	U	78-59-1	Isophorone	0.78	U
208-96-8	Acenaphthylene	0.78	U	**91-20-3**	**Naphthalene**	**0.78**	**0.47 J**
98-86-2	Acetophenone	0.78	U	98-95-3	Nitrobenzene	0.78	U
621-64-7	N-Nitroso-di-n-propylamine	0.78	U	120-12-7	Anthracene	0.78	U
86-30-6	n-Nitrosodiphenylamine	0.78	U	1912-24-9	Atrazine	0.78	U
87-86-5	Pentachlorophenol	0.78	U	100-52-7	Benzaldehyde	0.78	U
85-01-8	**Phenanthrene**	**0.78**	**0.48 J**	**56-55-3**	**Benzo[a]anthracene**	**0.78**	**0.20 J**
108-95-2	Phenol	0.78	U	**50-32-8**	**Benzo[a]pyrene**	**0.78**	**0.18 J**
129-00-0	**Pyrene**	**0.78**	**0.61 J**				

Worksheet #: 57379

Total Target Concentration 12.87

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-005
Client Id: CPC-SOIL-Duplicate
Data File: 9M05049.D
Analysis Date: 08/23/07 00:21
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 92

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.36	U	205-99-2	Benzo[b]fluoranthene	0.36	U
95-95-4	2,4,5-Trichlorophenol	0.36	U	191-24-2	Benzo[g,h,i]perylene	0.36	U
88-06-2	2,4,6-Trichlorophenol	0.36	U	207-08-9	Benzo[k]fluoranthene	0.36	U
120-83-2	2,4-Dichlorophenol	0.36	U	111-91-1	bis(2-Chloroethoxy)methan	0.36	U
105-67-9	2,4-Dimethylphenol	0.36	U	111-44-4	bis(2-Chloroethyl)ether	0.36	U
51-28-5	2,4-Dinitrophenol	0.36	U	39638-32-9	bis(2-chloroisopropyl)ether	0.36	U
121-14-2	2,4-Dinitrotoluene	0.36	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.36**	**0.085 J**
606-20-2	2,6-Dinitrotoluene	0.36	U	85-68-7	Butylbenzylphthalate	0.36	U
91-58-7	2-Chloronaphthalene	0.36	U	105-60-2	Caprolactam	0.36	U
95-57-8	2-Chlorophenol	0.36	U	86-74-8	Carbazole	0.36	U
91-57-6	2-Methylnaphthalene	0.36	U	218-01-9	Chrysene	0.36	U
95-48-7	2-Methylphenol	0.36	U	53-70-3	Dibenzo[a,h]anthracene	0.36	U
88-74-4	2-Nitroaniline	0.36	U	132-64-9	Dibenzofuran	0.36	U
88-75-5	2-Nitrophenol	0.36	U	84-66-2	Diethylphthalate	0.36	U
106-44-5	3&4-Methylphenol	0.36	U	131-11-3	Dimethylphthalate	0.36	U
91-94-1	3,3'-Dichlorobenzidine	0.36	U	84-74-2	Di-n-butylphthalate	0.36	U
99-09-2	3-Nitroaniline	0.36	U	117-84-0	Di-n-octylphthalate	0.36	U
534-52-1	4,6-Dinitro-2-methylphenol	0.36	U	206-44-0	Fluoranthene	0.36	U
101-55-3	4-Bromophenyl-phenylether	0.36	U	86-73-7	Fluorene	0.36	U
59-50-7	4-Chloro-3-methylphenol	0.36	U	118-74-1	Hexachlorobenzene	0.36	U
106-47-8	4-Chloroaniline	0.36	U	87-68-3	Hexachlorobutadiene	0.36	U
7005-72-3	4-Chlorophenyl-phenylether	0.36	U	77-47-4	Hexachlorocyclopentadiene	0.36	U
100-01-6	4-Nitroaniline	0.36	U	67-72-1	Hexachloroethane	0.36	U
100-02-7	4-Nitrophenol	0.36	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.36	U
83-32-9	Acenaphthene	0.36	U	78-59-1	Isophorone	0.36	U
208-96-8	Acenaphthylene	0.36	U	91-20-3	Naphthalene	0.36	U
98-86-2	Acetophenone	0.36	U	98-95-3	Nitrobenzene	0.36	U
621-64-7	N-Nitroso-di-n-propylamine	0.36	U	120-12-7	Anthracene	0.36	U
86-30-6	n-Nitrosodiphenylamine	0.36	U	1912-24-9	Atrazine	0.36	U
87-86-5	Pentachlorophenol	0.36	U	100-52-7	Benzaldehyde	0.36	U
85-01-8	Phenanthrene	0.36	U	56-55-3	Benzo[a]anthracene	0.36	U
108-95-2	Phenol	0.36	U	50-32-8	Benzo[a]pyrene	0.36	U
129-00-0	Pyrene	0.36	U				

Worksheet #: 57379

Total Target Concentration 0.085

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-006
Client Id: CPC-SB01-1.5-2.5
Data File: 9M05050.D
Analysis Date: 08/23/07 00:45
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 95

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.35	U	205-99-2	Benzo[b]fluoranthene	0.35	U
95-95-4	2,4,5-Trichlorophenol	0.35	U	191-24-2	Benzo[g,h,i]perylene	0.35	U
88-06-2	2,4,6-Trichlorophenol	0.35	U	207-08-9	Benzo[k]fluoranthene	0.35	U
120-83-2	2,4-Dichlorophenol	0.35	U	111-91-1	bis(2-Chloroethoxy)methan	0.35	U
105-67-9	2,4-Dimethylphenol	0.35	U	111-44-4	bis(2-Chloroethyl)ether	0.35	U
51-28-5	2,4-Dinitrophenol	0.35	U	39638-32-9	bis(2-chloroisopropyl)ether	0.35	U
121-14-2	2,4-Dinitrotoluene	0.35	U	117-81-7	bis(2-Ethylhexyl)phthalate	0.35	U
606-20-2	2,6-Dinitrotoluene	0.35	U	85-68-7	Butylbenzylphthalate	0.35	U
91-58-7	2-Chloronaphthalene	0.35	U	105-60-2	Caprolactam	0.35	U
95-57-8	2-Chlorophenol	0.35	U	86-74-8	Carbazole	0.35	U
91-57-6	2-Methylnaphthalene	0.35	U	218-01-9	Chrysene	0.35	U
95-48-7	2-Methylphenol	0.35	U	53-70-3	Dibenzo[a,h]anthracene	0.35	U
88-74-4	2-Nitroaniline	0.35	U	132-64-9	Dibenzofuran	0.35	U
88-75-5	2-Nitrophenol	0.35	U	84-66-2	Diethylphthalate	0.35	U
106-44-5	3&4-Methylphenol	0.35	U	131-11-3	Dimethylphthalate	0.35	U
91-94-1	3,3'-Dichlorobenzidine	0.35	U	84-74-2	Di-n-butylphthalate	0.35	U
99-09-2	3-Nitroaniline	0.35	U	117-84-0	Di-n-octylphthalate	0.35	U
534-52-1	4,6-Dinitro-2-methylphenol	0.35	U	206-44-0	Fluoranthene	0.35	U
101-55-3	4-Bromophenyl-phenylether	0.35	U	86-73-7	Fluorene	0.35	U
59-50-7	4-Chloro-3-methylphenol	0.35	U	118-74-1	Hexachlorobenzene	0.35	U
106-47-8	4-Chloroaniline	0.35	U	87-68-3	Hexachlorobutadiene	0.35	U
7005-72-3	4-Chlorophenyl-phenylether	0.35	U	77-47-4	Hexachlorocyclopentadiene	0.35	U
100-01-6	4-Nitroaniline	0.35	U	67-72-1	Hexachloroethane	0.35	U
100-02-7	4-Nitrophenol	0.35	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.35	U
83-32-9	Acenaphthene	0.35	U	78-59-1	Isophorone	0.35	U
208-96-8	Acenaphthylene	0.35	U	91-20-3	Naphthalene	0.35	U
98-86-2	Acetophenone	0.35	U	98-95-3	Nitrobenzene	0.35	U
621-64-7	N-Nitroso-di-n-propylamine	0.35	U	120-12-7	Anthracene	0.35	U
86-30-6	n-Nitrosodiphenylamine	0.35	U	1912-24-9	Atrazine	0.35	U
87-86-5	Pentachlorophenol	0.35	U	100-52-7	Benzaldehyde	0.35	U
85-01-8	Phenanthrene	0.35	U	56-55-3	Benzo[a]anthracene	0.35	U
108-95-2	Phenol	0.35	U	50-32-8	Benzo[a]pyrene	0.35	U
129-00-0	Pyrene	0.35	U				

Worksheet #: 57379

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-007
Client Id: CPC-SB02-1-5
Data File: 9M05051.D
Analysis Date: 08/23/07 01:08
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 94

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.35	U	205-99-2	Benzo[b]fluoranthene	0.35	U
95-95-4	2,4,5-Trichlorophenol	0.35	U	191-24-2	Benzo[g,h,i]perylene	0.35	U
88-06-2	2,4,6-Trichlorophenol	0.35	U	207-08-9	Benzo[k]fluoranthene	0.35	U
120-83-2	2,4-Dichlorophenol	0.35	U	111-91-1	bis(2-Chloroethoxy)methan	0.35	U
105-67-9	2,4-Dimethylphenol	0.35	U	111-44-4	bis(2-Chloroethyl)ether	0.35	U
51-28-5	2,4-Dinitrophenol	0.35	U	39638-32-9	bis(2-chloroisopropyl)ether	0.35	U
121-14-2	2,4-Dinitrotoluene	0.35	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.35**	**0.084 J**
606-20-2	2,6-Dinitrotoluene	0.35	U	85-68-7	Butylbenzylphthalate	0.35	U
91-58-7	2-Chloronaphthalene	0.35	U	105-60-2	Caprolactam	0.35	U
95-57-8	2-Chlorophenol	0.35	U	86-74-8	Carbazole	0.35	U
91-57-6	2-Methylnaphthalene	0.35	U	218-01-9	Chrysene	0.35	U
95-48-7	2-Methylphenol	0.35	U	53-70-3	Dibenzo[a,h]anthracene	0.35	U
88-74-4	2-Nitroaniline	0.35	U	132-64-9	Dibenzofuran	0.35	U
88-75-5	2-Nitrophenol	0.35	U	84-66-2	Diethylphthalate	0.35	U
106-44-5	3&4-Methylphenol	0.35	U	131-11-3	Dimethylphthalate	0.35	U
91-94-1	3,3'-Dichlorobenzidine	0.35	U	**84-74-2**	**Di-n-butylphthalate**	**0.35**	**0.038 J**
99-09-2	3-Nitroaniline	0.35	U	117-84-0	Di-n-octylphthalate	0.35	U
534-52-1	4,6-Dinitro-2-methylphenol	0.35	U	206-44-0	Fluoranthene	0.35	U
101-55-3	4-Bromophenyl-phenylether	0.35	U	86-73-7	Fluorene	0.35	U
59-50-7	4-Chloro-3-methylphenol	0.35	U	118-74-1	Hexachlorobenzene	0.35	U
106-47-8	4-Chloroaniline	0.35	U	87-68-3	Hexachlorobutadiene	0.35	U
7005-72-3	4-Chlorophenyl-phenylether	0.35	U	77-47-4	Hexachlorocyclopentadiene	0.35	U
100-01-6	4-Nitroaniline	0.35	U	67-72-1	Hexachloroethane	0.35	U
100-02-7	4-Nitrophenol	0.35	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.35	U
83-32-9	Acenaphthene	0.35	U	78-59-1	Isophorone	0.35	U
208-96-8	Acenaphthylene	0.35	U	91-20-3	Naphthalene	0.35	U
98-86-2	Acetophenone	0.35	U	98-95-3	Nitrobenzene	0.35	U
621-64-7	N-Nitroso-di-n-propylamine	0.35	U	120-12-7	Anthracene	0.35	U
86-30-6	n-Nitrosodiphenylamine	0.35	U	1912-24-9	Atrazine	0.35	U
87-86-5	Pentachlorophenol	0.35	U	100-52-7	Benzaldehyde	0.35	U
85-01-8	Phenanthrene	0.35	U	56-55-3	Benzo[a]anthracene	0.35	U
108-95-2	Phenol	0.35	U	50-32-8	Benzo[a]pyrene	0.35	U
129-00-0	Pyrene	0.35	U				

Worksheet #: 57379

Total Target Concentration 0.122

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-008
Client Id: CPC-SB03-1.5-3.5
Data File: 5M34220.D
Analysis Date: 08/23/07 17:56
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 93

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.36	U	205-99-2	Benzo[b]fluoranthene	0.36	U
95-95-4	2,4,5-Trichlorophenol	0.36	U	191-24-2	Benzo[g,h,i]perylene	0.36	U
88-06-2	2,4,6-Trichlorophenol	0.36	U	207-08-9	Benzo[k]fluoranthene	0.36	U
120-83-2	2,4-Dichlorophenol	0.36	U	111-91-1	bis(2-Chloroethoxy)methan	0.36	U
105-67-9	2,4-Dimethylphenol	0.36	U	111-44-4	bis(2-Chloroethyl)ether	0.36	U
51-28-5	2,4-Dinitrophenol	0.36	U	39638-32-9	bis(2-chloroisopropyl)ether	0.36	U
121-14-2	2,4-Dinitrotoluene	0.36	U	117-81-7	bis(2-Ethylhexyl)phthalate	0.36	U
606-20-2	2,6-Dinitrotoluene	0.36	U	85-68-7	Butylbenzylphthalate	0.36	U
91-58-7	2-Chloronaphthalene	0.36	U	105-60-2	Caprolactam	0.36	U
95-57-8	2-Chlorophenol	0.36	U	86-74-8	Carbazole	0.36	U
91-57-6	2-Methylnaphthalene	0.36	U	218-01-9	Chrysene	0.36	U
95-48-7	2-Methylphenol	0.36	U	53-70-3	Dibenzo[a,h]anthracene	0.36	U
88-74-4	2-Nitroaniline	0.36	U	132-64-9	Dibenzofuran	0.36	U
88-75-5	2-Nitrophenol	0.36	U	84-66-2	Diethylphthalate	0.36	U
106-44-5	3&4-Methylphenol	0.36	U	131-11-3	Dimethylphthalate	0.36	U
91-94-1	3,3'-Dichlorobenzidine	0.36	U	84-74-2	Di-n-butylphthalate	0.36	U
99-09-2	3-Nitroaniline	0.36	U	117-84-0	Di-n-octylphthalate	0.36	U
534-52-1	4,6-Dinitro-2-methylphenol	0.36	U	206-44-0	Fluoranthene	0.36	U
101-55-3	4-Bromophenyl-phenylether	0.36	U	86-73-7	Fluorene	0.36	U
59-50-7	4-Chloro-3-methylphenol	0.36	U	118-74-1	Hexachlorobenzene	0.36	U
106-47-8	4-Chloroaniline	0.36	U	87-68-3	Hexachlorobutadiene	0.36	U
7005-72-3	4-Chlorophenyl-phenylether	0.36	U	77-47-4	Hexachlorocyclopentadiene	0.90	U
100-01-6	4-Nitroaniline	0.36	U	67-72-1	Hexachloroethane	0.36	U
100-02-7	4-Nitrophenol	0.36	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.36	U
83-32-9	Acenaphthene	0.36	U	78-59-1	Isophorone	0.36	U
208-96-8	Acenaphthylene	0.36	U	91-20-3	Naphthalene	0.36	U
98-86-2	Acetophenone	0.36	U	98-95-3	Nitrobenzene	0.36	U
621-64-7	N-Nitroso-di-n-propylamine	0.36	U	120-12-7	Anthracene	0.36	U
86-30-6	n-Nitrosodiphenylamine	0.36	U	1912-24-9	Atrazine	0.36	U
87-86-5	Pentachlorophenol	0.36	U	100-52-7	Benzaldehyde	0.36	U
85-01-8	Phenanthrene	0.36	U	56-55-3	Benzo[a]anthracene	0.36	U
108-95-2	Phenol	0.36	U	50-32-8	Benzo[a]pyrene	0.36	U
129-00-0	Pyrene	0.36	U				

Worksheet #: 57379

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-011
Client Id: CPC-SB04-1.5-2.5
Data File: 9M05052.D
Analysis Date: 08/23/07 01:32
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 90

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.37	U	205-99-2	Benzo[b]fluoranthene	0.37	U
95-95-4	2,4,5-Trichlorophenol	0.37	U	191-24-2	Benzo[g,h,i]perylene	0.37	U
88-06-2	2,4,6-Trichlorophenol	0.37	U	207-08-9	Benzo[k]fluoranthene	0.37	U
120-83-2	2,4-Dichlorophenol	0.37	U	111-91-1	bis(2-Chloroethoxy)methan	0.37	U
105-67-9	2,4-Dimethylphenol	0.37	U	111-44-4	bis(2-Chloroethyl)ether	0.37	U
51-28-5	2,4-Dinitrophenol	0.37	U	39638-32-9	bis(2-chloroisopropyl)ether	0.37	U
121-14-2	2,4-Dinitrotoluene	0.37	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.37**	**0.17 J**
606-20-2	2,6-Dinitrotoluene	0.37	U	85-68-7	Butylbenzylphthalate	0.37	U
91-58-7	2-Chloronaphthalene	0.37	U	105-60-2	Caprolactam	0.37	U
95-57-8	2-Chlorophenol	0.37	U	86-74-8	Carbazole	0.37	U
91-57-6	2-Methylnaphthalene	0.37	U	218-01-9	Chrysene	0.37	U
95-48-7	2-Methylphenol	0.37	U	53-70-3	Dibenzo[a,h]anthracene	0.37	U
88-74-4	2-Nitroaniline	0.37	U	132-64-9	Dibenzofuran	0.37	U
88-75-5	2-Nitrophenol	0.37	U	84-66-2	Diethylphthalate	0.37	U
106-44-5	3&4-Methylphenol	0.37	U	131-11-3	Dimethylphthalate	0.37	U
91-94-1	3,3'-Dichlorobenzidine	0.37	U	84-74-2	Di-n-butylphthalate	0.37	U
99-09-2	3-Nitroaniline	0.37	U	117-84-0	Di-n-octylphthalate	0.37	U
534-52-1	4,6-Dinitro-2-methylphenol	0.37	U	206-44-0	Fluoranthene	0.37	U
101-55-3	4-Bromophenyl-phenylether	0.37	U	86-73-7	Fluorene	0.37	U
59-50-7	4-Chloro-3-methylphenol	0.37	U	118-74-1	Hexachlorobenzene	0.37	U
106-47-8	4-Chloroaniline	0.37	U	87-68-3	Hexachlorobutadiene	0.37	U
7005-72-3	4-Chlorophenyl-phenylether	0.37	U	77-47-4	Hexachlorocyclopentadiene	0.37	U
100-01-6	4-Nitroaniline	0.37	U	67-72-1	Hexachloroethane	0.37	U
100-02-7	4-Nitrophenol	0.37	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.37	U
83-32-9	Acenaphthene	0.37	U	78-59-1	Isophorone	0.37	U
208-96-8	Acenaphthylene	0.37	U	91-20-3	Naphthalene	0.37	U
98-86-2	Acetophenone	0.37	U	98-95-3	Nitrobenzene	0.37	U
621-64-7	N-Nitroso-di-n-propylamine	0.37	U	120-12-7	Anthracene	0.37	U
86-30-6	n-Nitrosodiphenylamine	0.37	U	1912-24-9	Atrazine	0.37	U
87-86-5	Pentachlorophenol	0.37	U	100-52-7	Benzaldehyde	0.37	U
85-01-8	Phenanthrene	0.37	U	56-55-3	Benzo[a]anthracene	0.37	U
108-95-2	Phenol	0.37	U	50-32-8	Benzo[a]pyrene	0.37	U
129-00-0	Pyrene	0.37	U				

Worksheet #: 57379

Total Target Concentration 0.17

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-012
Client Id: CPC-SB06-1.5-2.5
Data File: 9M05053.D
Analysis Date: 08/23/07 01:56
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 93

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.36	U	205-99-2	Benzo[b]fluoranthene	0.36	U
95-95-4	2,4,5-Trichlorophenol	0.36	U	191-24-2	Benzo[g,h,i]perylene	0.36	U
88-06-2	2,4,6-Trichlorophenol	0.36	U	207-08-9	Benzo[k]fluoranthene	0.36	U
120-83-2	2,4-Dichlorophenol	0.36	U	111-91-1	bis(2-Chloroethoxy)methan	0.36	U
105-67-9	2,4-Dimethylphenol	0.36	U	111-44-4	bis(2-Chloroethyl)ether	0.36	U
51-28-5	2,4-Dinitrophenol	0.36	U	39638-32-9	bis(2-chloroisopropyl)ether	0.36	U
121-14-2	2,4-Dinitrotoluene	0.36	U	117-81-7	bis(2-Ethylhexyl)phthalate	0.36	0.13 J
606-20-2	2,6-Dinitrotoluene	0.36	U	85-68-7	Butylbenzylphthalate	0.36	U
91-58-7	2-Chloronaphthalene	0.36	U	105-60-2	Caprolactam	0.36	U
95-57-8	2-Chlorophenol	0.36	U	86-74-8	Carbazole	0.36	U
91-57-6	2-Methylnaphthalene	0.36	U	218-01-9	Chrysene	0.36	U
95-48-7	2-Methylphenol	0.36	U	53-70-3	Dibenzo[a,h]anthracene	0.36	U
88-74-4	2-Nitroaniline	0.36	U	132-64-9	Dibenzofuran	0.36	U
88-75-5	2-Nitrophenol	0.36	U	84-66-2	Diethylphthalate	0.36	U
106-44-5	3&4-Methylphenol	0.36	U	131-11-3	Dimethylphthalate	0.36	U
91-94-1	3,3'-Dichlorobenzidine	0.36	U	84-74-2	Di-n-butylphthalate	0.36	U
99-09-2	3-Nitroaniline	0.36	U	117-84-0	Di-n-octylphthalate	0.36	U
534-52-1	4,6-Dinitro-2-methylphenol	0.36	U	206-44-0	Fluoranthene	0.36	U
101-55-3	4-Bromophenyl-phenylether	0.36	U	86-73-7	Fluorene	0.36	U
59-50-7	4-Chloro-3-methylphenol	0.36	U	118-74-1	Hexachlorobenzene	0.36	U
106-47-8	4-Chloroaniline	0.36	U	87-68-3	Hexachlorobutadiene	0.36	U
7005-72-3	4-Chlorophenyl-phenylether	0.36	U	77-47-4	Hexachlorocyclopentadiene	0.36	U
100-01-6	4-Nitroaniline	0.36	U	67-72-1	Hexachloroethane	0.36	U
100-02-7	4-Nitrophenol	0.36	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.36	U
83-32-9	Acenaphthene	0.36	U	78-59-1	Isophorone	0.36	U
208-96-8	Acenaphthylene	0.36	U	91-20-3	Naphthalene	0.36	U
98-86-2	Acetophenone	0.36	U	98-95-3	Nitrobenzene	0.36	U
621-64-7	N-Nitroso-di-n-propylamine	0.36	U	120-12-7	Anthracene	0.36	U
86-30-6	n-Nitrosodiphenylamine	0.36	U	1912-24-9	Atrazine	0.36	U
87-86-5	Pentachlorophenol	0.36	U	100-52-7	Benzaldehyde	0.36	U
85-01-8	Phenanthrene	0.36	U	56-55-3	Benzo[a]anthracene	0.36	U
108-95-2	Phenol	0.36	U	50-32-8	Benzo[a]pyrene	0.36	U
129-00-0	Pyrene	0.36	U				

Worksheet #: 57379

Total Target Concentration 0.13

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-013
Client Id: CPC-SB05-1-3
Data File: 9M05054.D
Analysis Date: 08/23/07 02:20
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 94

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.35	U	205-99-2	Benzo[b]fluoranthene	0.35	U
95-95-4	2,4,5-Trichlorophenol	0.35	U	191-24-2	Benzo[g,h,i]perylene	0.35	U
88-06-2	2,4,6-Trichlorophenol	0.35	U	207-08-9	Benzo[k]fluoranthene	0.35	U
120-83-2	2,4-Dichlorophenol	0.35	U	111-91-1	bis(2-Chloroethoxy)methan	0.35	U
105-67-9	2,4-Dimethylphenol	0.35	U	111-44-4	bis(2-Chloroethyl)ether	0.35	U
51-28-5	2,4-Dinitrophenol	0.35	U	39638-32-9	bis(2-chloroisopropyl)ether	0.35	U
121-14-2	2,4-Dinitrotoluene	0.35	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.35**	**0.11 J**
606-20-2	2,6-Dinitrotoluene	0.35	U	85-68-7	Butylbenzylphthalate	0.35	U
91-58-7	2-Chloronaphthalene	0.35	U	105-60-2	Caprolactam	0.35	U
95-57-8	2-Chlorophenol	0.35	U	86-74-8	Carbazole	0.35	U
91-57-6	2-Methylnaphthalene	0.35	U	218-01-9	Chrysene	0.35	U
95-48-7	2-Methylphenol	0.35	U	53-70-3	Dibenzo[a,h]anthracene	0.35	U
88-74-4	2-Nitroaniline	0.35	U	132-64-9	Dibenzofuran	0.35	U
88-75-5	2-Nitrophenol	0.35	U	84-66-2	Diethylphthalate	0.35	U
106-44-5	3&4-Methylphenol	0.35	U	131-11-3	Dimethylphthalate	0.35	U
91-94-1	3,3'-Dichlorobenzidine	0.35	U	84-74-2	Di-n-butylphthalate	0.35	U
99-09-2	3-Nitroaniline	0.35	U	117-84-0	Di-n-octylphthalate	0.35	U
534-52-1	4,6-Dinitro-2-methylphenol	0.35	U	206-44-0	Fluoranthene	0.35	U
101-55-3	4-Bromophenyl-phenylether	0.35	U	86-73-7	Fluorene	0.35	U
59-50-7	4-Chloro-3-methylphenol	0.35	U	118-74-1	Hexachlorobenzene	0.35	U
106-47-8	4-Chloroaniline	0.35	U	87-68-3	Hexachlorobutadiene	0.35	U
7005-72-3	4-Chlorophenyl-phenylether	0.35	U	77-47-4	Hexachlorocyclopentadiene	0.35	U
100-01-6	4-Nitroaniline	0.35	U	67-72-1	Hexachloroethane	0.35	U
100-02-7	4-Nitrophenol	0.35	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.35	U
83-32-9	Acenaphthene	0.35	U	78-59-1	Isophorone	0.35	U
208-96-8	Acenaphthylene	0.35	U	91-20-3	Naphthalene	0.35	U
98-86-2	Acetophenone	0.35	U	98-95-3	Nitrobenzene	0.35	U
62-53-3	Aniline	0.35	U	621-64-7	N-Nitroso-di-n-propylamine	0.35	U
120-12-7	Anthracene	0.35	U	86-30-6	n-Nitrosodiphenylamine	0.35	U
87-86-5	Pentachlorophenol	0.35	U	100-52-7	Benzaldehyde	0.35	U
85-01-8	Phenanthrene	0.35	U	56-55-3	Benzo[a]anthracene	0.35	U
108-95-2	Phenol	0.35	U	50-32-8	Benzo[a]pyrene	0.35	U
129-00-0	Pyrene	0.35	U				

Worksheet #: 57379

Total Target Concentration 0.11

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-014
Client Id: CPC-SB05-3-5
Data File: 9M05091.D
Analysis Date: 08/23/07 23:04
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 96

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.35	U	**205-99-2**	**Benzo[b]fluoranthene**	**0.35**	**0.071 J**
95-95-4	2,4,5-Trichlorophenol	0.35	U	**191-24-2**	**Benzo[g,h,i]perylene**	**0.35**	**0.041 J**
88-06-2	2,4,6-Trichlorophenol	0.35	U	207-08-9	Benzo[k]fluoranthene	0.35	U
120-83-2	2,4-Dichlorophenol	0.35	U	111-91-1	bis(2-Chloroethoxy)methan	0.35	U
105-67-9	2,4-Dimethylphenol	0.35	U	111-44-4	bis(2-Chloroethyl)ether	0.35	U
51-28-5	2,4-Dinitrophenol	0.35	U	39638-32-9	bis(2-chloroisopropyl)ether	0.35	U
121-14-2	2,4-Dinitrotoluene	0.35	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.35**	**0.14 J**
606-20-2	2,6-Dinitrotoluene	0.35	U	85-68-7	Butylbenzylphthalate	0.35	U
91-58-7	2-Chloronaphthalene	0.35	U	105-60-2	Caprolactam	0.35	U
95-57-8	2-Chlorophenol	0.35	U	86-74-8	Carbazole	0.35	U
91-57-6	2-Methylnaphthalene	0.35	U	**218-01-9**	**Chrysene**	**0.35**	**0.068 J**
95-48-7	2-Methylphenol	0.35	U	53-70-3	Dibenzo[a,h]anthracene	0.35	U
88-74-4	2-Nitroaniline	0.35	U	132-64-9	Dibenzofuran	0.35	U
88-75-5	2-Nitrophenol	0.35	U	84-66-2	Diethylphthalate	0.35	U
106-44-5	3&4-Methylphenol	0.35	U	131-11-3	Dimethylphthalate	0.35	U
91-94-1	3,3'-Dichlorobenzidine	0.35	U	84-74-2	Di-n-butylphthalate	0.35	U
99-09-2	3-Nitroaniline	0.35	U	117-84-0	Di-n-octylphthalate	0.35	U
534-52-1	4,6-Dinitro-2-methylphenol	0.35	U	**206-44-0**	**Fluoranthene**	**0.35**	**0.12 J**
101-55-3	4-Bromophenyl-phenylether	0.35	U	86-73-7	Fluorene	0.35	U
59-50-7	4-Chloro-3-methylphenol	0.35	U	118-74-1	Hexachlorobenzene	0.35	U
106-47-8	4-Chloroaniline	0.35	U	87-68-3	Hexachlorobutadiene	0.35	U
7005-72-3	4-Chlorophenyl-phenylether	0.35	U	77-47-4	Hexachlorocyclopentadiene	0.35	U
100-01-6	4-Nitroaniline	0.35	U	67-72-1	Hexachloroethane	0.35	U
100-02-7	4-Nitrophenol	0.35	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.35	U
83-32-9	Acenaphthene	0.35	U	78-59-1	Isophorone	0.35	U
208-96-8	Acenaphthylene	0.35	U	91-20-3	Naphthalene	0.35	U
98-86-2	Acetophenone	0.35	U	98-95-3	Nitrobenzene	0.35	U
621-64-7	N-Nitroso-di-n-propylamine	0.35	U	120-12-7	Anthracene	0.35	U
86-30-6	n-Nitrosodiphenylamine	0.35	U	1912-24-9	Atrazine	0.35	U
87-86-5	Pentachlorophenol	0.35	U	100-52-7	Benzaldehyde	0.35	U
85-01-8	**Phenanthrene**	**0.35**	**0.070 J**	**56-55-3**	**Benzo[a]anthracene**	**0.35**	**0.059 J**
108-95-2	Phenol	0.35	U	**50-32-8**	**Benzo[a]pyrene**	**0.35**	**0.049 J**
129-00-0	**Pyrene**	**0.35**	**0.13 J**				

Worksheet #: 57379

Total Target Concentration 0.748

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-015
Client Id: CPC-SB05-5-7
Data File: 9M05058.D
Analysis Date: 08/23/07 03:55
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 98

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.34	U	205-99-2	Benzo[b]fluoranthene	0.34	U
95-95-4	2,4,5-Trichlorophenol	0.34	U	191-24-2	Benzo[g,h,i]perylene	0.34	U
88-06-2	2,4,6-Trichlorophenol	0.34	U	207-08-9	Benzo[k]fluoranthene	0.34	U
120-83-2	2,4-Dichlorophenol	0.34	U	111-91-1	bis(2-Chloroethoxy)methan	0.34	U
105-67-9	2,4-Dimethylphenol	0.34	U	111-44-4	bis(2-Chloroethyl)ether	0.34	U
51-28-5	2,4-Dinitrophenol	0.34	U	39638-32-9	bis(2-chloroisopropyl)ether	0.34	U
121-14-2	2,4-Dinitrotoluene	0.34	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.34**	**0.060 J**
606-20-2	2,6-Dinitrotoluene	0.34	U	85-68-7	Butylbenzylphthalate	0.34	U
91-58-7	2-Chloronaphthalene	0.34	U	105-60-2	Caprolactam	0.34	U
95-57-8	2-Chlorophenol	0.34	U	86-74-8	Carbazole	0.34	U
91-57-6	2-Methylnaphthalene	0.34	U	218-01-9	Chrysene	0.34	U
95-48-7	2-Methylphenol	0.34	U	53-70-3	Dibenzo[a,h]anthracene	0.34	U
88-74-4	2-Nitroaniline	0.34	U	132-64-9	Dibenzofuran	0.34	U
88-75-5	2-Nitrophenol	0.34	U	84-66-2	Diethylphthalate	0.34	U
106-44-5	3&4-Methylphenol	0.34	U	131-11-3	Dimethylphthalate	0.34	U
91-94-1	3,3'-Dichlorobenzidine	0.34	U	84-74-2	Di-n-butylphthalate	0.34	U
99-09-2	3-Nitroaniline	0.34	U	117-84-0	Di-n-octylphthalate	0.34	U
534-52-1	4,6-Dinitro-2-methylphenol	0.34	U	206-44-0	Fluoranthene	0.34	U
101-55-3	4-Bromophenyl-phenylether	0.34	U	86-73-7	Fluorene	0.34	U
59-50-7	4-Chloro-3-methylphenol	0.34	U	118-74-1	Hexachlorobenzene	0.34	U
106-47-8	4-Chloroaniline	0.34	U	87-68-3	Hexachlorobutadiene	0.34	U
7005-72-3	4-Chlorophenyl-phenylether	0.34	U	77-47-4	Hexachlorocyclopentadiene	0.34	U
100-01-6	4-Nitroaniline	0.34	U	67-72-1	Hexachloroethane	0.34	U
100-02-7	4-Nitrophenol	0.34	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.34	U
83-32-9	Acenaphthene	0.34	U	78-59-1	Isophorone	0.34	U
208-96-8	Acenaphthylene	0.34	U	91-20-3	Naphthalene	0.34	U
98-86-2	Acetophenone	0.34	U	98-95-3	Nitrobenzene	0.34	U
621-64-7	N-Nitroso-di-n-propylamine	0.34	U	120-12-7	Anthracene	0.34	U
86-30-6	n-Nitrosodiphenylamine	0.34	U	1912-24-9	Atrazine	0.34	U
87-86-5	Pentachlorophenol	0.34	U	100-52-7	Benzaldehyde	0.34	U
85-01-8	Phenanthrene	0.34	U	56-55-3	Benzo[a]anthracene	0.34	U
108-95-2	Phenol	0.34	U	50-32-8	Benzo[a]pyrene	0.34	U
129-00-0	Pyrene	0.34	U				

Worksheet #: 57379

Total Target Concentration 0.06

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-016
Client Id: CPC-SB05-7-9
Data File: 5M34187.D
Analysis Date: 08/22/07 21:58
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 98

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.34	U	205-99-2	Benzo[b]fluoranthene	0.34	U
95-95-4	2,4,5-Trichlorophenol	0.34	U	191-24-2	Benzo[g,h,i]perylene	0.34	U
88-06-2	2,4,6-Trichlorophenol	0.34	U	207-08-9	Benzo[k]fluoranthene	0.34	U
120-83-2	2,4-Dichlorophenol	0.34	U	111-91-1	bis(2-Chloroethoxy)methan	0.34	U
105-67-9	2,4-Dimethylphenol	0.34	U	111-44-4	bis(2-Chloroethyl)ether	0.34	U
51-28-5	2,4-Dinitrophenol	0.34	U	39638-32-9	bis(2-chloroisopropyl)ether	0.34	U
121-14-2	2,4-Dinitrotoluene	0.34	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.34**	**0.051 J**
606-20-2	2,6-Dinitrotoluene	0.34	U	85-68-7	Butylbenzylphthalate	0.34	U
91-58-7	2-Chloronaphthalene	0.34	U	105-60-2	Caprolactam	0.34	U
95-57-8	2-Chlorophenol	0.34	U	86-74-8	Carbazole	0.34	U
91-57-6	2-Methylnaphthalene	0.34	U	218-01-9	Chrysene	0.34	U
95-48-7	2-Methylphenol	0.34	U	53-70-3	Dibenzo[a,h]anthracene	0.34	U
88-74-4	2-Nitroaniline	0.34	U	132-64-9	Dibenzofuran	0.34	U
88-75-5	2-Nitrophenol	0.34	U	84-66-2	Diethylphthalate	0.34	U
106-44-5	3&4-Methylphenol	0.34	U	131-11-3	Dimethylphthalate	0.34	U
91-94-1	3,3'-Dichlorobenzidine	0.34	U	84-74-2	Di-n-butylphthalate	0.34	U
99-09-2	3-Nitroaniline	0.34	U	117-84-0	Di-n-octylphthalate	0.34	U
534-52-1	4,6-Dinitro-2-methylphenol	0.34	U	206-44-0	Fluoranthene	0.34	U
101-55-3	4-Bromophenyl-phenylether	0.34	U	86-73-7	Fluorene	0.34	U
59-50-7	4-Chloro-3-methylphenol	0.34	U	118-74-1	Hexachlorobenzene	0.34	U
106-47-8	4-Chloroaniline	0.34	U	87-68-3	Hexachlorobutadiene	0.34	U
7005-72-3	4-Chlorophenyl-phenylether	0.34	U	77-47-4	Hexachlorocyclopentadiene	0.85	U
100-01-6	4-Nitroaniline	0.34	U	67-72-1	Hexachloroethane	0.34	U
100-02-7	4-Nitrophenol	0.34	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.34	U
83-32-9	Acenaphthene	0.34	U	78-59-1	Isophorone	0.34	U
208-96-8	Acenaphthylene	0.34	U	91-20-3	Naphthalene	0.34	U
98-86-2	Acetophenone	0.34	U	98-95-3	Nitrobenzene	0.34	U
621-64-7	N-Nitroso-di-n-propylamine	0.34	U	120-12-7	Anthracene	0.34	U
86-30-6	n-Nitrosodiphenylamine	0.34	U	1912-24-9	Atrazine	0.34	U
87-86-5	Pentachlorophenol	0.34	U	100-52-7	Benzaldehyde	0.34	U
85-01-8	Phenanthrene	0.34	U	56-55-3	Benzo[a]anthracene	0.34	U
108-95-2	Phenol	0.34	U	50-32-8	Benzo[a]pyrene	0.34	U
129-00-0	Pyrene	0.34	U				

Worksheet #: 57379

Total Target Concentration 0.051

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-017
Client Id: CPC-SB02-5-7
Data File: 9M05055.D
Analysis Date: 08/23/07 02:44
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 96

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.35	U	205-99-2	Benzo[b]fluoranthene	0.35	U
95-95-4	2,4,5-Trichlorophenol	0.35	U	191-24-2	Benzo[g,h,i]perylene	0.35	U
88-06-2	2,4,6-Trichlorophenol	0.35	U	207-08-9	Benzo[k]fluoranthene	0.35	U
120-83-2	2,4-Dichlorophenol	0.35	U	111-91-1	bis(2-Chloroethoxy)methan	0.35	U
105-67-9	2,4-Dimethylphenol	0.35	U	111-44-4	bis(2-Chloroethyl)ether	0.35	U
51-28-5	2,4-Dinitrophenol	0.35	U	39638-32-9	bis(2-chloroisopropyl)ether	0.35	U
121-14-2	2,4-Dinitrotoluene	0.35	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.35**	**0.55**
606-20-2	2,6-Dinitrotoluene	0.35	U	85-68-7	Butylbenzylphthalate	0.35	U
91-58-7	2-Chloronaphthalene	0.35	U	105-60-2	Caprolactam	0.35	U
95-57-8	2-Chlorophenol	0.35	U	86-74-8	Carbazole	0.35	U
91-57-6	2-Methylnaphthalene	0.35	U	218-01-9	Chrysene	0.35	U
95-48-7	2-Methylphenol	0.35	U	53-70-3	Dibenzo[a,h]anthracene	0.35	U
88-74-4	2-Nitroaniline	0.35	U	132-64-9	Dibenzofuran	0.35	U
88-75-5	2-Nitrophenol	0.35	U	84-66-2	Diethylphthalate	0.35	U
106-44-5	3&4-Methylphenol	0.35	U	131-11-3	Dimethylphthalate	0.35	U
91-94-1	3,3'-Dichlorobenzidine	0.35	U	84-74-2	Di-n-butylphthalate	0.35	U
99-09-2	3-Nitroaniline	0.35	U	117-84-0	Di-n-octylphthalate	0.35	U
534-52-1	4,6-Dinitro-2-methylphenol	0.35	U	206-44-0	Fluoranthene	0.35	U
101-55-3	4-Bromophenyl-phenylether	0.35	U	86-73-7	Fluorene	0.35	U
59-50-7	4-Chloro-3-methylphenol	0.35	U	118-74-1	Hexachlorobenzene	0.35	U
106-47-8	4-Chloroaniline	0.35	U	87-68-3	Hexachlorobutadiene	0.35	U
7005-72-3	4-Chlorophenyl-phenylether	0.35	U	77-47-4	Hexachlorocyclopentadiene	0.35	U
100-01-6	4-Nitroaniline	0.35	U	67-72-1	Hexachloroethane	0.35	U
100-02-7	4-Nitrophenol	0.35	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.35	U
83-32-9	Acenaphthene	0.35	U	78-59-1	Isophorone	0.35	U
208-96-8	Acenaphthylene	0.35	U	91-20-3	Naphthalene	0.35	U
98-86-2	Acetophenone	0.35	U	98-95-3	Nitrobenzene	0.35	U
621-64-7	N-Nitroso-di-n-propylamine	0.35	U	120-12-7	Anthracene	0.35	U
86-30-6	n-Nitrosodiphenylamine	0.35	U	1912-24-9	Atrazine	0.35	U
87-86-5	Pentachlorophenol	0.35	U	100-52-7	Benzaldehyde	0.35	U
85-01-8	Phenanthrene	0.35	U	56-55-3	Benzo[a]anthracene	0.35	U
108-95-2	Phenol	0.35	U	50-32-8	Benzo[a]pyrene	0.35	U
129-00-0	Pyrene	0.35	U				

Worksheet #: 57379

Total Target Concentration 0.55

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-018
Client Id: CPC-SB02-7-9
Data File: 9M05056.D
Analysis Date: 08/23/07 03:08
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 97

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.34	U	205-99-2	Benzo[b]fluoranthene	0.34	U
95-95-4	2,4,5-Trichlorophenol	0.34	U	191-24-2	Benzo[g,h,i]perylene	0.34	U
88-06-2	2,4,6-Trichlorophenol	0.34	U	207-08-9	Benzo[k]fluoranthene	0.34	U
120-83-2	2,4-Dichlorophenol	0.34	U	111-91-1	bis(2-Chloroethoxy)methan	0.34	U
105-67-9	2,4-Dimethylphenol	0.34	U	111-44-4	bis(2-Chloroethyl)ether	0.34	U
51-28-5	2,4-Dinitrophenol	0.34	U	39638-32-9	bis(2-chloroisopropyl)ether	0.34	U
121-14-2	2,4-Dinitrotoluene	0.34	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.34**	**0.067 J**
606-20-2	2,6-Dinitrotoluene	0.34	U	85-68-7	Butylbenzylphthalate	0.34	U
91-58-7	2-Chloronaphthalene	0.34	U	105-60-2	Caprolactam	0.34	U
95-57-8	2-Chlorophenol	0.34	U	86-74-8	Carbazole	0.34	U
91-57-6	2-Methylnaphthalene	0.34	U	218-01-9	Chrysene	0.34	U
95-48-7	2-Methylphenol	0.34	U	53-70-3	Dibenzo[a,h]anthracene	0.34	U
88-74-4	2-Nitroaniline	0.34	U	132-64-9	Dibenzofuran	0.34	U
88-75-5	2-Nitrophenol	0.34	U	84-66-2	Diethylphthalate	0.34	U
106-44-5	3&4-Methylphenol	0.34	U	131-11-3	Dimethylphthalate	0.34	U
91-94-1	3,3'-Dichlorobenzidine	0.34	U	84-74-2	Di-n-butylphthalate	0.34	U
99-09-2	3-Nitroaniline	0.34	U	117-84-0	Di-n-octylphthalate	0.34	U
534-52-1	4,6-Dinitro-2-methylphenol	0.34	U	206-44-0	Fluoranthene	0.34	U
101-55-3	4-Bromophenyl-phenylether	0.34	U	86-73-7	Fluorene	0.34	U
59-50-7	4-Chloro-3-methylphenol	0.34	U	118-74-1	Hexachlorobenzene	0.34	U
106-47-8	4-Chloroaniline	0.34	U	87-68-3	Hexachlorobutadiene	0.34	U
7005-72-3	4-Chlorophenyl-phenylether	0.34	U	77-47-4	Hexachlorocyclopentadiene	0.34	U
100-01-6	4-Nitroaniline	0.34	U	67-72-1	Hexachloroethane	0.34	U
100-02-7	4-Nitrophenol	0.34	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.34	U
83-32-9	Acenaphthene	0.34	U	78-59-1	Isophorone	0.34	U
208-96-8	Acenaphthylene	0.34	U	91-20-3	Naphthalene	0.34	U
98-86-2	Acetophenone	0.34	U	98-95-3	Nitrobenzene	0.34	U
621-64-7	N-Nitroso-di-n-propylamine	0.34	U	120-12-7	Anthracene	0.34	U
86-30-6	n-Nitrosodiphenylamine	0.34	U	1912-24-9	Atrazine	0.34	U
87-86-5	Pentachlorophenol	0.34	U	100-52-7	Benzaldehyde	0.34	U
85-01-8	Phenanthrene	0.34	U	56-55-3	Benzo[a]anthracene	0.34	U
108-95-2	Phenol	0.34	U	50-32-8	Benzo[a]pyrene	0.34	U
129-00-0	Pyrene	0.34	U				

Worksheet #: 57379

Total Target Concentration 0.067

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-019
Client Id: CPC-DW04-13-14
Data File: 9M05057.D
Analysis Date: 08/23/07 03:31
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 30g
Final Vol: 1ml
Dilution: 1
Solids: 82

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	0.41	U	205-99-2	Benzo[b]fluoranthene	0.41	U
95-95-4	2,4,5-Trichlorophenol	0.41	U	191-24-2	Benzo[g,h,i]perylene	0.41	U
88-06-2	2,4,6-Trichlorophenol	0.41	U	207-08-9	Benzo[k]fluoranthene	0.41	U
120-83-2	2,4-Dichlorophenol	0.41	U	111-91-1	bis(2-Chloroethoxy)methan	0.41	U
105-67-9	2,4-Dimethylphenol	0.41	U	111-44-4	bis(2-Chloroethyl)ether	0.41	U
51-28-5	2,4-Dinitrophenol	0.41	U	39638-32-9	bis(2-chloroisopropyl)ether	0.41	U
121-14-2	2,4-Dinitrotoluene	0.41	U	**117-81-7**	**bis(2-Ethylhexyl)phthalate**	**0.41**	**0.053 J**
606-20-2	2,6-Dinitrotoluene	0.41	U	85-68-7	Butylbenzylphthalate	0.41	U
91-58-7	2-Chloronaphthalene	0.41	U	105-60-2	Caprolactam	0.41	U
95-57-8	2-Chlorophenol	0.41	U	86-74-8	Carbazole	0.41	U
91-57-6	2-Methylnaphthalene	0.41	U	218-01-9	Chrysene	0.41	U
95-48-7	2-Methylphenol	0.41	U	53-70-3	Dibenzo[a,h]anthracene	0.41	U
88-74-4	2-Nitroaniline	0.41	U	132-64-9	Dibenzofuran	0.41	U
88-75-5	2-Nitrophenol	0.41	U	84-66-2	Diethylphthalate	0.41	U
106-44-5	3&4-Methylphenol	0.41	U	131-11-3	Dimethylphthalate	0.41	U
91-94-1	3,3'-Dichlorobenzidine	0.41	U	84-74-2	Di-n-butylphthalate	0.41	U
99-09-2	3-Nitroaniline	0.41	U	117-84-0	Di-n-octylphthalate	0.41	U
534-52-1	4,6-Dinitro-2-methylphenol	0.41	U	206-44-0	Fluoranthene	0.41	U
101-55-3	4-Bromophenyl-phenylether	0.41	U	86-73-7	Fluorene	0.41	U
59-50-7	4-Chloro-3-methylphenol	0.41	U	118-74-1	Hexachlorobenzene	0.41	U
106-47-8	4-Chloroaniline	0.41	U	87-68-3	Hexachlorobutadiene	0.41	U
7005-72-3	4-Chlorophenyl-phenylether	0.41	U	77-47-4	Hexachlorocyclopentadiene	0.41	U
100-01-6	4-Nitroaniline	0.41	U	67-72-1	Hexachloroethane	0.41	U
100-02-7	4-Nitrophenol	0.41	U	193-39-5	Indeno[1,2,3-cd]pyrene	0.41	U
83-32-9	Acenaphthene	0.41	U	78-59-1	Isophorone	0.41	U
208-96-8	Acenaphthylene	0.41	U	91-20-3	Naphthalene	0.41	U
98-86-2	Acetophenone	0.41	U	98-95-3	Nitrobenzene	0.41	U
621-64-7	N-Nitroso-di-n-propylamine	0.41	U	120-12-7	Anthracene	0.41	U
86-30-6	n-Nitrosodiphenylamine	0.41	U	1912-24-9	Atrazine	0.41	U
87-86-5	Pentachlorophenol	0.41	U	100-52-7	Benzaldehyde	0.41	U
85-01-8	Phenanthrene	0.41	U	56-55-3	Benzo[a]anthracene	0.41	U
108-95-2	Phenol	0.41	U	50-32-8	Benzo[a]pyrene	0.41	U
129-00-0	Pyrene	0.41	U				

Worksheet #: 57379

Total Target Concentration 0.053

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC32378-020
Client Id: CPC-Rinsate
Data File: 5M34128.D
Analysis Date: 08/21/07 15:02
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 1ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
92-52-4	1,1'-Biphenyl	10	U	205-99-2	Benzo[b]fluoranthene	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	207-08-9	Benzo[k]fluoranthene	10	U
120-83-2	2,4-Dichlorophenol	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
105-67-9	2,4-Dimethylphenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
51-28-5	2,4-Dinitrophenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
121-14-2	2,4-Dinitrotoluene	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
606-20-2	2,6-Dinitrotoluene	10	U	85-68-7	Butylbenzylphthalate	10	U
91-58-7	2-Chloronaphthalene	10	U	105-60-2	Caprolactam	10	U
95-57-8	2-Chlorophenol	10	U	86-74-8	Carbazole	10	U
91-57-6	2-Methylnaphthalene	10	U	218-01-9	Chrysene	10	U
95-48-7	2-Methylphenol	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
88-74-4	2-Nitroaniline	10	U	132-64-9	Dibenzofuran	10	U
88-75-5	2-Nitrophenol	10	U	84-66-2	Diethylphthalate	10	U
106-44-5	3&4-Methylphenol	10	U	131-11-3	Dimethylphthalate	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	84-74-2	Di-n-butylphthalate	10	U
99-09-2	3-Nitroaniline	10	U	117-84-0	Di-n-octylphthalate	10	U
534-52-1	4,6-Dinitro-2-methylphenol	10	U	206-44-0	Fluoranthene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	86-73-7	Fluorene	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	118-74-1	Hexachlorobenzene	10	U
106-47-8	4-Chloroaniline	10	U	87-68-3	Hexachlorobutadiene	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	77-47-4	Hexachlorocyclopentadiene	25	U
100-01-6	4-Nitroaniline	10	U	67-72-1	Hexachloroethane	10	U
100-02-7	4-Nitrophenol	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
83-32-9	Acenaphthene	10	U	78-59-1	Isophorone	10	U
208-96-8	Acenaphthylene	10	U	91-20-3	Naphthalene	10	U
98-86-2	Acetophenone	10	U	98-95-3	Nitrobenzene	10	U
621-64-7	N-Nitroso-di-n-propylamine	10	U	120-12-7	Anthracene	10	U
86-30-6	n-Nitrosodiphenylamine	10	U	1912-24-9	Atrazine	10	U
87-86-5	Pentachlorophenol	10	U	100-52-7	Benzaldehyde	10	U
85-01-8	Phenanthrene	10	U	56-55-3	Benzo[a]anthracene	10	U
108-95-2	Phenol	10	U	50-32-8	Benzo[a]pyrene	10	U
129-00-0	Pyrene	10	U				

Worksheet #: 57379

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-001
Client Id: CPC-MW01
Data File: 5G14260.D
Analysis Date: 08/17/07 08:59
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.050	U	7421-93-4	Endrin Aldehyde	0.050	U J
319-84-6	alpha-BHC	0.050	U	53494-70-5	Endrin Ketone	0.050	U J
319-85-7	beta-BHC	0.050	U	58-89-9	gamma-BHC	0.050	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.050	U
319-86-8	delta-BHC	0.050	U J	1024-57-3	Heptachlor Epoxide	0.050	U
60-57-1	Dieldrin	0.050	U	72-43-5	Methoxychlor	0.050	U J
959-98-8	Endosulfan I	0.050	U	72-54-8	p,p'-DDD	0.025	U J
33213-65-9	Endosulfan II	0.050	U	72-55-9	p,p'-DDE	0.025	U J
1031-07-8	Endosulfan Sulfate	0.050	U J	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.050	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-002
Client Id: CPC-MW08
Data File: 5G14268.D
Analysis Date: 08/17/07 11:25
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.051	U	7421-93-4	Endrin Aldehyde	0.051	U J
319-84-6	alpha-BHC	0.051	U	53494-70-5	Endrin Ketone	0.051	U J
319-85-7	beta-BHC	0.051	U	58-89-9	gamma-BHC	0.051	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.051	U
319-86-8	delta-BHC	0.051	U J	1024-57-3	Heptachlor Epoxide	0.051	U
60-57-1	Dieldrin	0.051	U	72-43-5	Methoxychlor	0.051	U J
959-98-8	Endosulfan I	0.051	U	72-54-8	p,p'-DDD	0.025	U J
33213-65-9	Endosulfan II	0.051	U	72-55-9	p,p'-DDE	0.025	U J
1031-07-8	Endosulfan Sulfate	0.051	U J	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.051	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-003
Client Id: CPC-MW03
Data File: 5G14261.D
Analysis Date: 08/17/07 09:17
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.050	U	7421-93-4	Endrin Aldehyde	0.050	U J
319-84-6	alpha-BHC	0.050	U	53494-70-5	Endrin Ketone	0.050	U J
319-85-7	beta-BHC	0.050	U	58-89-9	gamma-BHC	0.050	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.050	U
319-86-8	delta-BHC	0.050	U J	1024-57-3	Heptachlor Epoxide	0.050	U
60-57-1	Dieldrin	0.050	U	72-43-5	Methoxychlor	0.050	U J
959-98-8	Endosulfan I	0.050	U	72-54-8	p,p'-DDD	0.025	U J
33213-65-9	Endosulfan II	0.050	U	72-55-9	p,p'-DDE	0.025	U J
1031-07-8	Endosulfan Sulfate	0.050	U J	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.050	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-004
Client Id: CPC-MW04
Data File: 5G14262.D
Analysis Date: 08/17/07 09:36
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.051	U	7421-93-4	Endrin Aldehyde	0.051	U UJ
319-84-6	alpha-BHC	0.051	U	53494-70-5	Endrin Ketone	0.051	U UJ
319-85-7	beta-BHC	0.051	U	58-89-9	gamma-BHC	0.051	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.051	U
319-86-8	delta-BHC	0.051	U J	1024-57-3	Heptachlor Epoxide	0.051	U
60-57-1	Dieldrin	0.051	U	72-43-5	Methoxychlor	0.051	U UJ
959-98-8	Endosulfan I	0.051	U	72-54-8	p,p'-DDD	0.025	U UJ
33213-65-9	Endosulfan II	0.051	U	72-55-9	p,p'-DDE	0.025	U UJ
1031-07-8	Endosulfan Sulfate	0.051	U J	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.051	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-005
Client Id: CPC-MW05
Data File: 5G14269.D
Analysis Date: 08/17/07 11:43
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.050	U	7421-93-4	Endrin Aldehyde	0.050	U J
319-84-6	alpha-BHC	0.050	U	53494-70-5	Endrin Ketone	0.050	U J
319-85-7	beta-BHC	0.050	U	58-89-9	gamma-BHC	0.050	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.050	U
319-86-8	delta-BHC	0.050	U J	1024-57-3	Heptachlor Epoxide	0.050	U
60-57-1	Dieldrin	0.050	U	72-43-5	Methoxychlor	0.050	U J
959-98-8	Endosulfan I	0.050	U	72-54-8	p,p'-DDD	0.025	U J
33213-65-9	Endosulfan II	0.050	U	72-55-9	p,p'-DDE	0.025	U J
1031-07-8	Endosulfan Sulfate	0.050	U J	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.050	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-006
Client Id: CPC-MW06
Data File: 5G14270.D
Analysis Date: 08/17/07 12:02
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.051	U	7421-93-4	Endrin Aldehyde	0.051	UJ
319-84-6	alpha-BHC	0.051	U	53494-70-5	Endrin Ketone	0.051	UJ
319-85-7	beta-BHC	0.051	U	58-89-9	gamma-BHC	0.051	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.051	U
319-86-8	delta-BHC	0.051	UJ	1024-57-3	Heptachlor Epoxide	0.051	U
60-57-1	Dieldrin	0.051	U	72-43-5	Methoxychlor	0.051	UJ
959-98-8	Endosulfan I	0.051	U	72-54-8	p,p'-DDD	0.025	UJ
33213-65-9	Endosulfan II	0.051	U	72-55-9	p,p'-DDE	0.025	UJ
1031-07-8	Endosulfan Sulfate	0.051	UJ	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.051	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

" - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-007
Client Id: CPC-MW07
Data File: 5G14271.D
Analysis Date: 08/17/07 12:20
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 980ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.051	U	7421-93-4	Endrin Aldehyde	0.051	U J
319-84-6	alpha-BHC	0.051	U	53494-70-5	Endrin Ketone	0.051	U J
319-85-7	beta-BHC	0.051	U	58-89-9	gamma-BHC	0.051	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.051	U
319-86-8	delta-BHC	0.051	U J	1024-57-3	Heptachlor Epoxide	0.051	U
60-57-1	Dieldrin	0.051	U	72-43-5	Methoxychlor	0.051	U J
959-98-8	Endosulfan I	0.051	U	72-54-8	p,p'-DDD	0.026	U J
33213-65-9	Endosulfan II	0.051	U	72-55-9	p,p'-DDE	0.026	U J
1031-07-8	Endosulfan Sulfate	0.051	U J	50-29-3	p,p'-DDT	0.026	U
72-20-8	Endrin	0.051	U	8001-35-2	Toxaphene	0.26	U

Worksheet #: 55946

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-008
Client Id: CPC-MW09
Data File: 5G14266.D
Analysis Date: 08/17/07 10:49
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.051	U	7421-93-4	Endrin Aldehyde	0.051	U J
319-84-6	alpha-BHC	0.051	U	53494-70-5	Endrin Ketone	0.051	U J
319-85-7	beta-BHC	0.051	U	58-89-9	gamma-BHC	0.051	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.051	U
319-86-8	delta-BHC	0.051	U J	1024-57-3	Heptachlor Epoxide	0.051	U
60-57-1	Dieldrin	0.051	U	72-43-5	Methoxychlor	0.051	U J
959-98-8	Endosulfan I	0.051	U	72-54-8	p,p'-DDD	0.025	U J
33213-65-9	Endosulfan II	0.051	U	72-55-9	p,p'-DDE	0.025	U J
1031-07-8	Endosulfan Sulfate	0.051	U J	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.051	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

" - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-009
Client Id: CPC-MW10
Data File: 5G14263.D
Analysis Date: 08/17/07 09:54
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.050	U	7421-93-4	Endrin Aldehyde	0.050	U J
319-84-6	alpha-BHC	0.050	U	53494-70-5	Endrin Ketone	0.050	U J
319-85-7	beta-BHC	0.050	U	58-89-9	gamma-BHC	0.050	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.050	U
319-86-8	delta-BHC	0.050	U J	1024-57-3	Heptachlor Epoxide	0.050	U
60-57-1	Dieldrin	0.050	U	72-43-5	Methoxychlor	0.050	U J
959-98-8	Endosulfan I	0.050	U	72-54-8	p,p'-DDD	0.025	U J
33213-65-9	Endosulfan II	0.050	U	72-55-9	p,p'-DDE	0.025	U J
1031-07-8	Endosulfan Sulfate	0.050	U J	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.050	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

" - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32351-012
Client Id: CPC-DUPLICATE
Data File: 5G14267.D
Analysis Date: 08/17/07 11:07
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.051	U	7421-93-4	Endrin Aldehyde	0.051	U UJ
319-84-6	alpha-BHC	0.051	U	53494-70-5	Endrin Ketone	0.051	U UJ
319-85-7	beta-BHC	0.051	U	58-89-9	gamma-BHC	0.051	U
57-74-9	Chlordane	0.10	U	76-44-8	Heptachlor	0.051	U
319-86-8	delta-BHC	0.051	U UJ	1024-57-3	Heptachlor Epoxide	0.051	U
60-57-1	Dieldrin	0.051	U	72-43-5	Methoxychlor	0.051	U UJ
959-98-8	Endosulfan I	0.051	U	72-54-8	p,p'-DDD	0.025	U UJ
33213-65-9	Endosulfan II	0.051	U	72-55-9	p,p'-DDE	0.025	U UJ
1031-07-8	Endosulfan Sulfate	0.051	U UJ	50-29-3	p,p'-DDT	0.025	U
72-20-8	Endrin	0.051	U	8001-35-2	Toxaphene	0.25	U

Worksheet #: 55946

Total Target Concentration 0

*** - Indicates the compound was analyzed but not detected.*
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-001
Client Id: CPC-DW01-10-11
Data File: 6G03701.D
Analysis Date: 08/20/07 08:21
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 86

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0058	U	7421-93-4	Endrin Aldehyde	0.0058	U
319-84-6	alpha-BHC	0.0058	U	53494-70-5	Endrin Ketone	0.0058	U UJ
319-85-7	beta-BHC	0.0058	U	58-89-9	gamma-BHC	0.0058	U UJ
57-74-9	Chlordane	0.012	U	76-44-8	Heptachlor	0.0058	U UJ
319-86-8	delta-BHC	0.0058	U UJ	1024-57-3	Heptachlor Epoxide	0.0058	U
60-57-1	Dieldrin	0.0058	U	72-43-5	Methoxychlor	0.0058	U UJ
959-98-8	Endosulfan I	0.0058	U UJ	72-54-8	p,p'-DDD	0.0029	U
33213-65-9	Endosulfan II	0.0058	U	72-55-9	p,p'-DDE	0.0029	U
1031-07-8	Endosulfan Sulfate	0.0058	U UJ	50-29-3	p,p'-DDT	0.0029	U UJ
72-20-8	Endrin	0.0058	U	8001-35-2	Toxaphene	0.029	U

Worksheet #: 56166

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-002
Client Id: CPC-DW02-10-11
Data File: 6G03702.D
Analysis Date: 08/20/07 08:35
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 83

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0060	U	7421-93-4	Endrin Aldehyde	0.0060	U
319-84-6	alpha-BHC	0.0060	U	53494-70-5	Endrin Ketone	0.0060	UJ
319-85-7	beta-BHC	0.0060	U	58-89-9	gamma-BHC	0.0060	UJ
57-74-9	**Chlordane**	**0.012**	**0.021**	76-44-8	Heptachlor	0.0060	UJ
319-86-8	delta-BHC	0.0060	UJ	1024-57-3	Heptachlor Epoxide	0.0060	U
60-57-1	Dieldrin	0.0060	U	72-43-5	Methoxychlor	0.0060	UJ
959-98-8	Endosulfan I	0.0060	UJ	72-54-8	p,p'-DDD	0.0030	U
33213-65-9	Endosulfan II	0.0060	U	72-55-9	p,p'-DDE	0.0030	U
1031-07-8	Endosulfan Sulfate	0.0060	UJ	50-29-3	p,p'-DDT	0.0030	UJ
72-20-8	Endrin	0.0060	U	8001-35-2	Toxaphene	0.030	U

Worksheet #: 56166

Total Target Concentration 0.021

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-003
Client Id: CPC-DW03-10-11
Data File: 5G14287.D
Analysis Date: 08/20/07 07:55
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 90

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0056	U	7421-93-4	Endrin Aldehyde	0.0056	U
319-84-6	alpha-BHC	0.0056	U	53494-70-5	Endrin Ketone	0.0056	U J
319-85-7	beta-BHC	0.0056	U	58-89-9	gamma-BHC	0.0056	U J
57-74-9	Chlordane	0.011	U	76-44-8	Heptachlor	0.0056	U J
319-86-8	delta-BHC	0.0056	U J	1024-57-3	Heptachlor Epoxide	0.0056	U
60-57-1	Dieldrin	0.0056	U	72-43-5	Methoxychlor	0.0056	U J
959-98-8	Endosulfan I	0.0056	U J	72-54-8	p,p'-DDD	0.0028	U
33213-65-9	Endosulfan II	0.0056	U	72-55-9	p,p'-DDE	0.0028	U
1031-07-8	Endosulfan Sulfate	0.0056	U J	50-29-3	p,p'-DDT	0.0028	U J
72-20-8	Endrin	0.0056	U	8001-35-2	Toxaphene	0.028	U

Worksheet #: 56166

Total Target Concentration 0

ndicates the compound was analyzed but not detected.
ndicates the analyte was found in the blank as well as in the sample.
.ndicates the analyte concentration exceeds the calibration range of le instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-004
Client Id: CPC-DW04-8-9
Data File: 6G03703.D
Analysis Date: 08/20/07 08:50
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 85

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0059	U	7421-93-4	Endrin Aldehyde	0.0059	U
319-84-6	alpha-BHC	0.0059	U	53494-70-5	Endrin Ketone	0.0059	U J
319-85-7	beta-BHC	0.0059	U	58-89-9	gamma-BHC	0.0059	U J
57-74-9	**Chlordane**	**0.012**	**0.050**	76-44-8	Heptachlor	0.0059	U J
319-86-8	delta-BHC	0.0059	U J	1024-57-3	Heptachlor Epoxide	0.0059	U
60-57-1	Dieldrin	0.0059	U	72-43-5	Methoxychlor	0.0059	U J
959-98-8	Endosulfan I	0.0059	U J	**72-54-8**	**p,p'-DDD**	**0.0029**	**0.019**
33213-65-9	Endosulfan II	0.0059	U	**72-55-9**	**p,p'-DDE**	**0.0029**	**0.0036 d**
1031-07-8	Endosulfan Sulfate	0.0059	U J	**50-29-3**	**p,p'-DDT**	**0.0029**	**0.075 J**
72-20-8	Endrin	0.0059	U	8001-35-2	Toxaphene	0.029	U

Worksheet #: 56166

Total Target Concentration 0.1476

U - Indicates the compound was analyzed but not detected.
' - Indicates the analyte was found in the blank as well as in the sample.
' - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-005
Client Id: CPC-SOIL-Duplicate
Data File: 6G03704.D
Analysis Date: 08/20/07 09:05
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 92

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0054	U	7421-93-4	Endrin Aldehyde	0.0054	U
319-84-6	alpha-BHC	0.0054	U	53494-70-5	Endrin Ketone	0.0054	UJ
319-85-7	beta-BHC	0.0054	U	58-89-9	gamma-BHC	0.0054	UJ
57-74-9	Chlordane	0.011	U	76-44-8	Heptachlor	0.0054	UJ
319-86-8	delta-BHC	0.0054	UJ	1024-57-3	Heptachlor Epoxide	0.0054	U
60-57-1	Dieldrin	0.0054	U	72-43-5	Methoxychlor	0.0054	UJ
959-98-8	Endosulfan I	0.0054	UJ	72-54-8	p,p'-DDD	0.0027	U
33213-65-9	Endosulfan II	0.0054	U	72-55-9	p,p'-DDE	0.0027	U
1031-07-8	Endosulfan Sulfate	0.0054	UJ	50-29-3	p,p'-DDT	0.0027	UJ
72-20-8	Endrin	0.0054	U	8001-35-2	Toxaphene	0.027	U

Worksheet #: 56166

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
' - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-006
Client Id: CPC-SB01-1.5-2.5
Data File: 5G14281.D
Analysis Date: 08/20/07 06:06
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 95

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0053	U	7421-93-4	Endrin Aldehyde	0.0053	U
319-84-6	alpha-BHC	0.0053	U	53494-70-5	Endrin Ketone	0.0053	U J
319-85-7	beta-BHC	0.0053	U	58-89-9	gamma-BHC	0.0053	U J
57-74-9	Chlordane	0.011	U	76-44-8	Heptachlor	0.0053	U J
319-86-8	delta-BHC	0.0053	U J	1024-57-3	Heptachlor Epoxide	0.0053	U
60-57-1	Dieldrin	0.0053	U	72-43-5	Methoxychlor	0.0053	U J
959-98-8	Endosulfan I	0.0053	U J	72-54-8	p,p'-DDD	0.0026	U
33213-65-9	Endosulfan II	0.0053	U	72-55-9	p,p'-DDE	0.0026	U
1031-07-8	Endosulfan Sulfate	0.0053	U J	50-29-3	p,p'-DDT	0.0026	U J
72-20-8	Endrin	0.0053	U	8001-35-2	Toxaphene	0.026	U

Worksheet #: 56166

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
' - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-007
Client Id: CPC-SB02-1-5
Data File: 5G14282.D
Analysis Date: 08/20/07 06:24
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 94

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0053	U	7421-93-4	Endrin Aldehyde	0.0053	U
319-84-6	alpha-BHC	0.0053	U	53494-70-5	Endrin Ketone	0.0053	U J
319-85-7	beta-BHC	0.0053	U	58-89-9	gamma-BHC	0.0053	U J
57-74-9	Chlordane	0.011	U	76-44-8	Heptachlor	0.0053	U
319-86-8	delta-BHC	0.0053	U J	1024-57-3	Heptachlor Epoxide	0.0053	U J
60-57-1	Dieldrin	0.0053	U	72-43-5	Methoxychlor	0.0053	U J
959-98-8	Endosulfan I	0.0053	U J	72-54-8	p,p'-DDD	0.0027	U
33213-65-9	Endosulfan II	0.0053	U	72-55-9	p,p'-DDE	0.0027	U
1031-07-8	Endosulfan Sulfate	0.0053	U J	50-29-3	p,p'-DDT	0.0027	U J
72-20-8	Endrin	0.0053	U	8001-35-2	Toxaphene	0.027	U

Worksheet #: 56166

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
: - Indicates the analyte was found in the blank as well as in the sample.
' - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-008
Client Id: CPC-SB03-1.5-3.5
Data File: 6G03696.D
Analysis Date: 08/20/07 07:06
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 93

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0054	U	7421-93-4	Endrin Aldehyde	0.0054	U
319-84-6	alpha-BHC	0.0054	U	53494-70-5	Endrin Ketone	0.0054	UJ
319-85-7	beta-BHC	0.0054	U	58-89-9	gamma-BHC	0.0054	UJ
57-74-9	Chlordane	0.011	U	76-44-8	Heptachlor	0.0054	UJ
319-86-8	delta-BHC	0.0054	U	1024-57-3	Heptachlor Epoxide	0.0054	U
60-57-1	Dieldrin	0.0054	U	72-43-5	Methoxychlor	0.0054	UJ
959-98-8	Endosulfan I	0.0054	UJ	72-54-8	p,p'-DDD	0.0027	U
33213-65-9	Endosulfan II	0.0054	U	72-55-9	p,p'-DDE	0.0027	U
1031-07-8	Endosulfan Sulfate	0.0054	UJ	50-29-3	p,p'-DDT	0.0027	UJ
72-20-8	Endrin	0.0054	U	8001-35-2	Toxaphene	0.027	U

Worksheet #: 56166

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
I - Indicates the analyte was found in the blank as well as in the sample.
' - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-011
Client Id: CPC-SB04-1.5-2.5
Data File: 5G14283.D
Analysis Date: 08/20/07 06:42
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 90

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0056	U	7421-93-4	Endrin Aldehyde	0.0056	U
319-84-6	alpha-BHC	0.0056	U	53494-70-5	Endrin Ketone	0.0056	U J
319-85-7	beta-BHC	0.0056	U	58-89-9	gamma-BHC	0.0056	U J
57-74-9	Chlordane	0.011	U	76-44-8	Heptachlor	0.0056	U J
319-86-8	delta-BHC	0.0056	U J	1024-57-3	Heptachlor Epoxide	0.0056	U
60-57-1	Dieldrin	0.0056	U	72-43-5	Methoxychlor	0.0056	U J
959-98-8	Endosulfan I	0.0056	U J	72-54-8	p,p'-DDD	0.0028	U
33213-65-9	Endosulfan II	0.0056	U	72-55-9	p,p'-DDE	0.0028	U
1031-07-8	Endosulfan Sulfate	0.0056	U J	50-29-3	p,p'-DDT	0.0028	U J
72-20-8	Endrin	0.0056	U	8001-35-2	Toxaphene	0.028	U

Worksheet #: 56166

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
I - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-012
Client Id: CPC-SB06-1.5-2.5
Data File: 6G03705.D
Analysis Date: 08/20/07 09:20
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 93

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0054	U	7421-93-4	Endrin Aldehyde	0.0054	U
319-84-6	alpha-BHC	0.0054	U	53494-70-5	Endrin Ketone	0.0054	UJ
319-85-7	beta-BHC	0.0054	U	58-89-9	gamma-BHC	0.0054	UJ
57-74-9	Chlordane	0.011	U	76-44-8	Heptachlor	0.0054	UJ
319-86-8	delta-BHC	0.0054	UJ	1024-57-3	Heptachlor Epoxide	0.0054	U
60-57-1	Dieldrin	0.0054	U	72-43-5	Methoxychlor	0.0054	UJ
959-98-8	Endosulfan I	0.0054	UJ	72-54-8	p,p'-DDD	0.0027	U
33213-65-9	Endosulfan II	0.0054	U	72-55-9	p,p'-DDE	0.0027	U
1031-07-8	Endosulfan Sulfate	0.0054	UJ	50-29-3	p,p'-DDT	0.0027	UJ
72-20-8	Endrin	0.0054	U	8001-35-2	Toxaphene	0.027	U

Worksheet #: 56166

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-013
Client Id: CPC-SB05-1-3
Data File: 6G03706.D
Analysis Date: 08/20/07 09:35
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 94

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0053	U	7421-93-4	Endrin Aldehyde	0.0053	U
319-84-6	alpha-BHC	0.0053	U	53494-70-5	Endrin Ketone	0.0053	UJ
319-85-7	beta-BHC	0.0053	U	58-89-9	gamma-BHC	0.0053	UJ
57-74-9	Chlordane	0.011	U	76-44-8	Heptachlor	0.0053	UJ
319-86-8	delta-BHC	0.0053	UJ	1024-57-3	Heptachlor Epoxide	0.0053	U
60-57-1	Dieldrin	0.0053	U	72-43-5	Methoxychlor	0.0053	UJ
959-98-8	Endosulfan I	0.0053	UJ	72-54-8	p,p'-DDD	0.0027	U
33213-65-9	Endosulfan II	0.0053	U	72-55-9	p,p'-DDE	0.0027	U
1031-07-8	Endosulfan Sulfate	0.0053	UJ	50-29-3	p,p'-DDT	0.0027	UJ
72-20-8	Endrin	0.0053	U	8001-35-2	Toxaphene	0.027	U

Worksheet #: 56166

Total Target Concentration 0

" - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-014
Client Id: CPC-SB05-3-5
Data File: 6G03707.D
Analysis Date: 08/20/07 09:50
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 96

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0052	U	7421-93-4	Endrin Aldehyde	0.0052	U
319-84-6	alpha-BHC	0.0052	U	53494-70-5	Endrin Ketone	0.0052	U J
319-85-7	beta-BHC	0.0052	U	58-89-9	gamma-BHC	0.0052	U J
57-74-9	Chlordane	0.010	U	76-44-8	Heptachlor	0.0052	U J
319-86-8	delta-BHC	0.0052	U J	1024-57-3	Heptachlor Epoxide	0.0052	U
60-57-1	Dieldrin	0.0052	U	72-43-5	Methoxychlor	0.0052	U J
959-98-8	Endosulfan I	0.0052	U J	72-54-8	p,p'-DDD	0.0026	U
33213-65-9	Endosulfan II	0.0052	U	72-55-9	p,p'-DDE	0.0026	U
1031-07-8	Endosulfan Sulfate	0.0052	U J	50-29-3	p,p'-DDT	0.0026	U J
72-20-8	Endrin	0.0052	U	8001-35-2	Toxaphene	0.026	U

Worksheet #: 56166

Total Target Concentration 0

" - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-015
Client Id: CPC-SB05-5-7
Data File: 6G03708.D
Analysis Date: 08/20/07 10:05
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 98

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0051	U	7421-93-4	Endrin Aldehyde	0.0051	U
319-84-6	alpha-BHC	0.0051	U	53494-70-5	Endrin Ketone	0.0051	UJ
319-85-7	beta-BHC	0.0051	U	58-89-9	gamma-BHC	0.0051	UJ
57-74-9	Chlordane	0.010	U	76-44-8	Heptachlor	0.0051	UJ
319-86-8	delta-BHC	0.0051	UJ	1024-57-3	Heptachlor Epoxide	0.0051	U
60-57-1	Dieldrin	0.0051	U	72-43-5	Methoxychlor	0.0051	UJ
959-98-8	Endosulfan I	0.0051	UJ	72-54-8	p,p'-DDD	0.0026	U
33213-65-9	Endosulfan II	0.0051	U	72-55-9	p,p'-DDE	0.0026	U
1031-07-8	Endosulfan Sulfate	0.0051	UJ	50-29-3	p,p'-DDT	0.0026	UJ
72-20-8	Endrin	0.0051	U	8001-35-2	Toxaphene	0.026	U

Worksheet #: 56166

Total Target Concentration 0

" - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-016
Client Id: CPC-SB05-7-9
Data File: 5G14284.D
Analysis Date: 08/20/07 07:00
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 98

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0051	U	7421-93-4	Endrin Aldehyde	0.0051	U
319-84-6	alpha-BHC	0.0051	U	53494-70-5	Endrin Ketone	0.0051	U J
319-85-7	beta-BHC	0.0051	U	58-89-9	gamma-BHC	0.0051	U J
57-74-9	Chlordane	0.010	U	76-44-8	Heptachlor	0.0051	U J
319-86-8	delta-BHC	0.0051	U J	1024-57-3	Heptachlor Epoxide	0.0051	U
60-57-1	Dieldrin	0.0051	U	72-43-5	Methoxychlor	0.0051	U J
959-98-8	Endosulfan I	0.0051	U J	72-54-8	p,p'-DDD	0.0026	U
33213-65-9	Endosulfan II	0.0051	U	72-55-9	p,p'-DDE	0.0026	U
1031-07-8	Endosulfan Sulfate	0.0051	U J	50-29-3	p,p'-DDT	0.0026	U J
72-20-8	Endrin	0.0051	U	8001-35-2	Toxaphene	0.026	U

Worksheet #: 56166

Total Target Concentration 0

"U" - Indicates the compound was analyzed but not detected.
' - Indicates the analyte was found in the blank as well as in the sample.
' - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-017
Client Id: CPC-SB02-5-7
Data File: 6G03709.D
Analysis Date: 08/20/07 10:20
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 96

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0052	U	7421-93-4	Endrin Aldehyde	0.0052	U
319-84-6	alpha-BHC	0.0052	U	53494-70-5	Endrin Ketone	0.0052	U J
319-85-7	beta-BHC	0.0052	U	58-89-9	gamma-BHC	0.0052	U J
57-74-9	**Chlordane**	**0.010**	**0.021**	76-44-8	Heptachlor	0.0052	U J
319-86-8	delta-BHC	0.0052	U J	1024-57-3	Heptachlor Epoxide	0.0052	U
60-57-1	Dieldrin	0.0052	U	72-43-5	Methoxychlor	0.0052	U J
959-98-8	Endosulfan I	0.0052	U J	72-54-8	p,p'-DDD	0.0026	U
33213-65-9	Endosulfan II	0.0052	U	72-55-9	p,p'-DDE	0.0026	U
1031-07-8	Endosulfan Sulfate	0.0052	U J	50-29-3	p,p'-DDT	0.0026	U J
72-20-8	Endrin	0.0052	U	8001-35-2	Toxaphene	0.026	U

Worksheet #: 56166

Total Target Concentration 0.021

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-018
Client Id: CPC-SB02-7-9
Data File: 5G14285.D
Analysis Date: 08/20/07 07:18
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 97

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0052	U	7421-93-4	Endrin Aldehyde	0.0052	U
319-84-6	alpha-BHC	0.0052	U	53494-70-5	Endrin Ketone	0.0052	UJ
319-85-7	beta-BHC	0.0052	U	58-89-9	gamma-BHC	0.0052	UJ
57-74-9	Chlordane	0.010	U	76-44-8	Heptachlor	0.0052	UJ
319-86-8	delta-BHC	0.0052	UJ	1024-57-3	Heptachlor Epoxide	0.0052	U
60-57-1	Dieldrin	0.0052	U	72-43-5	Methoxychlor	0.0052	UJ
959-98-8	Endosulfan I	0.0052	UJ	72-54-8	p,p'-DDD	0.0026	U
33213-65-9	Endosulfan II	0.0052	U	72-55-9	p,p'-DDE	0.0026	U
1031-07-8	Endosulfan Sulfate	0.0052	UJ	50-29-3	p,p'-DDT	0.0026	UJ
72-20-8	Endrin	0.0052	U	8001-35-2	Toxaphene	0.026	U

Worksheet #: 56166

Total Target Concentration 0

'' - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-019
Client Id: CPC-DW04-13-14
Data File: 5G14286.D
Analysis Date: 08/20/07 07:37
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 82

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.0061	U	7421-93-4	Endrin Aldehyde	0.0061	U
319-84-6	alpha-BHC	0.0061	U	53494-70-5	Endrin Ketone	0.0061	U J
319-85-7	beta-BHC	0.0061	U	58-89-9	gamma-BHC	0.0061	U J
57-74-9	Chlordane	0.012	U	76-44-8	Heptachlor	0.0061	U J
319-86-8	delta-BHC	0.0061	U J	1024-57-3	Heptachlor Epoxide	0.0061	U
60-57-1	Dieldrin	0.0061	U	72-43-5	Methoxychlor	0.0061	U J
959-98-8	Endosulfan I	0.0061	U J	72-54-8	p,p'-DDD	0.0030	U
33213-65-9	Endosulfan II	0.0061	U	72-55-9	p,p'-DDE	0.0030	U
1031-07-8	Endosulfan Sulfate	0.0061	U J	50-29-3	p,p'-DDT	0.0030	U J
72-20-8	Endrin	0.0061	U	8001-35-2	Toxaphene	0.030	U

Worksheet #: 56166

Total Target Concentration 0

" - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PESTICIDE REPORT

Sample Number: AC32378-020
Client Id: CPC-Rinsate
Data File: 6G03731.D
Analysis Date: 08/21/07 09:54
Date Rec/Extracted: 08/16/07-08/20/07

Matrix: Aqueous
Initial Vol: 890ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
309-00-2	Aldrin	0.056	U	7421-93-4	Endrin Aldehyde	0.056	U
319-84-6	alpha-BHC	0.056	U J	53494-70-5	Endrin Ketone	0.056	U
319-85-7	beta-BHC	0.056	U	58-89-9	gamma-BHC	0.056	U
57-74-9	Chlordane	0.11	U	76-44-8	Heptachlor	0.056	U J
319-86-8	delta-BHC	0.056	U J	1024-57-3	Heptachlor Epoxide	0.056	U
60-57-1	Dieldrin	0.056	U	72-43-5	Methoxychlor	0.056	U
959-98-8	Endosulfan I	0.056	U	72-54-8	p,p'-DDD	0.028	U J
33213-65-9	Endosulfan II	0.056	U	72-55-9	p,p'-DDE	0.028	U
1031-07-8	Endosulfan Sulfate	0.056	U	50-29-3	p,p'-DDT	0.028	U
72-20-8	Endrin	0.056	U	8001-35-2	Toxaphene	0.28	U

Worksheet #: 56166

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-001
Client Id: CPC-MW01
Data File: 2G32826.D
Analysis Date: 08/17/07 06:38
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-002
Client Id: CPC-MW08
Data File: 2G32827.D
Analysis Date: 08/17/07 06:52
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-003
Client Id: CPC-MW03
Data File: 2G32828.D
Analysis Date: 08/17/07 07:06
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-004
Client Id: CPC-MW04
Data File: 2G32829.D
Analysis Date: 08/17/07 07:20
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-005
Client Id: CPC-MW05
Data File: 2G32830.D
Analysis Date: 08/17/07 07:35
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-006
Client Id: CPC-MW06
Data File: 2G32831.D
Analysis Date: 08/17/07 07:49
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-007
Client Id: CPC-MW07
Data File: 2G32832.D
Analysis Date: 08/17/07 08:03
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 980ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.26	U	12672-29-6	Aroclor-1248	0.26	U
11104-28-2	Aroclor-1221	0.26	U	11097-69-1	Aroclor-1254	0.26	U
11141-16-5	Aroclor-1232	0.26	U	11096-82-5	Aroclor-1260	0.26	U
53469-21-9	Aroclor-1242	0.26	U	37324-23-5	Aroclor-1262	0.26	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-008
Client Id: CPC-MW09
Data File: 2G32833.D
Analysis Date: 08/17/07 08:18
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-009
Client Id: CPC-MW10
Data File: 2G32824.D
Analysis Date: 08/17/07 06:09
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 1000ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32351-012
Client Id: CPC-DUPLICATE
Data File: 2G32834.D
Analysis Date: 08/17/07 08:32
Date Rec/Extracted: 08/15/07-08/16/07

Matrix: Aqueous
Initial Vol: 990ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.25	U	12672-29-6	Aroclor-1248	0.25	U
11104-28-2	Aroclor-1221	0.25	U	11097-69-1	Aroclor-1254	0.25	U
11141-16-5	Aroclor-1232	0.25	U	11096-82-5	Aroclor-1260	0.25	U
53469-21-9	Aroclor-1242	0.25	U	37324-23-5	Aroclor-1262	0.25	U

Worksheet #: 55977

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration us

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-001
Client Id: CPC-DW01-10-11
Data File: 3G29841.D
Analysis Date: 08/20/07 07:47
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 86

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.029	U	12672-29-6	Aroclor-1248	0.029	U
11104-28-2	Aroclor-1221	0.029	U	11097-69-1	Aroclor-1254	0.029	U
11141-16-5	Aroclor-1232	0.029	U	11096-82-5	Aroclor-1260	0.029	U
53469-21-9	Aroclor-1242	0.029	U	37324-23-5	Aroclor-1262	0.029	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-002
Client Id: CPC-DW02-10-11
Data File: 3G29842.D
Analysis Date: 08/20/07 08:03
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 83

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.030	U	12672-29-6	Aroclor-1248	0.030	U
11104-28-2	Aroclor-1221	0.030	U	11097-69-1	Aroclor-1254	0.030	U
11141-16-5	Aroclor-1232	0.030	U	11096-82-5	Aroclor-1260	0.030	U
53469-21-9	Aroclor-1242	0.030	U	37324-23-5	Aroclor-1262	0.030	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-003
Client Id: CPC-DW03-10-11
Data File: 3G29843.D
Analysis Date: 08/20/07 08:19
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 90

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.028	U	12672-29-6	Aroclor-1248	0.028	U
11104-28-2	Aroclor-1221	0.028	U	11097-69-1	Aroclor-1254	0.028	U
11141-16-5	Aroclor-1232	0.028	U	11096-82-5	Aroclor-1260	0.028	U
53469-21-9	Aroclor-1242	0.028	U	37324-23-5	Aroclor-1262	0.028	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-004
Client Id: CPC-DW04-8-9
Data File: 3G29844.D
Analysis Date: 08/20/07 08:35
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 85

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.029	U	12672-29-6	Aroclor-1248	0.029	U
11104-28-2	Aroclor-1221	0.029	U	11097-69-1	Aroclor-1254	0.029	U
11141-16-5	Aroclor-1232	0.029	U	11096-82-5	Aroclor-1260	0.029	U
53469-21-9	Aroclor-1242	0.029	U	37324-23-5	Aroclor-1262	0.029	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-005
Client Id: CPC-SOIL-Duplicate
Data File: 3G29845.D
Analysis Date: 08/20/07 08:51
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 92

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.027	U	12672-29-6	Aroclor-1248	0.027	U
1104-28-2	Aroclor-1221	0.027	U	11097-69-1	Aroclor-1254	0.027	U
1141-16-5	Aroclor-1232	0.027	U	11096-82-5	Aroclor-1260	0.027	U
53469-21-9	Aroclor-1242	0.027	U	37324-23-5	Aroclor-1262	0.027	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-006
Client Id: CPC-SB01-1.5-2.5
Data File: 3G29846.D
Analysis Date: 08/20/07 09:07
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 95

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.026	U	12672-29-6	Aroclor-1248	0.026	U
1104-28-2	Aroclor-1221	0.026	U	11097-69-1	Aroclor-1254	0.026	U
1141-16-5	Aroclor-1232	0.026	U	11096-82-5	Aroclor-1260	0.026	U
53469-21-9	Aroclor-1242	0.026	U	37324-23-5	Aroclor-1262	0.026	U

Worksheet #: 55995

Total Target Concentration 0

" - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-007
Client Id: CPC-SB02-1-5
Data File: 3G29847.D
Analysis Date: 08/20/07 09:23
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 94

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.027	U	12672-29-6	Aroclor-1248	0.027	U
1104-28-2	Aroclor-1221	0.027	U	11097-69-1	Aroclor-1254	0.027	U
1141-16-5	Aroclor-1232	0.027	U	11096-82-5	Aroclor-1260	0.027	U
53469-21-9	Aroclor-1242	0.027	U	37324-23-5	Aroclor-1262	0.027	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-008
Client Id: CPC-SB03-1.5-3.5
Data File: 3G29835.D
Analysis Date: 08/20/07 06:10
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 93

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.027	U	12672-29-6	Aroclor-1248	0.027	U
1104-28-2	Aroclor-1221	0.027	U	11097-69-1	Aroclor-1254	0.027	U
1141-16-5	Aroclor-1232	0.027	U	11096-82-5	Aroclor-1260	0.027	U
53469-21-9	Aroclor-1242	0.027	U	37324-23-5	Aroclor-1262	0.027	U

Worksheet #: 55995

Total Target Concentration 0

ᵀᵀ - Indicates the compound was analyzed but not detected.
- Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-012
Client Id: CPC-SB06-1.5-2.5
Data File: 3G29849.D
Analysis Date: 08/20/07 09:55
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 93

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.027	U	12672-29-6	Aroclor-1248	0.027	U
11104-28-2	Aroclor-1221	0.027	U	11097-69-1	Aroclor-1254	0.027	U
1141-16-5	Aroclor-1232	0.027	U	11096-82-5	Aroclor-1260	0.027	U
J3469-21-9	Aroclor-1242	0.027	U	37324-23-5	Aroclor-1262	0.027	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-013
Client Id: CPC-SB05-1-3
Data File: 3G29850.D
Analysis Date: 08/20/07 10:11
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 94

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.027	U	12672-29-6	Aroclor-1248	0.027	U
11104-28-2	Aroclor-1221	0.027	U	11097-69-1	Aroclor-1254	0.027	U
1141-16-5	Aroclor-1232	0.027	U	11096-82-5	Aroclor-1260	0.027	U
53469-21-9	Aroclor-1242	0.027	U	37324-23-5	Aroclor-1262	0.027	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
R - Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of e instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-014
Client Id: CPC-SB05-3-5
Data File: 3G29851.D
Analysis Date: 08/20/07 10:28
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 96

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.026	U	12672-29-6	Aroclor-1248	0.026	U
11104-28-2	Aroclor-1221	0.026	U	11097-69-1	Aroclor-1254	0.026	U
11141-16-5	Aroclor-1232	0.026	U	11096-82-5	Aroclor-1260	0.026	U
53469-21-9	Aroclor-1242	0.026	U	37324-23-5	Aroclor-1262	0.026	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.
R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-017
Client Id: CPC-SB02-5-7
Data File: 3G29855.D
Analysis Date: 08/20/07 11:35
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 96

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.026	U	12672-29-6	Aroclor-1248	0.026	U
11104-28-2	Aroclor-1221	0.026	U	11097-69-1	Aroclor-1254	0.026	U
11141-16-5	Aroclor-1232	0.026	U	11096-82-5	Aroclor-1260	0.026	U
53469-21-9	Aroclor-1242	0.026	U	37324-23-5	Aroclor-1262	0.026	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
- Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-018
Client Id: CPC-SB02-7-9
Data File: 3G29856.D
Analysis Date: 08/20/07 11:51
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 97

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.026	U	12672-29-6	Aroclor-1248	0.026	U
1104-28-2	Aroclor-1221	0.026	U	11097-69-1	Aroclor-1254	0.026	U
1141-16-5	Aroclor-1232	0.026	U	11096-82-5	Aroclor-1260	0.026	U
53469-21-9	Aroclor-1242	0.026	U	37324-23-5	Aroclor-1262	0.026	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-019
Client Id: CPC-DW04-13-14
Data File: 3G29857.D
Analysis Date: 08/20/07 12:08
Date Rec/Extracted: 08/16/07-08/19/07

Matrix: Soil
Initial Vol: 20g
Final Vol: 10ml
Dilution: 1
Solids: 82

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.030	U	12672-29-6	Aroclor-1248	0.030	U
11104-28-2	Aroclor-1221	0.030	U	11097-69-1	Aroclor-1254	0.030	U
11141-16-5	Aroclor-1232	0.030	U	11096-82-5	Aroclor-1260	0.030	U
53469-21-9	Aroclor-1242	0.030	U	37324-23-5	Aroclor-1262	0.030	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1

ORGANICS PCB REPORT

Sample Number: AC32378-020
Client Id: CPC-Rinsate
Data File: 3G29890.D
Analysis Date: 08/21/07 06:17
Date Rec/Extracted: 08/16/07-08/20/07

Matrix: Aqueous
Initial Vol: 890ml
Final Vol: 5ml
Dilution: 1
Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
12674-11-2	Aroclor-1016	0.28	U	12672-29-6	Aroclor-1248	0.28	U
11104-28-2	Aroclor-1221	0.28	U	11097-69-1	Aroclor-1254	0.28	U
11141-16-5	Aroclor-1232	0.28	U	11096-82-5	Aroclor-1260	0.28	U
53469-21-9	Aroclor-1242	0.28	U	37324-23-5	Aroclor-1262	0.28	U

Worksheet #: 55995

Total Target Concentration 0

U - Indicates the compound was analyzed but not detected.
B - Indicates the analyte was found in the blank as well as in the sample.
E - Indicates the analyte concentration exceeds the calibration range of the instrument.

R - Retention Time Out
J - Indicates an estimated value when a compound is detected at less than the specified detection limit.
d - Pesticide %Diff>50% between columns due to coelution. Lower concentration used.

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-001	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW01	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-39-3	Barium	50	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-70-2	Calcium	2000	24000	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7439-89-6	Iron	280	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7439-92-1	Lead	4.0	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7439-95-4	Magnesium	2000	2300	1	08/23/07	8241	SW8241A	20	P	PEICP1
7439-96-5	Manganese	40	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	17	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-09-7	Potassium	5000	26000	1	08/23/07	8241	SW8241B	20	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-23-5	Sodium	5000	22000	1	08/23/07	8241	SW8241B	20	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	20	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-002	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW08	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	250	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-39-3	Barium	50	210	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-70-2	Calcium	2000	25000	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7439-89-6	Iron	280	530	1	08/23/07	8241	SW8241A	22	P	PEICP1
7439-92-1	Lead	4.0	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7439-95-4	Magnesium	2000	2500	1	08/23/07	8241	SW8241A	22	P	PEICP1
7439-96-5	Manganese	40	81	1	08/23/07	8241	SW8241A	22	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	18	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-09-7	Potassium	5000	13000	1	08/23/07	8241	SW8241B	22	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-23-5	Sodium	5000	20000	1	08/23/07	8241	SW8241B	22	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	22	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-003	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW03	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-39-3	Barium	50	97	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-70-2	Calcium	2000	16000	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7439-89-6	Iron	280	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7439-92-1	Lead	4.0	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7439-95-4	Magnesium	2000	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7439-96-5	Manganese	40	74	1	08/23/07	8241	SW8241A	23	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	19	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-09-7	Potassium	5000	ND	1	08/23/07	8241	SW8241B	23	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-23-5	Sodium	5000	ND	1	08/23/07	8241	SW8241B	23	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	23	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-004	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW04	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-39-3	Barium	50	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-70-2	Calcium	2000	56000	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7439-89-6	Iron	280	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7439-92-1	Lead	4.0	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7439-95-4	Magnesium	2000	4600	1	08/23/07	8241	SW8241A	24	P	PEICP1
7439-96-5	Manganese	40	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	22	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-09-7	Potassium	5000	11000	1	08/23/07	8241	SW8241B	24	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-23-5	Sodium	5000	25000	1	08/23/07	8241	SW8241B	24	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	24	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-005	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW05	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-39-3	Barium	50	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-70-2	Calcium	2000	58000	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7439-89-6	Iron	280	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7439-92-1	Lead	4.0	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7439-95-4	Magnesium	2000	3100	1	08/23/07	8241	SW8241A	25	P	PEICP1
7439-96-5	Manganese	40	83	1	08/23/07	8241	SW8241A	25	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	23	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-09-7	Potassium	5000	26000	1	08/23/07	8241	SW8241B	25	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-23-5	Sodium	5000	22000	1	08/23/07	8241	SW8241B	25	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	25	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-006	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW06	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-39-3	Barium	50	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-70-2	Calcium	2000	25000	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7439-89-6	Iron	280	1100	1	08/23/07	8241	SW8241A	30	P	PEICP1
7439-92-1	Lead	4.0	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7439-95-4	Magnesium	2000	2200	1	08/23/07	8241	SW8241A	30	P	PEICP1
7439-96-5	Manganese	40	110	1	08/23/07	8241	SW8241A	30	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	24	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-09-7	Potassium	5000	8700	1	08/23/07	8241	SW8241B	30	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-23-5	Sodium	5000	10000	1	08/23/07	8241	SW8241B	30	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	30	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-007	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW07	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-39-3	Barium	50	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-70-2	Calcium	2000	29000	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7439-89-6	Iron	280	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7439-92-1	Lead	4.0	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7439-95-4	Magnesium	2000	2900	1	08/23/07	8241	SW8241A	31	P	PEICP1
7439-96-5	Manganese	40	130	1	08/23/07	8241	SW8241A	31	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	25	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-09-7	Potassium	5000	13000	1	08/23/07	8241	SW8241B	31	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-23-5	Sodium	5000	15000	1	08/23/07	8241	SW8241B	31	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	31	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-008	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW09	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-39-3	Barium	50	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-70-2	Calcium	2000	30000	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7439-89-6	Iron	280	340	1	08/23/07	8241	SW8241A	32	P	PEICP1
7439-92-1	Lead	4.0	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7439-95-4	Magnesium	2000	2400	1	08/23/07	8241	SW8241A	32	P	PEICP1
7439-96-5	Manganese	40	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	26	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-09-7	Potassium	5000	16000	1	08/23/07	8241	SW8241B	32	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-23-5	Sodium	5000	27000	1	08/23/07	8241	SW8241B	32	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	32	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-009	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-MW10	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	300	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-39-3	Barium	50	58	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-70-2	Calcium	2000	25000	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7439-89-6	Iron	280	710	1	08/23/07	8241	SW8241A	13	P	PEICP1
7439-92-1	Lead	4.0	10	1	08/23/07	8241	SW8241A	13	P	PEICP1
7439-95-4	Magnesium	2000	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7439-96-5	Manganese	40	330 J	1	08/23/07	8241	SW8241A	13	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	13	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-09-7	Potassium	5000	17000	1	08/23/07	8241	SW8241B	13	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-23-5	Sodium	5000	15000	1	08/23/07	8241	SW8241B	13	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	13	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32351-012	% Solid:	0	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-DUPLICATE	Units:	UG/L	Lab Code:		Sdg No:	
Matrix:	AQUEOUS	Date Rec:	8/16/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	180	320	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-36-0	Antimony	12	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-38-2	Arsenic	7.5	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-39-3	Barium	50	210	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-41-7	Beryllium	4.0	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-43-9	Cadmium	3.5	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-70-2	Calcium	2000	26000	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-47-3	Chromium	50	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-48-4	Cobalt	20	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-50-8	Copper	50	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7439-89-6	Iron	280	660	1	08/23/07	8241	SW8241A	33	P	PEICP1
7439-92-1	Lead	4.0	5.6	1	08/23/07	8241	SW8241A	33	P	PEICP1
7439-95-4	Magnesium	2000	2600	1	08/23/07	8241	SW8241A	33	P	PEICP1
7439-96-5	Manganese	40	83	1	08/23/07	8241	SW8241A	33	P	PEICP1
7439-97-6	Mercury	0.70	ND	1	08/23/07	8241	H8241SW	27	CV	HGCV1
7440-02-0	Nickel	50	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-09-7	Potassium	5000	13000	1	08/23/07	8241	SW8241B	33	P	PEICPRAD1
7782-49-2	Selenium	40	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-22-4	Silver	20	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-23-5	Sodium	5000	20000	1	08/23/07	8241	SW8241B	33	P	PEICPRAD1
7440-28-0	Thallium	10	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-62-2	Vanadium	50	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1
7440-66-6	Zinc	50	ND	1	08/23/07	8241	SW8241A	33	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-001	% Solid:	86	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-DW01-10-11	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	230	530	100	08/22/07	8234	S8234B	13	P	PEICPRAD1
7440-36-0	Antimony	2.3	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-38-2	Arsenic	2.3	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-39-3	Barium	12	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-41-7	Beryllium	0.70	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-43-9	Cadmium	0.70	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-70-2	Calcium	1200	ND	100	08/22/07	8234	S8234B	13	P	PEICPRAD1
7440-47-3	Chromium	5.8	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-48-4	Cobalt	2.9	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-50-8	Copper	5.8	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7439-89-6	Iron	230	1400	100	08/22/07	8234	S8234B	13	P	PEICPRAD1
7439-92-1	Lead	5.8	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7439-95-4	Magnesium	580	ND	100	08/22/07	8234	S8234B	13	P	PEICPRAD1
7439-96-5	Manganese	12	30	100	08/21/07	8234	S8234A	13	P	PEICP1
7439-97-6	Mercury	0.097	ND	167	08/21/07	8234	h8234s	13	CV	HGCV1
7440-02-0	Nickel	5.8	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-09-7	Potassium	580	ND	100	08/22/07	8234	S8234B	13	P	PEICPRAD1
7782-49-2	Selenium	2.1	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-22-4	Silver	1.7	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-23-5	Sodium	290	ND	100	08/22/07	8234	S8234B	13	P	PEICPRAD1
7440-28-0	Thallium	1.4	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-62-2	Vanadium	12	ND	100	08/21/07	8234	S8234A	13	P	PEICP1
7440-66-6	Zinc	12	ND	100	08/21/07	8234	S8234A	13	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-002	% Solid:	83	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-DW02-10-11	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	240	1600	100	08/22/07	8234	S8234B	20	P	PEICPRAD1
7440-36-0	Antimony	2.4	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-38-2	Arsenic	2.4	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-39-3	Barium	12	24	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-41-7	Beryllium	0.72	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-43-9	Cadmium	0.72	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-70-2	Calcium	1200	2400	100	08/22/07	8234	S8234B	20	P	PEICPRAD1
7440-47-3	Chromium	6.0	6.7 J	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-48-4	Cobalt	3.0	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-50-8	Copper	6.0	13	100	08/21/07	8234	S8234A	20	P	PEICP1
7439-89-6	Iron	240	19000	100	08/22/07	8234	S8234B	20	P	PEICPRAD1
7439-92-1	Lead	6.0	16	100	08/21/07	8234	S8234A	20	P	PEICP1
7439-95-4	Magnesium	600	930	100	08/22/07	8234	S8234B	20	P	PEICPRAD1
7439-96-5	Manganese	12	85	100	08/21/07	8234	S8234A	20	P	PEICP1
7439-97-6	Mercury	0.10	ND	167	08/21/07	8234	h8234s	17	CV	HGCV1
7440-02-0	Nickel	6.0	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-09-7	Potassium	600	ND	100	08/22/07	8234	S8234B	20	P	PEICPRAD1
7782-49-2	Selenium	2.2	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-22-4	Silver	1.8	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-23-5	Sodium	300	ND	100	08/22/07	8234	S8234B	20	P	PEICPRAD1
7440-28-0	Thallium	1.4	ND	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-62-2	Vanadium	12	17	100	08/21/07	8234	S8234A	20	P	PEICP1
7440-66-6	Zinc	12	68	100	08/21/07	8234	S8234A	20	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-003	% Solid:	90	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-DW03-10-11	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	220	530	100	08/22/07	8234	S8234B	22	P	PEICPRAD1
7440-36-0	Antimony	2.2	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-38-2	Arsenic	2.2	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-39-3	Barium	11	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-41-7	Beryllium	0.67	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-43-9	Cadmium	0.67	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-70-2	Calcium	1100	ND	100	08/22/07	8234	S8234B	22	P	PEICPRAD1
7440-47-3	Chromium	5.6	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-48-4	Cobalt	2.8	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-50-8	Copper	5.6	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7439-89-6	Iron	220	1700	100	08/22/07	8234	S8234B	22	P	PEICPRAD1
7439-92-1	Lead	5.6	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7439-95-4	Magnesium	560	ND	100	08/22/07	8234	S8234B	22	P	PEICPRAD1
7439-96-5	Manganese	11	20	100	08/21/07	8234	S8234A	22	P	PEICP1
7439-97-6	Mercury	0.093	ND	167	08/21/07	8234	h8234s	18	CV	HGCV1
7440-02-0	Nickel	5.6	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-09-7	Potassium	560	ND	100	08/22/07	8234	S8234B	22	P	PEICPRAD1
7782-49-2	Selenium	2.0	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-22-4	Silver	1.7	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-23-5	Sodium	280	ND	100	08/22/07	8234	S8234B	22	P	PEICPRAD1
7440-28-0	Thallium	1.3	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-62-2	Vanadium	11	ND	100	08/21/07	8234	S8234A	22	P	PEICP1
7440-66-6	Zinc	11	ND	100	08/21/07	8234	S8234A	22	P	PEICP1

Comments: _____

Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-004	% Solid:	85	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-DW04-8-9	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	240	1800	100	08/22/07	8234	S8234B	23	P	PEICPRAD1
7440-36-0	Antimony	2.4	ND	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-38-2	Arsenic	2.4	2.7	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-39-3	Barium	12	51	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-41-7	Beryllium	0.71	ND	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-43-9	Cadmium	0.71	1.7	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-70-2	Calcium	1200	1700	100	08/22/07	8234	S8234B	23	P	PEICPRAD1
7440-47-3	Chromium	5.9	28	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-48-4	Cobalt	2.9	ND	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-50-8	Copper	5.9	42	100	08/21/07	8234	S8234A	23	P	PEICP1
7439-89-6	Iron	240	5600	100	08/22/07	8234	S8234B	23	P	PEICPRAD1
7439-92-1	Lead	5.9	390	100	08/21/07	8234	S8234A	23	P	PEICP1
7439-95-4	Magnesium	590	1200	100	08/22/07	8234	S8234B	23	P	PEICPRAD1
7439-96-5	Manganese	12	28	100	08/21/07	8234	S8234A	23	P	PEICP1
7439-97-6	Mercury	0.098	ND	167	08/21/07	8234	h8234s	19	CV	HGCV1
7440-02-0	Nickel	5.9	6.2	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-09-7	Potassium	590	ND	100	08/22/07	8234	S8234B	23	P	PEICPRAD1
7782-49-2	Selenium	2.1	ND	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-22-4	Silver	1.8	17	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-23-5	Sodium	290	ND	100	08/22/07	8234	S8234B	23	P	PEICPRAD1
7440-28-0	Thallium	1.4	ND	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-62-2	Vanadium	12	ND	100	08/21/07	8234	S8234A	23	P	PEICP1
7440-66-6	Zinc	12	140	100	08/21/07	8234	S8234A	23	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-005	% Solid:	92	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SOIL-Duplicate	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	220	1100	100	08/22/07	8234	S8234B	24	P	PEICPRAD1
7440-36-0	Antimony	2.2	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-38-2	Arsenic	4.3	ND	200	08/22/07	8234	S8190B	17	P	PEICP1
7440-39-3	Barium	11	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-41-7	Beryllium	0.65	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-43-9	Cadmium	0.65	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-70-2	Calcium	2200	130000	200	08/22/07	8234	S8234C	10	P	PEICPRAD1
7440-47-3	Chromium	5.4	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-48-4	Cobalt	2.7	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-50-8	Copper	11	ND	200	08/22/07	8234	S8190B	17	P	PEICP1
7439-89-6	Iron	220	3800	100	08/22/07	8234	S8234B	24	P	PEICPRAD1
7439-92-1	Lead	11	ND	200	08/22/07	8234	S8190B	17	P	PEICP1
7439-95-4	Magnesium	540	78000	100	08/22/07	8234	S8234B	24	P	PEICPRAD1
7439-96-5	Manganese	11	120	100	08/21/07	8234	S8234A	24	P	PEICP1
7439-97-6	Mercury	0.091	ND	167	08/21/07	8234	h8234s	22	CV	HGCV1
7440-02-0	Nickel	5.4	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-09-7	Potassium	540	ND	100	08/22/07	8234	S8234B	24	P	PEICPRAD1
7782-49-2	Selenium	2.0	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-22-4	Silver	3.3	ND	200	08/22/07	8234	S8190B	17	P	PEICP1
7440-23-5	Sodium	270	ND	100	08/22/07	8234	S8234B	24	P	PEICPRAD1
7440-28-0	Thallium	1.3	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-62-2	Vanadium	11	ND	100	08/21/07	8234	S8234A	24	P	PEICP1
7440-66-6	Zinc	11	ND	100	08/21/07	8234	S8234A	24	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-006	% Solid:	95	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB01-1.5-2.5	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	210	3900	100	08/22/07	8234	S8234B	25	P	PEICPRAD1
7440-36-0	Antimony	2.1	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-38-2	Arsenic	2.1	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-39-3	Barium	11	13	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-41-7	Beryllium	0.63	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-43-9	Cadmium	0.63	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-70-2	Calcium	1100	ND	100	08/22/07	8234	S8234B	25	P	PEICPRAD1
7440-47-3	Chromium	5.3	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-48-4	Cobalt	2.6	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-50-8	Copper	5.3	6.0	100	08/21/07	8234	S8234A	25	P	PEICP1
7439-89-6	Iron	210	5800	100	08/22/07	8234	S8234B	25	P	PEICPRAD1
7439-92-1	Lead	5.3	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7439-95-4	Magnesium	530	770	100	08/22/07	8234	S8234B	25	P	PEICPRAD1
7439-96-5	Manganese	11	120	100	08/21/07	8234	S8234A	25	P	PEICP1
7439-97-6	Mercury	0.088	ND	167	08/21/07	8234	h8234s	23	CV	HGCV1
7440-02-0	Nickel	5.3	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-09-7	Potassium	530	ND	100	08/22/07	8234	S8234B	25	P	PEICPRAD1
7782-49-2	Selenium	1.9	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-22-4	Silver	1.6	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-23-5	Sodium	260	ND	100	08/22/07	8234	S8234B	25	P	PEICPRAD1
7440-28-0	Thallium	1.3	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-62-2	Vanadium	11	ND	100	08/21/07	8234	S8234A	25	P	PEICP1
7440-66-6	Zinc	11	ND	100	08/21/07	8234	S8234A	25	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-007	% Solid:	94	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB02-1-5	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	210	2100	100	08/22/07	8234	S8234B	26	P	PEICPRAD1
7440-36-0	Antimony	2.1	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-38-2	Arsenic	2.1	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-39-3	Barium	11	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-41-7	Beryllium	0.64	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-43-9	Cadmium	0.64	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-70-2	Calcium	1100	ND	100	08/22/07	8234	S8234B	26	P	PEICPRAD1
7440-47-3	Chromium	5.3	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-48-4	Cobalt	2.7	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-50-8	Copper	5.3	8.5	100	08/21/07	8234	S8234A	26	P	PEICP1
7439-89-6	Iron	210	4300	100	08/22/07	8234	S8234B	26	P	PEICPRAD1
7439-92-1	Lead	5.3	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7439-95-4	Magnesium	530	ND	100	08/22/07	8234	S8234B	26	P	PEICPRAD1
7439-96-5	Manganese	11	64	100	08/21/07	8234	S8234A	26	P	PEICP1
7439-97-6	Mercury	0.089	ND	167	08/21/07	8234	h8234s	24	CV	HGCV1
7440-02-0	Nickel	5.3	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-09-7	Potassium	530	ND	100	08/22/07	8234	S8234B	26	P	PEICPRAD1
7782-49-2	Selenium	1.9	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-22-4	Silver	1.6	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-23-5	Sodium	270	ND	100	08/22/07	8234	S8234B	26	P	PEICPRAD1
7440-28-0	Thallium	1.3	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-62-2	Vanadium	11	ND	100	08/21/07	8234	S8234A	26	P	PEICP1
7440-66-6	Zinc	11	ND	100	08/21/07	8234	S8234A	26	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-008	% Solid:	93	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB03-1.5-3.5	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	220	2700	100	08/30/07	8269	S8269B	13	P	PEICPRAD1
7440-36-0	Antimony	2.2	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-38-2	Arsenic	2.2	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-39-3	Barium	11	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-41-7	Beryllium	0.65	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-43-9	Cadmium	0.65	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-70-2	Calcium	1100	ND	100	08/30/07	8269	S8269B	13	P	PEICPRAD1
7440-47-3	Chromium	5.4	180 J	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-48-4	Cobalt	2.7	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-50-8	Copper	5.4	31	100	08/29/07	8269	S8269A	13	P	PEICP1
7439-89-6	Iron	220	4400	100	08/30/07	8269	S8269B	13	P	PEICPRAD1
7439-92-1	Lead	5.4	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7439-95-4	Magnesium	540	ND	100	08/30/07	8269	S8269B	13	P	PEICPRAD1
7439-96-5	Manganese	11	70 J	100	08/29/07	8269	S8269A	13	P	PEICP1
7439-97-6	Mercury	0.090	ND	167	08/30/07	8269	H8269S	13	CV	HGCV2
7440-02-0	Nickel	5.4	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-09-7	Potassium	540	ND	100	08/30/07	8269	S8269B	13	P	PEICPRAD1
7782-49-2	Selenium	1.9	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-22-4	Silver	1.6	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-23-5	Sodium	270	ND	100	08/30/07	8269	S8269B	13	P	PEICPRAD1
7440-28-0	Thallium	1.3	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-62-2	Vanadium	11	ND	100	08/29/07	8269	S8269A	13	P	PEICP1
7440-66-6	Zinc	11	ND	100	08/29/07	8269	S8269A	13	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-011	% Solid:	90	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB04-1.5-2.5	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	220	9000	100	08/22/07	8234	S8234B	34	P	PEICPRAD1
7440-36-0	Antimony	2.2	ND	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-38-2	Arsenic	2.2	3.2	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-39-3	Barium	11	18	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-41-7	Beryllium	0.67	ND	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-43-9	Cadmium	0.67	ND	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-70-2	Calcium	1100	ND	100	08/22/07	8234	S8234B	34	P	PEICPRAD1
7440-47-3	Chromium	5.6	11	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-48-4	Cobalt	2.8	4.4	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-50-8	Copper	5.6	8.7	100	08/21/07	8234	S8234A	34	P	PEICP1
7439-89-6	Iron	220	16000	100	08/22/07	8234	S8234B	34	P	PEICPRAD1
7439-92-1	Lead	5.6	5.7	100	08/21/07	8234	S8234A	34	P	PEICP1
7439-95-4	Magnesium	560	1400	100	08/22/07	8234	S8234B	34	P	PEICPRAD1
7439-96-5	Manganese	11	98	100	08/21/07	8234	S8234A	34	P	PEICP1
7439-97-6	Mercury	0.093	ND	167	08/21/07	8234	h8234s	28	CV	HGCV1
7440-02-0	Nickel	5.6	ND	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-09-7	Potassium	560	ND	100	08/22/07	8234	S8234B	34	P	PEICPRAD1
7782-49-2	Selenium	2.0	ND	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-22-4	Silver	1.7	ND	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-23-5	Sodium	280	ND	100	08/22/07	8234	S8234B	34	P	PEICPRAD1
7440-28-0	Thallium	1.3	ND	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-62-2	Vanadium	11	16	100	08/21/07	8234	S8234A	34	P	PEICP1
7440-66-6	Zinc	11	18	100	08/21/07	8234	S8234A	34	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-012	% Solid:	93	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB06-1.5-2.5	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	220	5800	100	08/22/07	8234	S8234B	35	P	PEICPRAD1
7440-36-0	Antimony	2.2	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-38-2	Arsenic	2.2	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-39-3	Barium	11	13	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-41-7	Beryllium	0.65	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-43-9	Cadmium	0.65	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-70-2	Calcium	1100	ND	100	08/22/07	8234	S8234B	35	P	PEICPRAD1
7440-47-3	Chromium	5.4	6.8	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-48-4	Cobalt	2.7	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-50-8	Copper	5.4	6.3	100	08/21/07	8234	S8234A	35	P	PEICP1
7439-89-6	Iron	220	6700	100	08/22/07	8234	S8234B	35	P	PEICPRAD1
7439-92-1	Lead	5.4	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7439-95-4	Magnesium	540	850	100	08/22/07	8234	S8234B	35	P	PEICPRAD1
7439-96-5	Manganese	11	83	100	08/21/07	8234	S8234A	35	P	PEICP1
7439-97-6	Mercury	0.090	ND	167	08/21/07	8234	h8234s	29	CV	HGCV1
7440-02-0	Nickel	5.4	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-09-7	Potassium	540	ND	100	08/22/07	8234	S8234B	35	P	PEICPRAD1
7782-49-2	Selenium	1.9	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-22-4	Silver	1.6	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-23-5	Sodium	270	ND	100	08/22/07	8234	S8234B	35	P	PEICPRAD1
7440-28-0	Thallium	1.3	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-62-2	Vanadium	11	ND	100	08/21/07	8234	S8234A	35	P	PEICP1
7440-66-6	Zinc	11	11	100	08/21/07	8234	S8234A	35	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-013	% Solid:	94	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB05-1-3	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	210	5800	100	08/22/07	8234	S8234B	36	P	PEICPRAD1
7440-36-0	Antimony	2.1	ND	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-38-2	Arsenic	2.1	2.5	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-39-3	Barium	11	13	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-41-7	Beryllium	0.64	ND	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-43-9	Cadmium	0.64	ND	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-70-2	Calcium	1100	ND	100	08/22/07	8234	S8234B	36	P	PEICPRAD1
7440-47-3	Chromium	5.3	9.8	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-48-4	Cobalt	2.7	2.9	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-50-8	Copper	5.3	7.7	100	08/21/07	8234	S8234A	36	P	PEICP1
7439-89-6	Iron	210	6900	100	08/22/07	8234	S8234B	36	P	PEICPRAD1
7439-92-1	Lead	5.3	5.6	100	08/21/07	8234	S8234A	36	P	PEICP1
7439-95-4	Magnesium	530	930	100	08/22/07	8234	S8234B	36	P	PEICPRAD1
7439-96-5	Manganese	11	75	100	08/21/07	8234	S8234A	36	P	PEICP1
7439-97-6	Mercury	0.089	ND	167	08/21/07	8234	h8234s	30	CV	HGCV1
7440-02-0	Nickel	5.3	ND	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-09-7	Potassium	530	ND	100	08/22/07	8234	S8234B	36	P	PEICPRAD1
7782-49-2	Selenium	1.9	ND	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-22-4	Silver	1.6	ND	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-23-5	Sodium	270	ND	100	08/22/07	8234	S8234B	36	P	PEICPRAD1
7440-28-0	Thallium	1.3	ND	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-62-2	Vanadium	11	11	100	08/21/07	8234	S8234A	36	P	PEICP1
7440-66-6	Zinc	11	13	100	08/21/07	8234	S8234A	36	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-014	% Solid:	96	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB05-3-5	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	210	2700	100	08/22/07	8234	S8234B	37	P	PEICPRAD1
7440-36-0	Antimony	2.1	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-38-2	Arsenic	2.1	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-39-3	Barium	10	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-41-7	Beryllium	0.62	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-43-9	Cadmium	0.62	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-70-2	Calcium	1000	ND	100	08/22/07	8234	S8234B	37	P	PEICPRAD1
7440-47-3	Chromium	5.2	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-48-4	Cobalt	2.6	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-50-8	Copper	5.2	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7439-89-6	Iron	210	3700	100	08/22/07	8234	S8234B	37	P	PEICPRAD1
7439-92-1	Lead	5.2	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7439-95-4	Magnesium	520	ND	100	08/22/07	8234	S8234B	37	P	PEICPRAD1
7439-96-5	Manganese	10	53	100	08/21/07	8234	S8234A	37	P	PEICP1
7439-97-6	Mercury	0.087	ND	167	08/21/07	8234	h8234s	31	CV	HGCV1
7440-02-0	Nickel	5.2	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-09-7	Potassium	520	ND	100	08/22/07	8234	S8234B	37	P	PEICPRAD1
7782-49-2	Selenium	1.9	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-22-4	Silver	1.6	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-23-5	Sodium	260	ND	100	08/22/07	8234	S8234B	37	P	PEICPRAD1
7440-28-0	Thallium	1.2	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-62-2	Vanadium	10	ND	100	08/21/07	8234	S8234A	37	P	PEICP1
7440-66-6	Zinc	10	ND	100	08/21/07	8234	S8234A	37	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-015	% Solid:	98	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB05-5-7	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	200	430	100	08/22/07	8234	S8234B	38	P	PEICPRAD1
7440-36-0	Antimony	2.0	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-38-2	Arsenic	2.0	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-39-3	Barium	10	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-41-7	Beryllium	0.61	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-43-9	Cadmium	0.61	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-70-2	Calcium	1000	ND	100	08/22/07	8234	S8234B	38	P	PEICPRAD1
7440-47-3	Chromium	5.1	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-48-4	Cobalt	2.6	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-50-8	Copper	5.1	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7439-89-6	Iron	200	2200	100	08/22/07	8234	S8234B	38	P	PEICPRAD1
7439-92-1	Lead	5.1	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7439-95-4	Magnesium	510	ND	100	08/22/07	8234	S8234B	38	P	PEICPRAD1
7439-96-5	Manganese	10	52	100	08/21/07	8234	S8234A	38	P	PEICP1
7439-97-6	Mercury	0.085	ND	167	08/21/07	8234	h8234s	34	CV	HGCV1
7440-02-0	Nickel	5.1	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-09-7	Potassium	510	ND	100	08/22/07	8234	S8234B	38	P	PEICPRAD1
7782-49-2	Selenium	1.8	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-22-4	Silver	1.5	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-23-5	Sodium	260	ND	100	08/22/07	8234	S8234B	38	P	PEICPRAD1
7440-28-0	Thallium	1.2	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-62-2	Vanadium	10	ND	100	08/21/07	8234	S8234A	38	P	PEICP1
7440-66-6	Zinc	10	ND	100	08/21/07	8234	S8234A	38	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-016	% Solid:	98	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB05-7-9	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	200	680	100	08/22/07	8234	S8234B	39	P	PEICPRAD1
7440-36-0	Antimony	2.0	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-38-2	Arsenic	2.0	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-39-3	Barium	10	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-41-7	Beryllium	0.61	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-43-9	Cadmium	0.61	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-70-2	Calcium	1000	ND	100	08/22/07	8234	S8234B	39	P	PEICPRAD1
7440-47-3	Chromium	5.1	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-48-4	Cobalt	2.6	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-50-8	Copper	5.1	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7439-89-6	Iron	200	2100	100	08/22/07	8234	S8234B	39	P	PEICPRAD1
7439-92-1	Lead	5.1	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7439-95-4	Magnesium	510	ND	100	08/22/07	8234	S8234B	39	P	PEICPRAD1
7439-96-5	Manganese	10	75	100	08/21/07	8234	S8234A	39	P	PEICP1
7439-97-6	Mercury	0.085	ND	167	08/21/07	8234	h8234s	35	CV	HGCV1
7440-02-0	Nickel	5.1	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-09-7	Potassium	510	ND	100	08/22/07	8234	S8234B	39	P	PEICPRAD1
7782-49-2	Selenium	1.8	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-22-4	Silver	1.5	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-23-5	Sodium	260	ND	100	08/22/07	8234	S8234B	39	P	PEICPRAD1
7440-28-0	Thallium	1.2	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-62-2	Vanadium	10	ND	100	08/21/07	8234	S8234A	39	P	PEICP1
7440-66-6	Zinc	10	ND	100	08/21/07	8234	S8234A	39	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-017	% Solid:	96	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB02-5-7	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	210	1900	100	08/22/07	8234	S8234B	42	P	PEICPRAD1
7440-36-0	Antimony	2.1	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-38-2	Arsenic	2.1	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-39-3	Barium	10	25	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-41-7	Beryllium	0.62	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-43-9	Cadmium	0.62	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-70-2	Calcium	1000	24000	100	08/22/07	8234	S8234B	42	P	PEICPRAD1
7440-47-3	Chromium	5.2	13	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-48-4	Cobalt	2.6	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-50-8	Copper	5.2	9.4	100	08/21/07	8234	S8234A	42	P	PEICP1
7439-89-6	Iron	210	3500	100	08/22/07	8234	S8234B	42	P	PEICPRAD1
7439-92-1	Lead	5.2	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7439-95-4	Magnesium	520	7800	100	08/22/07	8234	S8234B	42	P	PEICPRAD1
7439-96-5	Manganese	10	70	100	08/21/07	8234	S8234A	42	P	PEICP1
7439-97-6	Mercury	0.087	ND	167	08/21/07	8234	h8234s	36	CV	HGCV1
7440-02-0	Nickel	5.2	6.6	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-09-7	Potassium	520	ND	100	08/22/07	8234	S8234B	42	P	PEICPRAD1
7782-49-2	Selenium	1.9	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-22-4	Silver	1.6	3.4	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-23-5	Sodium	260	ND	100	08/22/07	8234	S8234B	42	P	PEICPRAD1
7440-28-0	Thallium	1.2	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-62-2	Vanadium	10	ND	100	08/21/07	8234	S8234A	42	P	PEICP1
7440-66-6	Zinc	10	ND	100	08/21/07	8234	S8234A	42	P	PEICP1

Comments: _____

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Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-018	% Solid:	97	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-SB02-7-9	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	210	650	100	08/22/07	8234	S8234B	43	P	PEICPRAD1
7440-36-0	Antimony	2.1	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-38-2	Arsenic	2.1	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-39-3	Barium	10	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-41-7	Beryllium	0.62	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-43-9	Cadmium	0.62	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-70-2	Calcium	1000	ND	100	08/22/07	8234	S8234B	43	P	PEICPRAD1
7440-47-3	Chromium	5.2	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-48-4	Cobalt	2.6	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-50-8	Copper	5.2	5.5	100	08/21/07	8234	S8234A	43	P	PEICP1
7439-89-6	Iron	210	1400	100	08/22/07	8234	S8234B	43	P	PEICPRAD1
7439-92-1	Lead	5.2	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7439-95-4	Magnesium	520	ND	100	08/22/07	8234	S8234B	43	P	PEICPRAD1
7439-96-5	Manganese	10	53	100	08/21/07	8234	S8234A	43	P	PEICP1
7439-97-6	Mercury	0.086	ND	167	08/21/07	8234	h8234s	37	CV	HGCV1
7440-02-0	Nickel	5.2	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-09-7	Potassium	520	ND	100	08/22/07	8234	S8234B	43	P	PEICPRAD1
7782-49-2	Selenium	1.9	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-22-4	Silver	1.5	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-23-5	Sodium	260	ND	100	08/22/07	8234	S8234B	43	P	PEICPRAD1
7440-28-0	Thallium	1.2	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-62-2	Vanadium	10	ND	100	08/21/07	8234	S8234A	43	P	PEICP1
7440-66-6	Zinc	10	ND	100	08/21/07	8234	S8234A	43	P	PEICP1

Comments: _____

_____

Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

Form1
Inorganic Analysis Data Sheet

Sample ID:	AC32378-019	% Solid:	82	Lab Name:	Veritech	Nras No:	
Client Id:	CPC-DW04-13-14	Units:	MG/KG	Lab Code:		Sdg No:	
Matrix:	SOIL	Date Rec:	8/17/2007	Contract:		Case No:	
Level:	LOW						

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	240	2300	100	08/22/07	8234	S8234B	44	P	PEICPRAD1
7440-36-0	Antimony	2.4	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-38-2	Arsenic	2.4	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-39-3	Barium	12	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-41-7	Beryllium	0.73	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-43-9	Cadmium	0.73	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-70-2	Calcium	1200	ND	100	08/22/07	8234	S8234B	44	P	PEICPRAD1
7440-47-3	Chromium	6.1	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-48-4	Cobalt	3.0	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-50-8	Copper	6.1	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7439-89-6	Iron	240	4400	100	08/22/07	8234	S8234B	44	P	PEICPRAD1
7439-92-1	Lead	6.1	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7439-95-4	Magnesium	610	1100	100	08/22/07	8234	S8234B	44	P	PEICPRAD1
7439-96-5	Manganese	12	36	100	08/21/07	8234	S8234A	44	P	PEICP1
7439-97-6	Mercury	0.10	ND	167	08/21/07	8234	h8234s	38	CV	HGCV1
7440-02-0	Nickel	6.1	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-09-7	Potassium	610	810	100	08/22/07	8234	S8234B	44	P	PEICPRAD1
7782-49-2	Selenium	2.2	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-22-4	Silver	1.8	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-23-5	Sodium	300	ND	100	08/22/07	8234	S8234B	44	P	PEICPRAD1
7440-28-0	Thallium	1.5	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-62-2	Vanadium	12	ND	100	08/21/07	8234	S8234A	44	P	PEICP1
7440-66-6	Zinc	12	17	100	08/21/07	8234	S8234A	44	P	PEICP1

Comments: _____

_____

Flag Codes:

U or ND - Indicates Compound was not found above the detection/reporting limit

VERITECH Wet Chem Form1 Analysis Summary

Lab#: AC32378-001
Matrix Soil
Client SampleID: CPC-DW01-10-11

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.29	08/25/07	08/25/07

Lab#: AC32378-002
Matrix Soil
Client SampleID: CPC-DW02-10-11

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.3	08/25/07	08/25/07

Lab#: AC32378-003
Matrix Soil
Client SampleID: CPC-DW03-10-11

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.28	08/25/07	08/25/07

Lab#: AC32378-004
Matrix Soil
Client SampleID: CPC-DW04-8-9

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.29	08/25/07	08/25/07

Lab#: AC32378-005
Matrix Soil
Client SampleID: CPC-SOIL-Duplicate

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.27	08/25/07	08/25/07

Lab#: AC32378-006
Matrix Soil
Client SampleID: CPC-SB01-1.5-2.5

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.26	08/25/07	08/25/07

Lab#: AC32378-007
Matrix Soil
Client SampleID: CPC-SB02-1-5

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.26	08/25/07	08/25/07

Lab#: AC32378-008
Matrix Soil
Client SampleID: CPC-SB03-1.5-3.5

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.27	08/25/07	08/25/07

Lab#: AC32378-009
Matrix Soil
Client SampleID: CPC-SB03-1.5-3.5MS

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	5.4	mg/kg	0.26	08/25/07	08/25/07

VERITECH Wet Chem Form1 Analysis Summary

Lab#: AC32378-010 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB03-1.5-3.5MSD | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	5.4	mg/kg	0.27	08/25/07	08/25/07

Lab#: AC32378-011 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB04-1.5-2.5 | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.28	08/25/07	08/25/07

Lab#: AC32378-012 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB06-1.5-2.5 | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.27	08/25/07	08/25/07

Lab#: AC32378-013 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB05-1-3 | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.26	08/25/07	08/25/07

Lab#: AC32378-014 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB05-3-5 | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.26	08/25/07	08/25/07

Lab#: AC32378-015 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB05-5-7 | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.25	08/25/07	08/25/07

Lab#: AC32378-016 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB05-7-9 | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.25	08/25/07	08/25/07

Lab#: AC32378-017 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB02-5-7 | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	0.29	mg/kg	0.26	08/25/07	08/25/07

Lab#: AC32378-018 | Project Number: 7081622
Matrix: Soil | Received Date: 8/16/2007
Client SampleID: CPC-SB02-7-9 | Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.26	08/25/07	08/25/07

VERITECH Wet Chem Form1 Analysis Summary

Lab#: AC32378-019
Matrix Soil
Client SampleID: CPC-DW04-13-14

Project Number: 7081622
Received Date: 8/16/2007
Collect Date: 8/16/2007

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Cyanide	CN-SOIL	1	ND	mg/kg	0.3	08/25/07	08/25/07

ATTACHMENT B

SDG Narrative

Client: EA Engineering Science & Technology
Project: CPC

Hampton-Clarke/Veritech (HC·V) received the following samples from EA Engineering on August 15, 2007:

EA#	HC·V #	Type	Analysis
CPC-MW01	AC32351-001	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW02	AC32351-002	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW03	AC32351-003	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW04	AC32351-004	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW05	AC32351-005	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW06	AC32351-006	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW07	AC32351-007	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW09	AC32351-008	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW10	AC32351-009	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW10MS	AC32351-010	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-MW10MSD	AC32351-011	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
CPC-Duplicate	AC32351-012	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)
TB	AC32351-013	Aqueous	VO (8260B)

Problems associated with the analysis of these samples:

Volatiles:

Methylene chloride was recovered in sample AC32351-013 due to possible laboratory contamination.

Sample AC32351-004 was analyzed at a 10x dilution.

Sample AC32351-006 was analyzed twice to confirm high surrogate recoveries. Also one surrogate exceeded QC limit in the MSD for batch MBS6026.

The RPD (29%) for Chloroform exceeded the QC criteria in batch MBS6026.
Trichloroethene was recovered outside the criteria in the MS (64%) and MSD (63%).

There were no other problems associated with this analysis.

Semi-Volatiles:

There were no problems associated with this analysis.

PCB:

There were no problems associated with this analysis.

Pesticides:

There were no problems associated with this analysis.

Metals:

Manganese was recovered below QC criteria in the MS (59%) and MSD (60%). The LCS and LCS MR met all QC criteria for batch 8241.

There were no other problems associated with this analysis.

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or his designee, as verified by the following signature.

Jeri Rossi
Quality Assurance Director

Or

Stanley Gilewicz
Laboratory Director

9/7/07
Date

SDG Narrative

Client: EA Engineering Science & Technology
Project: CPC

Hampton-Clarke/Veritech (HC·V) received the following samples from EA Engineering on August 16, 2007:

EA#	HC·V #	Type	Analysis
CPC-DW01-10-11	AC32378-001	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-DW02-10-11	AC32378-002	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-DW03-10-11	AC32378-003	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-DW04-8-9	AC32378-004	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-Soil-Dup	AC32378-005	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB01-1.5-2.5	AC32378-006	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB02-1-5	AC32378-007	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB03-1.5-3.5	AC32378-008	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC- SB03-1.5-3.5MS	AC32378-009	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC- SB03-1.5-3.5MSD	AC32378-010	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB04-1-2.5	AC32378-011	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB06-1.5-2.5	AC32378-012	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB-05-1-3	AC32378-013	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB-05-3-5	AC32378-014	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB-05-5-7	AC32378-015	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB-05-7-9	AC32378-016	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB-02-5-7	AC32378-017	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-SB-02-7-9	AC32378-018	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-DW04-13-14	AC32378-019	Soil	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A), CN (SM4500E)
CPC-Rinsate	AC32378-020	Aqueous	VO (8260B), BNA (8270C), PCB (8082), PEST (8081A), TAL Metals (6010B/7471A)

Problems associated with the analysis of these samples:

Volatiles:

Methylene chloride was recovered in method blanks 1M24245, 1M24283 and 2M23108 and in samples AC32378-001-008, 011-019 due to possible laboratory contamination.

Sample AC32378-012 was analyzed at a 5x dilution.

The RPD for most of the compounds exceeded the QC criteria in batch MBS6041. However, the MBS, MS and MSD were met all QC criteria in this batch.
Carbon Tetrachloride was recovered outside the criteria in the MS (193%) and MSD (180%) for batch MBS5923.

There were no other problems associated with this analysis.

Semi-Volatiles:

Sample AC32378-001was analyzed at a 2x dilution.

There were no other problems associated with this analysis.

PCB:

There were no problems associated with this analysis.

Pesticides:

There were no problems associated with this analysis.

Metals:

The following is a summary of batch QC:

8234: The MS (0%) for Iron and MSD (140%) for Aluminum were recovered outside the QC limit. The LCS and LSC MR were met all QC criteria for this batch.
The serial dilution exceeded the criteria for Beryllium (22%) and Chromium (19%) suggesting possible matrix interference.

8269: Two elements were recovered outside the QC criteria in the MS/MSD: Chromium (158%/64%) and Manganese (132%/128%). The LCS and LSC MR were met all QC criteria for this batch.
The serial dilution exceeded the criteria for Vanadium (21%) and Zinc (14%) suggesting possible matrix interference

There were no other problems associated with this analysis.

Wet Chemistry:

There were no other problems associated with this analysis.

JR
9/10/2

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or his designee, as verified by the following signature.

Jeri Rossi	Or	Stanley Gilewicz	9/10/02
Quality Assurance Director		Laboratory Director	Date

Veritech/Division of Hampton-Clarke
175 US Hwy 46 West, Fairfield, New Jersey 07004 Ph: 800-426-9992 fax:973-439-1458

CHAIN OF CUSTODY RECORD

Project#(Lab Use Only) 7081532

Page 1 of 1

Customer Information

1a) Customer: EA Engineering
Address: 6712 Brooklawn Pkwy
Syracuse NY 13211
1b) Email/Cell/Fax/Ph: JVonUderitz@EAEst.com
1c) Send Invoice To: EA
1d) Send Report To: Jim Hayward

Project Information

2a) Project: CPC
2b) Project Manager: Jim Hayward
2c) Location (City/State): Bay Shore, NY
2d) Quote#/PO# (If Applicable):

3) Reporting Requirements(please circle)

Turnaround Time	Report type	Electronic Deliv
24-Hour(100%)	Data Sum	Hazsite/Csv
48-Hour(75%)	Waste	Equis
72-Hour(50%)	Red-NJ/NY/PA	Excel-NJCC
1-Week(25%)	CLP	Excel-Nytagm
10 Days(10%)	Full/Cat-B	Excel-PAActII
Standard	Cat-A	PDF
Other:	Other:	Other:

Expedited TAT Not always available (Please check with lab)!

7) Analysis Request

FOR LAB USE ONLY
Batch# AC32351

Check if Contingent===> <===Check if Contingent

Matrix Codes: DW-Drinking Water, GW-Ground Water, WW-Waste Water, S-Soil, SL-Sludge, O-Oil, A-Air, Ot-Other

Lab Sample#	4) Customer Sample ID	5) Matrix	6) Sample Date	6) Sample Time	Sample Type: Composite(C)	Sample Type: Grab(G)	VOC	SVOC	PCB/Pest	TAL metals	8) # Of Bottles: None	MeOH	Encore	NaOH	HCl	H2SO4	HNO3	Other	9) Methanol Bottle Numbers (If applicable)
-001	CPC-MW01	GW	8/14/07	1300		X	X	X	X	X	4				2		1		
-002	CPC-MW08	GW	8/14/07	850		X	X	X	X	X	4				2		1		
-003	CPC-MW03	GW	8/14/07	1030		X	X	X	X	X	4				2		1		
-004	CPC-MW04	GW	8/14/07	740		X	X	X	X	X	4				2		1		
-005	CPC-MW05	GW	8/14/07	1140		X	X	X	X	X	4				2		1		
-006	CPC-MW06	GW	8/14/07	1130		X	X	X	X	X	4				2		1		
-007	CPC-MW07	GW	8/14/07	1300		X	X	X	X	X	4				2		1		
-008	CPC-MW09 (DUP)	GW	8/14/07	950		X	X	X	X	X	4				2		1		
-009	CPC-MW10 (MS/MSD)	GW	8/14/07	1210		X	X	X	X	X	4				2		1		
~~-010~~ 013	~~CPC-Duplicate~~ DUP	GW	8/14/07	—		X	X	X	X	X	4				2		1		

10) Relinquished By:	Accepted By	Date	Time
		8/15/07	1200
		8/15/07	1440

AC32351 Comments, Notes, Special Requirements, HAZARDS
~~010~~ 013 TB AQ 8/14/7 VD MS/MSD for S.VOA

See CLR
Cooler Tmp

11) Sampler: Date: 8/14/07

Please note NUMBERED items. If not completed your analytical work may be delayed.
A fee of $6/sample will be assessed for storage should sample not be activated for any analysis

Veritech/Division of Hampton-Clarke
175 US Hwy 46 West, Fairfield, New Jersey 07004 Ph: 800-426-9992 fax:973-439-1458

CHAIN OF CUSTODY RECORD

Project#(Lab Use Only) 7081622

Customer Information

1a) Customer: EA
Address: 6712 Brooklawn Pkw
Syracuse NY 13211
1b) Email/Cell/Fax/Ph: JVanLuvender2@EAEst.com
1c) Send Invoice To: EA 6712 Brooklawn Pk
Syracuse NY 13211
1d) Send Report To: Jim Hayward

Project Information

2a) Project: CPC
2b) Project Manager: Jim Hayward
2c) Location (City/State): Bay Shore, NY
2d) Quote#/PO# (If Applicable):

3) Reporting Requirements(please circle)

Tumaround Time	Report type	Electronic Deliv
24-Hour(100%)	Data Sum	Hazsite/Csv
48-Hour(75%)	Waste	Equis
72-Hour(50%)	Red-NJ/NY/PA	Excel-NJCC
1-Week(25%)	CLP	Excel-Nytagm
10 Days(10%)	Full/Cat-B	Excel-PAActII
Standard	Cat-A	PDF
Other:	Other:	Other:

Expedited TAT Not always available (Please check with lab)!

7) Analysis Request

Check if Contingent===> <===Check if Contingent

FOR LAB USE ONLY
Batch# AC32378

Matrix Codes: DW-Drinking Water, GW-Ground Water, WW-Waste Water, S-Soil, SL-Sludge, O-Oil, A-Air, Ot-Other

Lab Sample#	4) Customer Sample ID	5) Matrix	6) Sample Date	6) Sample Time	Sample Type: Composite(C)	Sample Type: Grab(G)	VOC	SVOC	Heavy Metals-TAL	PCBs	Pesticides	CN	8) # Of Bottles: None	MeOH	Encore	NaOH	HCl	H_2SO_4	HNO_3	Other:	9) Methanol Bottle Numbers (If applicable) / Comments
001	CPC-DW01-10-11	S	8/16/07	750		X	X	X	X	X	X	X	2								
002	CPC-DW02-10-11	S	8/16/07	830		X	X	X	X	X	X		2								
003	CPC-DW03-10-11	S	8/16/07	845		X	X	X	X	X	X		2								
004	CPC-DW04-8-9	S	8/16/07	1130		X	X	X	X	X	X		2								
005	CPC-SO1 Duplicate	S	8/16/07	—		X	X	X	X	X	X		2								
006	CPC-SB01-1.5-2.5	S	8/16/07	1115		X	X	X	X	X	X		2								
007	CPC-SB02-1-3	S	8/16/07	1045		X	X	X	X	X	X		2								
008 009 010	CPC-SB03-1.5-3.5 (MS/MSD)	S	8/16/07	1030		X	X	X	X	X	X		2								
011	CPC-SB04-1.5-2.5	S	8/16/07	915		X	X	X	X	X	X		2								
012	CPC-SB06-1.5-2.5	S	8/16/07	1000		X	X	X	X	X	X		2								

10) Relinquished By:	Accepted By	Date	Time
J. Hayward	Ricardo Lima	8/16/07	15:00

Comments, Notes, Special Requirements, HAZARDS

Cooler Tmp: 3°

11) Sampler: J. Hayward Date: 8/16/07

Please note NUMBERED items. If not completed your analytical work may be delayed.
A fee of $5/sample will be assessed for storage should sample not be activated for any analysis

Veritech/Division of Hampton-Clarke — **CHAIN OF CUSTODY RECORD** — Project#(Lab Use Only) — Page 2 of 2

175 US Hwy 46 West, Fairfield, New Jersey 07004 Ph: 800-426-9992 fax:973-439-1458

Customer Information

1a) Customer: EA
Address: 6712 Brooklawn Pkw Syracuse NY 13211
1b) Email/Cell/Fax/Ph: JVanUdenitz@EAEst.com
1c) Send Invoice To: EA 6712 Brooklawn Pkw Syracuse, NY 13211
1d) Send Report To: Jim Hayward

Project Information

2a) Project: CPC
2b) Project Manager: Jim Hayward
2c) Location (City/State): Bay Shore, NY
2d) Quote#/PO# (If Applicable):

3) Reporting Requirements(please circle)

Turnaround Time	Report type	Electronic Deliv
24-Hour(100%)	Data Sum	Hazsite/Csv
48-Hour(75%)	Waste	Equis
72-Hour(50%)	Red-NJ/NY/PA	Excel-NJCC
1-Week(25%)	CLP	Excel-Nytagm
10 Days(10%)	Full/Cat-B	Excel-PAActII
Standard	Cat-A	PDF
Other:	Other:	Other:

Expedited TAT Not always available (Please check with lab)!

7) Analysis Request

FOR LAB USE ONLY — Batch# AC32378

Check if Contingent===> <===Check if Contingent

Matrix Codes: DW-Drinking Water; GW-Ground Water; WW-Waste Water; S-Soil; SL-Sludge; O-Oil; A-Air; Ot-Other

Lab Sample#	4) Customer Sample ID	5) Matrix	6) Sample Date	6) Sample Time	Sample Type: Composite(C)	Sample Type: Grab(G)	VOC	SVOC	Heavy Metals	PCB's	Pesticides	ON HOLD 8/23/07 Per JVU	8) # Of Bottles: None	MeOH	Encore	NaOH	HCl	H2SO4	HNO3	Other	9) Methanol Bottle Numbers (If applicable) / Comments
-013	CPC-SB05-1-3	S	8/16/07	930		X	X	X	X	X	X	X	2								
-014	CPC-SB05-3-5	S	8/16/07	935		X	X	X	X	X	X		2								
-015	CPC-SB05-5-7	S	8/16/07	940		X	X	X	X	X	X		2								
-016	CPC-SB05-7-9	S	8/16/07	945		X	X	X	X	X	X		2								
-017	CPC-SB02-5-7	S	8/16/07	1050		X	X	X	X	X	X		2								
-018	CPC-SB02-7-9	S	8/16/07	1055		X	X	X	X	X	X		2								
-019	CPC-DW04-13-14	S	8/16/07	1145		X	X	X	X	X	X	1	2								
-020	CPC-Rinsate	OT	8/16/07	1200		X	X	X	X	X	X		4				2		1		
			8/17/07																		

10) Relinquished By:	Accepted By	Date	Time
J. Hayward	Records Clerk	8/16/07	15:00

Comments, Notes, Special Requirements, HAZARDS

11) Sampler: J. Hayward Date: 8/16/07

Cooler Tmp: 3.1

Please note NUMBERED items. If not completed your analytical work may be delayed.
A fee of $5/sample will be assessed for storage should sample not be activated for any analysis