



**NEW YORK STATE DEPARTMENT OF TRANSPORTATION
ALBANY, NEW YORK**

**HARRISON SUBRESIDENCY
WESTCHESTER COUNTY
POST-CLOSURE MONITORING RESULTS**

JULY 2007, FIFTH QUARTER SAMPLING

February 2008

HDR|LMS

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**HARRISON SUBRESIDENCY
WESTCHESTER COUNTY
POST-CLOSURE QUARTERLY MONITORING RESULTS
FIFTH QUARTER SAMPLING
JULY 2007**

EXECUTIVE SUMMARY

This report presents the results of the Fifth Quarter 2007 post-closure sampling and monitoring conducted at the Harrison Subresidency site located in the town of Harrison, Westchester County, New York on July 20, 2007. The objectives of the post-closure sampling and monitoring program are to; 1) evaluate the environmental impacts of the landfill, if any; 2) meet the post-closure monitoring requirements of the NYSDEC and; 3) provide the NYSDEC with data to evaluate and/or modify the existing sampling and monitoring program. Five wells (LMW-2, LMW-4, PC-1, PC-2, and PC-3), three surface water locations (SW-1, SW-2, and SW-4) and four sediment locations (SD-1, SD-2, SD-3 and SD-4) were sampled for Target Analyte List (TAL) metals and chloride, volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs) and cyanide. The stream at one surface water location, SW-3, was dry and therefore was not sampled.

Overall, based on groundwater and surface water on-site/background data, the landfill does not appear to be contributing significant levels of contaminants to the groundwater or surface water, as we have observed in the past. Iron, manganese, and sodium concentrations are elevated in the downgradient/on-site well samples, which indicates a contribution from the landfill. Concentrations of iron and manganese in the downgradient/off-site sample are much less, indicating these metals are attenuating as they are migrating off-site.

The landfill may locally contribute or have contributed to elevated levels of lead in the on-site sediment (SD-3); however, on-site lead concentrations downgradient (SD-4) are much lower; therefore, lead does not appear to be significantly migrating off-site.

1.0 INTRODUCTION

1.1 Background

This report presents the results of the July 2007 Fifth Quarter post-closure sampling and monitoring conducted at the Harrison Subresidency site located in the town of Harrison, Westchester County, New York (Figure 1). The site, once a seasonal highway maintenance support and salt storage facility operated by the New York State Department of Transportation (NYSDOT), is now occupied by the Town of Harrison. The site includes approximately 2.6 acres of landfill area (Figure 2) that was closed in December 1998. The Fifth Quarter sampling and monitoring was conducted to evaluate the environmental impacts of landfill closure, if any, through groundwater, surface water and sediment sampling, gas monitoring and a landfill inspection. Sampling was conducted in accordance with the NYSDEC-approved Operation and Maintenance Plan for the Harrison Sub-residency Landfill and Spill Area dated, February 2007 (NYSDOT 2007).

1.2 Monitoring Objectives

The objectives of the post-closure sampling and monitoring program are to; 1) evaluate the environmental impacts of the landfill, if any; 2) meet the post-closure monitoring requirements of the NYSDEC and; 3) provide NYSDEC with data to evaluate and/or modify the existing sampling and monitoring program.

2.0 FIELD INVESTIGATION

2.1 Monitoring Well Sampling

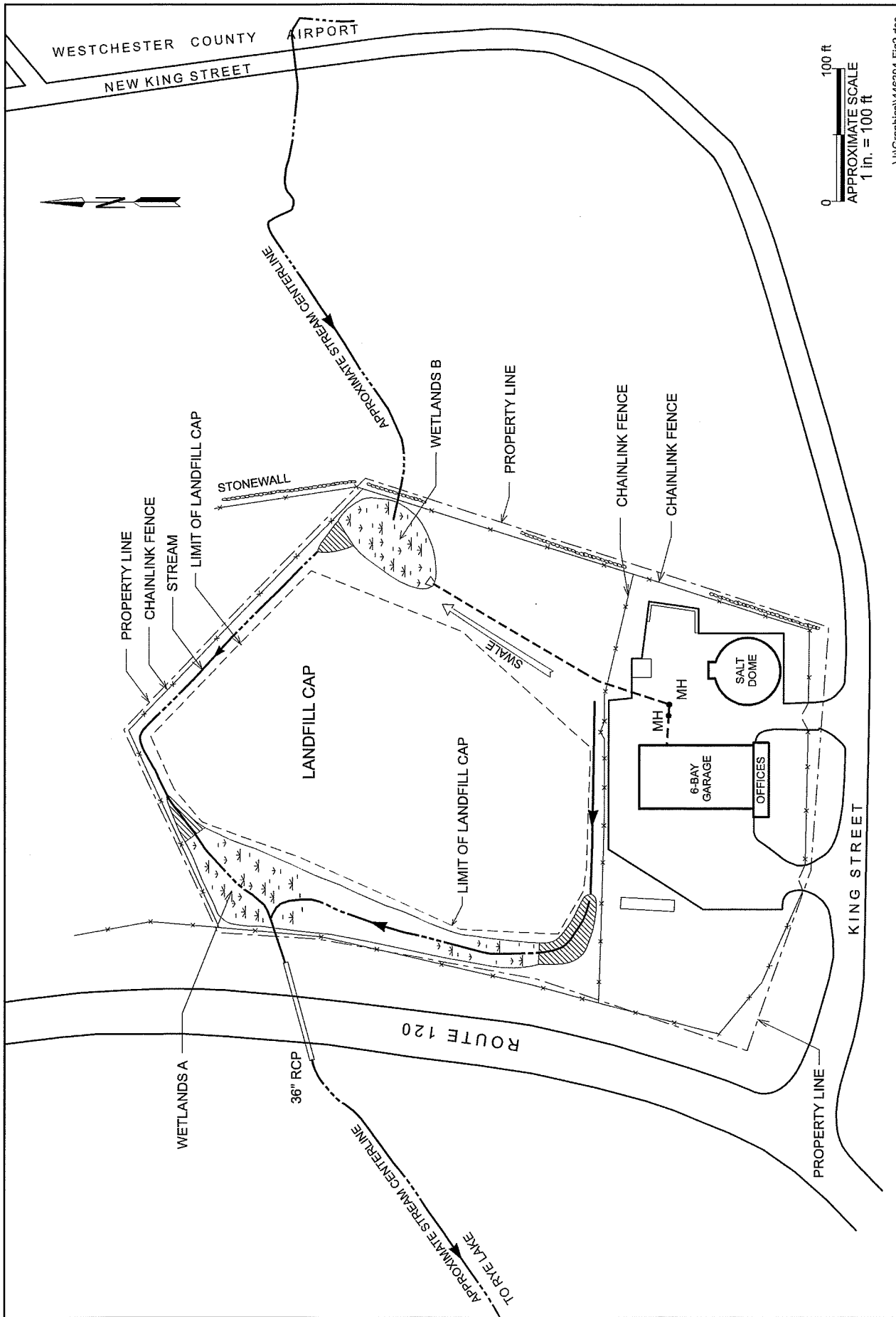
On July 20, 2007, as part of the quarterly post-closure sampling and monitoring program, five monitoring wells (Figure 3) were purged and sampled for TAL metals (filtered) and chloride, volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), and cyanide. Samples were collected using dedicated disposable bailers. The samples were preserved on ice to 4°C and sent, under chain of custody, to a New York State Department of Health (NYSDOH)-approved laboratory for analysis.

A round of static water level measurements was recorded for all wells prior to sampling activities. Groundwater chemistry measurements (temperature, pH, conductivity, and turbidity) were recorded before, during, and after purging, with a measurement recorded for approximately every well volume. An additional round of static water measurements



Name: GLENVILLE
 Date: 1/23/2008
 Scale: 1 inch equals 2000 feet

Location: 041° 04' 00.92" N 073° 42' 27.38" W NAD 27
 Caption: Figure 1 - Site Location
 Harrison Subresidency Site



0 100 ft
 APPROXIMATE SCALE
 1 in. = 100 ft

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Figure 2

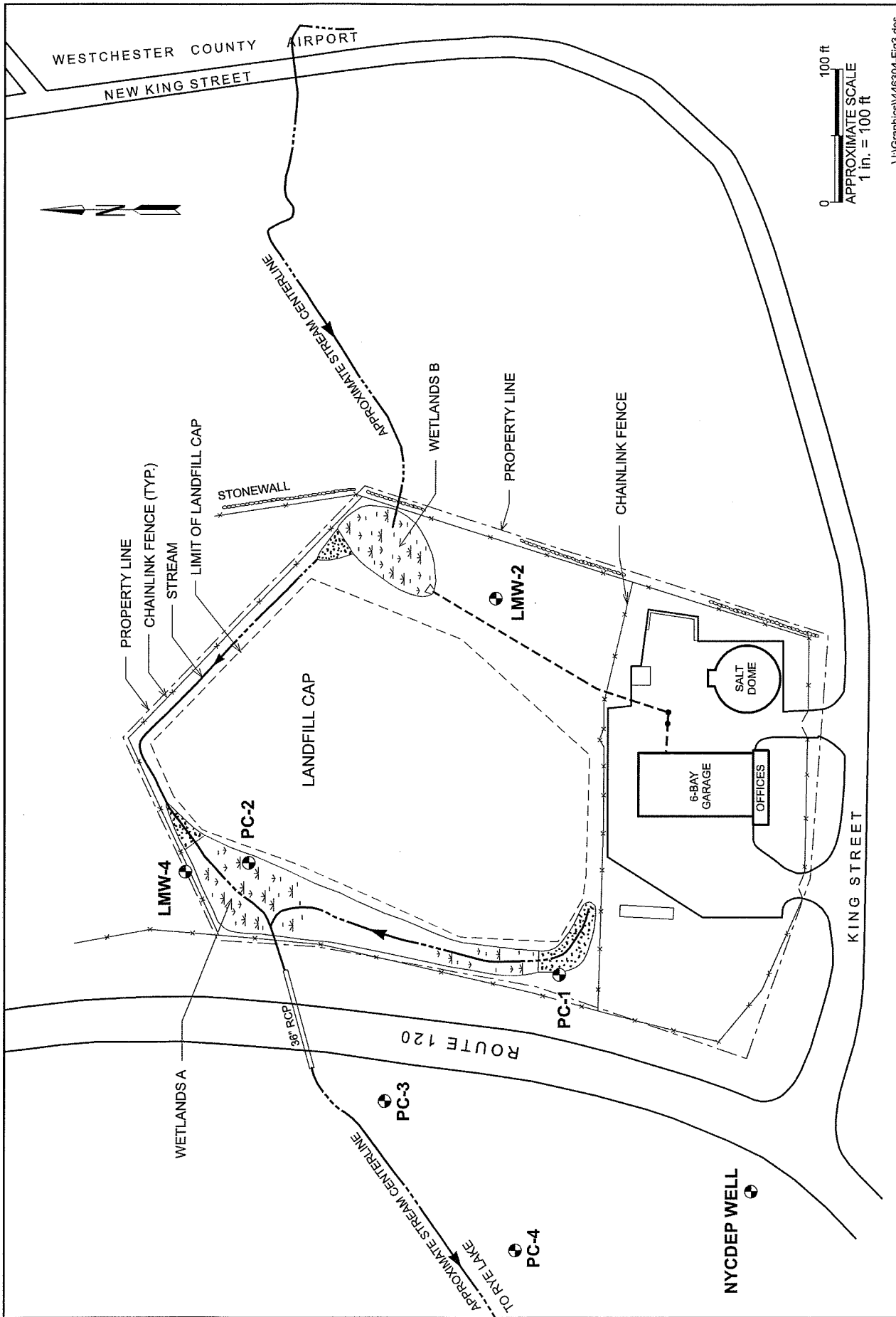
Site Map

NYS DOT PIN: 8806.51.301

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Figure 3

Groundwater Sample Locations

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was recorded for all wells prior to sampling. Groundwater purging information is recorded on the groundwater well sampling logs included in Attachment A.

2.2 Surface Water Sampling

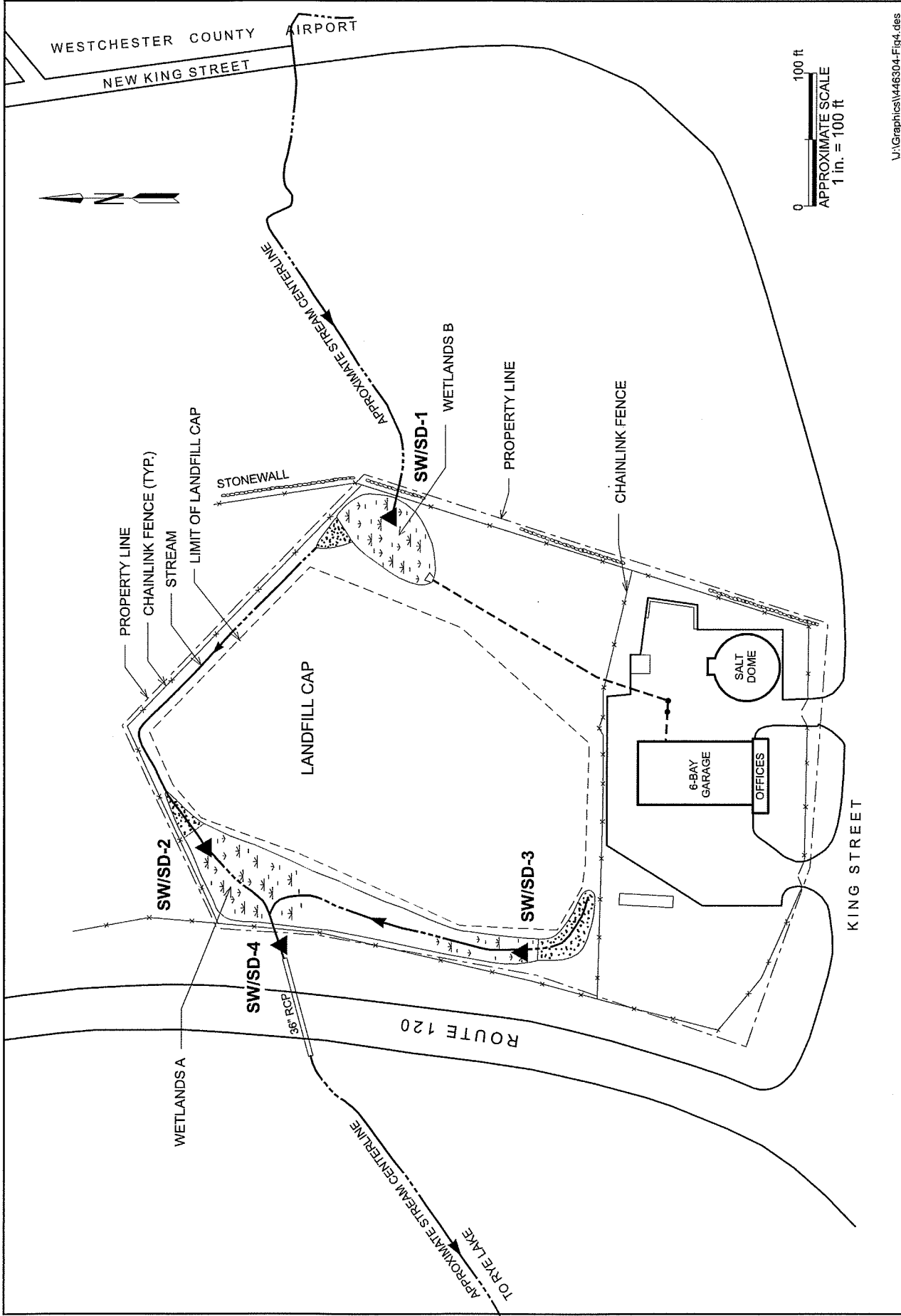
Three surface water locations (SW-1, SW-2 and SW-4) were sampled for TAL metals (total) and chloride, VOCs, SVOCs and cyanide. A sample was not collected from SW-3 (Figure 4) because the stream at this location was dry. All surface water samples were collected prior to sediment sampling to prevent increased turbidity in the surface water.

SW-1 was collected from a point located at the eastern section of wetlands B. SW-2 was collected from a point located at the northern portion of wetlands A (western side of the landfill). SW-4 was collected from a point located approximately 17 ft. northeast of the 36-in. reinforced concrete pipe (RCP) culvert that diverts the stream southwest under Route 120 to Rye Lake.

The samples were collected by dipping a dedicated laboratory-cleaned stainless steel ladle into the water and transferring the sample to the appropriate pre-cleaned laboratory-supplied container. The containers were iced to 4°C and sent, under chain of custody, to a NYSDOH-approved laboratory for analysis. Water chemistry measurements (temperature, pH, conductivity, and turbidity) were recorded during sample collection and are included in Attachment B.

2.3 Sediment Sampling

Four sediment samples (Figure 4) and one sediment field duplicate sample of SD-4 (labeled SD-5) were collected for TAL metals and chloride, VOCs, SVOCs, and cyanide analyses. Each sediment sample was collected subsequent to, and at the same location as, its corresponding surface water sample. The samples were collected using a dedicated laboratory-cleaned stainless steel trowel and were placed directly into the appropriate pre-cleaned laboratory-supplied sample container. Each sample container was iced to 4°C and sent, under chain of custody, to a NYSDOH-approved laboratory for analysis. The sample depth, texture, color and odor were noted and are included in Attachment B.



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Surface Water and Sediment Sample Locations

Figure 4

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2.4 Gas Monitoring

In accordance with the NYSDEC-approved Operation and Maintenance Plan for the Harrison Sub-residency Landfill and Spill Area, dated February 2007 (NYSDOT 2007), a gas monitoring program was instituted to verify that any gases, produced as a result of the natural decomposition of waste, do not pose a hazard to health or safety. The program includes the measurement of concentrations of methane or other explosive gases, hydrogen sulfide and volatile organic compounds (VOCs) at each of four gas vents and around the perimeter of the landfill (Figure 5).

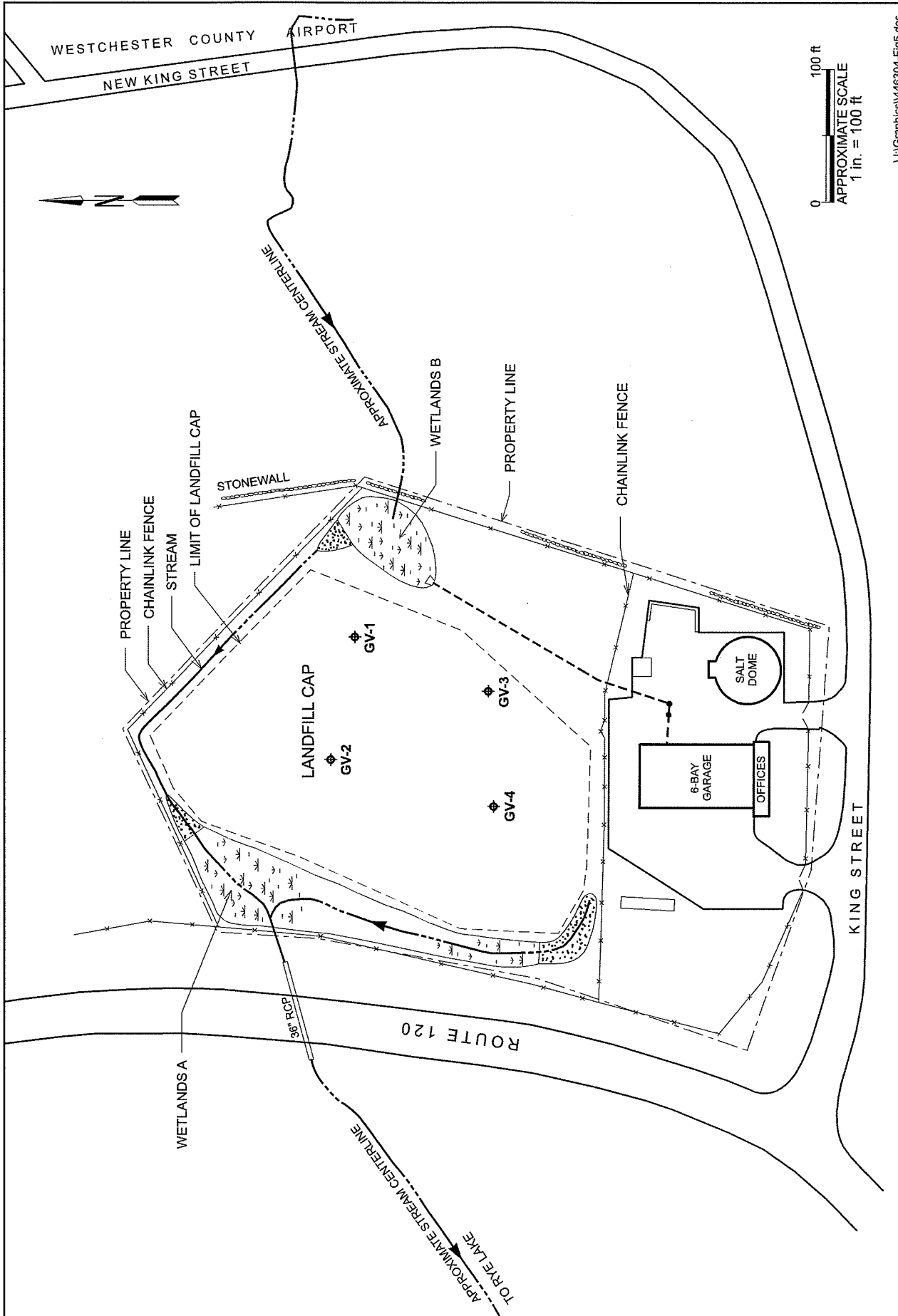
Each of the four gas vents, located around the perimeter of the landfill, were monitored for methane and other explosive gases with a combustible gas indicator (CGI). Gas vent readings were obtained by inserting the instrument detector probe into the vent. The CGI was set to sound an alarm if the readings exceeded 10 % of the lower explosive limit (LEL) of methane. In addition, alarms were set at 10% of the LEL of hydrogen sulfide, 25 parts per million (ppm) of carbon monoxide and 19.5% and 23.5% of oxygen. There were no readings exceeding the percent LEL for methane and preset alarm levels for hydrogen sulfide, carbon monoxide and oxygen.

A photoionization detector (PID) and a flame ionization detector (FID) (with and without a methane filter) were used to monitor for VOCs around the perimeter of the landfill and at each of the four gas vents. There were no readings above background levels.

An air monitoring field data sheet is included in Attachment C.

2.5 Inspections

All five groundwater monitoring wells were inspected and, with the exception of PC-2, were found to be in good condition. The well casing at PC-2 had previously sheared sideways allowing only a gap of approximately two centimeters down the well. The landfill cap, drainage swales, and gas vents were also inspected and with the exception of the presence of signs of erosion at the western side of the landfill, all appeared to be in good condition. No vermin or vector was observed on the landfill.



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Gas Vent Locations

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Figure 5

3.0 ANALYTICAL RESULTS

3.1 Groundwater Results

Groundwater samples were collected from four on-site monitoring wells (LMW-2, LMW-4, PC-1 and PC-2) and one off-site well (PC-3).

Groundwater samples were analyzed according to NYSDEC Analytical Services Protocol (ASP). Analytical results for filtered TAL metals and chloride, VOCs, SVOCs and cyanide are presented in Table 1. A copy of the analytical laboratory data is presented in Attachment D. Field parameters for temperature, pH, specific conductance, and turbidity are provided on the groundwater well sampling logs included in Attachment A.

3.2 Surface Water Results

Three surface water samples were collected and analyzed for TAL metals (unfiltered) and chloride, VOCs, SVOCs and cyanide according to NYSDEC ASP. Analytical results are presented in Table 2. A copy of the analytical laboratory data is presented in Attachment D. Field parameters for temperature, pH, specific conductance, and turbidity are provided on the surface water sampling logs included in Attachment B.

3.3 Sediment Data Results

Four sediment samples and one field duplicate sample were collected and analyzed for TAL metals and chloride, VOCs, SVOCs and cyanide according to NYSDEC ASP. Analytical results are presented in Table 3. A copy of the analytical laboratory data is presented in Attachment D. Sample depths and field observations are provided on the sediment sampling logs included in Attachment B.

3.31 Field Duplicate Results

The relative percent difference (RPD) of the duplicate data for SD-4 (Table 4) indicates that arsenic, barium, beryllium, calcium, copper, iron, lead, manganese, potassium, silver, acetone and carbon disulfide exceeded the normally accepted range of 15%. A review of the laboratory data indicates that the data for these parameters was accurately reported, as per standard reporting protocols. The field crew collected the original and duplicate samples according to the appropriate field duplicate sampling protocols (i.e. one sample

TABLE 1
GROUNDWATER DATA SUMMARY
 Fifth Quarter Sampling - Harrison Subresidency
 July 2007

PARAMETER	Site Background					FIL PC-1 7/20/07	FIL PC-2 7/20/07	FIL PC-3 7/20/07	Trip Blank 7/20/07	NATURAL AMBIENT GROUNDWATER RANGES (n)	NYSDEC CLASS GA STANDARDS (a)
	FIL LMW-2 7/20/07	FIL LMW-4 7/20/07	FIL PC-1 7/20/07	FIL PC-2 7/20/07	FIL PC-3 7/20/07						
TAL Metals (ug/L)											
Aluminum	ND	ND	ND	ND	ND	ND	ND	ND	<5.0 - 1000	NS	
Antimony	ND	ND	ND	ND	ND	ND	ND	ND	NA	3	
Arsenic	ND	ND	ND	ND	ND	ND	ND	ND	<1.0 - 30	25	
Barium	110	86	63	88	85	85	85	85	10 - 500	1000	
Beryllium	ND	ND	ND	ND	ND	ND	ND	ND	<10	3.0 GV	
Cadmium	ND	ND	ND	ND	ND	ND	ND	ND	<1.0	5	
Calcium	77000	51000	49000	69000	41000	41000	41000	41000	1000 - 150000	NS	
Chromium	ND	7.8	ND	ND	ND	ND	ND	ND	<1.0 - 5.0	50	
Cobalt	ND	33	ND	2.5	0.94	0.94	0.94	0.94	<10	NS	
Copper	4.3	ND	2.4	ND	ND	ND	ND	ND	<1.0 - 3	200	
Iron	ND	22000	ND	13000	2800	2800	2800	2800	10 - 10000	300 (m)	
Lead	ND	ND	ND	ND	ND	ND	ND	ND	<15	25	
Magnesium	29000	20000	5800	20000	12000	12000	12000	12000	1000 - 50000	35000 GV	
Manganese	17	25000	820	7700	1800	1800	1800	1800	<1.0 - 1000	300 (m)	
Mercury	ND	ND	ND	ND	ND	ND	ND	ND	<1.0	0.7	
Nickel	3.3	ND	1.3	ND	1.9	1.9	1.9	1.9	<10 - 50	100	
Potassium	3900	2900	4200	6300	3800	3800	3800	3800	1000 - 10000	NS	
Selenium	21	11	12	13	12	12	12	12	<1.0 - 10	10	
Silver	ND	5.6	2.0	2.0	0.61	0.61	0.61	0.61	<5	50	
Sodium	39000	35000	77000	45000	75000	75000	75000	75000	500 - 120000	20000	
Thallium	ND	ND	ND	ND	ND	ND	ND	ND	NA	0.5 GV	
Vanadium	3.9	ND	1.8	2.0	1.7	1.7	1.7	1.7	<1.0 - 10	NS	
Zinc	ND	ND	ND	ND	ND	ND	ND	ND	<10 - 2000	2000 GV	
Chloride (mg/l)	9.1	30	52	20	84	84	84	84	NA	250	
Cyanide (mg/l)	ND	ND	ND	ND	0.012	0.012	0.012	0.012	NA	200	
Volatile Organics (ug/L)											
Toluene	1.2	1.5	3.2	3.5	1.6	1.6	1.6	1.6	NA	5	
Semi-Volatile Organics (ug/L)											
Total SVOCs	ND	ND	ND	ND	ND	ND	ND	ND	NA	50	

(a) - NYSDEC Division of Water Technical and Operational Guidance Series (1.1.1). June 1998, revised April 2000.
 GV - Guidance Value.
 (m) - Sum of Iron and Manganese not to exceed 500 ug/L.
 (n) - Dragun, J., The Soil Chemistry of Hazardous Materials.
 NA - Not applicable.
 ND - Not detected at analytical detection limit.
 J- Detected below the detection limit.
 NS - No standard.
 * - Not analyzed.

TABLE 2
SURFACE WATER DATA SUMMARY
Fifth Quarter Sampling - Harrison Subresidency
July 2007

PARAMETER	Site Background				NATURAL AMBIENT GROUNDWATER RANGES (n)	NYSDEC CLASS GA STANDARDS (a)	NYSDEC CLASS A STANDARDS (a)
	SW-1 7/20/07	SW-2 7/20/07	SW-3 7/20/07	SW-4 7/20/07			
TAL Metals (ug/L)							
Aluminum	47000	93	*	2600	<5.0 - 1000	NS	100 ²
Antimony	ND	ND	*	ND	NA	3	3 ¹
Arsenic	10	ND	*	ND	<1.0 - 30	25	50 ¹ , 150 ² , 340 ³
Barium	360	26	*	94	10 - 500	1000	1,000 ¹
Beryllium	1	ND	*	ND	<10	3.0 GV	3 GV ¹
Cadmium	0.8	ND	*	ND	<1.0	5	5 ¹
Calcium	71000	44000	*	47000	1000 - 150000	NS	NS
Chromium	73	ND	*	6.9	<1.0 - 5.0	50	50 ¹
Cobalt	21	ND	*	2.5	<10	NS	5 ²
Copper	97	22	*	35	<1.0 - 3	200	200 ¹
Iron	78000	450	*	8200	10 - 10000	300 (m)	300 ^{2,4}
Lead	140	ND	*	15	<15	25	50 ¹
Magnesium	26000	13000	*	14000	1000 - 50000	35000 GV	35,000 ¹
Manganese	5400	520	*	2600	<1.0 - 1000	300 (m)	300 ⁴
Mercury	0.27	ND	*	ND	<1.0	0.7	0.7 ¹ , 7e-4 ⁵ , 0.77 ² , 1.4 ³ , 0.0026 ⁶
Nickel	46	ND	*	5.9	<10 - 50	100	100 ¹
Potassium	11000	3500	*	3800	1000 - 10000	NS	NS
Selenium	17	14	*	17	<1.0 - 10	10	10 ¹ , 4.6 ²
Silver	3.4	ND	*	0.78	<5	50	50 ¹
Sodium	45000	11000	*	12000	500 - 120000	20000	NS
Thallium	ND	ND	*	ND	NA	0.5 GV	0.5 GV ^{1,8}
Vanadium	110	2.4	*	9.7	<1.0 - 10	NS	14 ²
Zinc	270	ND	*	38	<10 - 2000	2000 GV	2,000 GV ¹ , 5,000 GV ⁴
Chloride (mg/l)	15	6.1	*	6.4	NA	250	250,000 ¹
Cyanide (mg/l)	ND	0.014	*	ND	NA	200	200 ¹ , 9,000 ⁵ , 5.2 ² , 22 ³
Volatile Organics (ug/L)							
Toluene	16	2.8	*	1.8	NA	5	5 ¹ , 6,000 ⁵ , 100 GV ² , 480 GV ³
Semi-Volatile Organics (ug/L)							
Total SVOCs	ND	ND	*	ND	NA	50	NA

(a) - NYSDEC Division of Water Technical and Operational Guidance Series (1.1.1). June 1998, revised April 2000.

(m) - Sum of Iron and Manganese not to exceed 500 ug/L.

(n) - Dragun, J., The Soil Chemistry of Hazardous Materials.

Class A Standards for Surface Water as a source of Drinking Water¹ as these are tributaries to Kensico Reservoir.

Other Class A Standards are for Fish Propagation² Fish Survival³ Aesthetic⁴ Human Consumption of Fish⁵, and Wildlife Protection⁶

GV - Guidance Value.

ND - Not detected at analytical detection limit.

* - Not analyzed.

NS - No Standard.

NA - Not applicable.

Note - results in bold exceed one or more of the standards.

- select Class A standards are specific to sample conditions and require hardness concentrations; therefore, these were not included.

TABLE 3
SEDIMENT DATA SUMMARY
Fifth Quarter Sampling - Harrison Subresidency
July 2007

PARAMETER	Site Background					Sediment Criteria (a)	
	SD-1 7/20/07	SD-2 7/20/07	SD-3 7/20/07	SD-4 7/20/07	SD-5 Dup. of SD-4 7/20/07	LEL ¹	SEL ²
TAL Metals (mg/kg)							
Aluminum	5700	3700	8800	3600	3700	NA	NA
Antimony	ND	ND	ND	ND	ND	2	<u>25</u>
Arsenic	1.5	2	3	2.7	1.3	6	<u>33</u>
Barium	34	52	54	36	47	NA	NA
Beryllium	0.15	0.042	0.37	0.088	0.057	NA	NA
Cadmium	0.13	0.18	0.55	0.24	0.27	0.6	<u>9</u>
Calcium	8800	42000	23000	9400	11000	NA	NA
Chromium	12	7.4	15	8.2	8.6	26	<u>110</u>
Cobalt	3.5	3.6	5.4	3.5	3.1	NA	NA
Copper	17	11	22	18	8.7	16	<u>110</u>
Iron	9900	11000	13000	12000	10000	20000	<u>40000</u>
Lead	18	11	120	19	16	31	<u>110</u>
Magnesium	5900	25000	12000	6700	7100	NA	NA
Manganese	130	1600	200	590	940	460	<u>1100</u>
Mercury	ND	ND	ND	ND	ND	0.15	<u>1.3</u>
Nickel	7.4	8.1	12	9.1	8.6	16	<u>50</u>
Potassium	710	920	880	560	760	NA	NA
Selenium	ND	ND	ND	ND	ND	NA	NA
Silver	ND	0.16	ND	0.088	0.12	1	<u>2.2</u>
Sodium	78	53	180	51	58	NA	NA
Thallium	ND	ND	ND	ND	ND	NA	NA
Vanadium	17	14	25	11	12	NA	NA
Zinc	35	33	81	56	58	120	<u>270</u>
Chloride	ND	ND	ND	ND	ND	NA	NA
Cyanide	ND	ND	0.47	ND	ND	NA	NA
Volatile Organics (mg/Kg)						Sediment Criteria (a) Water Qual.	
Acetone	0.13	ND	0.13	0.078	0.14	NA	NA
Carbon Disulfide	0.007J	0.0064J	0.002J	0.0032J	0.0052J	NA	NA
Methylene Chloride	0.055B	0.060B	0.041B	0.038B	0.036B	NA	NA
Semi-Volatile Organics (mg/Kg)							
Anthracene	ND	ND	0.072J	ND	ND	NA	NA
Benzo(a)anthracene	ND	ND	0.18J	ND	ND	1.3	1.3
Benzo(a)pyrene	ND	ND	0.12J	ND	ND	1.3	1.3
Benzo(b)fluoranthene	ND	ND	0.14J	ND	ND	1.3	1.3
Benzo(g,h,i)perylene	ND	ND	0.082J	ND	ND	NA	NA
Benzo(k)fluoranthene	ND	ND	0.080J	ND	ND	1.3	1.3
Chrysene	ND	ND	0.16J	ND	ND	1.3	1.3
Di-n-butylphthalate	0.071JB	ND	ND	ND	ND	NA	NA
Fluoranthene	ND	ND	0.34J	ND	0.051J	NA	NA
Indeno(1,2,3-cd)pyrene	ND	ND	0.078J	ND	ND	1.3	1.3
Phenanthrene	ND	ND	0.31J	ND	ND	NA	NA
Pyrene	ND	0.051J	0.43J	ND	0.070J	NA	NA

(a) - NYSDEC Technical Guidance for Screening Contaminated Sediments. November 1993, revised January 1999.

1 - Lowest Effect Level

2 - Severe Effect Level

ND - Not detected at analytical detection limit.

B - Detected in laboratory sample.

J - Detected below detection limit.

NA - No applicable criterion.

Note - results exceeding the LEL and SEL are shown in **bold** and underlined, respectively.

Table 4
Field Duplicate Relative Percent Difference
Fifth Quarter Sampling - Harrison Subresidency
July 2007

Parameters	SD-4	Duplicate SD-5	Relative Percent Difference
Metals (mg/Kg)			
Aluminum	3600	3700	2.74
Antimony	ND	ND	NA
Arsenic	2.7	1.3	70.00
Barium	36	47	26.51
Beryllium	0.088	0.057	42.76
Cadmium	0.24	0.27	11.76
Calcium	9400	11000	15.69
Chromium	8.2	8.6	4.76
Cobalt	3.5	3.1	12.12
Copper	18	8.7	69.66
Iron	12000	10000	18.18
Lead	19	16	17.14
Magnesium	6700	7100	5.80
Manganese	590	940	45.75
Mercury	ND	ND	NA
Nickel	9.1	8.6	5.65
Potassium	560	760	30.30
Selenium	ND	ND	NA
Silver	0.088	0.12	30.77
Sodium	51	58	12.84
Thallium	ND	ND	NA
Vanadium	11	12	8.70
Zinc	56	58	3.51
Chloride	ND	ND	NA
Cyanide	ND	ND	NA
Volatile Organics (mg/Kg)			
Actone	0.078	0.140	56.88
Carbon Disulfide	0.003	0.005	43.90
Methylene Chloride	0.038	0.036	5.41
Semi-Volatile Organics (mg/Kg)			
Fluoranthene	ND	0.051	NA
Pyrene	ND	0.07	NA

Notes:

- 1) ND indicates analyte not detected at analytical reporting limit.
- 2) NA indicates Not Applicable
- 3) Values in bold exceed the normally accepted range of 15%

collected and split between the original and duplicate sample bottles). Results for all other parameters analyzed are within the acceptable range for RPDs. HDR|LMS is confident that the original results for SD-4 are sufficient to assess sediment quality conditions taking into account the inherent variability in analyte concentrations. All analytical values reported can be utilized for the July 2007 Fifth Quarter Sampling period.

4.0 COMPARISON TO APPLICABLE CRITERIA

4.1 Groundwater

TAL Metals and Chloride

The results of the groundwater filtered TAL metal and chloride analyses were compared to current NYSDEC Ambient Water Quality Class GA Standards and Guidance Values (April 2000 Revision).

The TAL metals groundwater analysis revealed detectable concentrations of three non-RCRA metals (iron, manganese and sodium) and one RCRA metal (selenium) that exceeded their respective groundwater criteria (Table 1). LMW-2, the upgradient/site background well; however, had a comparable concentration of sodium and a higher concentration of selenium, which would indicate a contribution from an upgradient/off-site source and not necessarily a release from the landfill. Results for the remaining metals analyzed were either non-detect at the respective analytical reporting limits, or less than the respective Class GA standards or guidance values.

Chloride was not detected above the standard in the groundwater.

VOCs, SVOCs and Cyanide

The results of the VOCs, SVOCs and cyanide groundwater analyses (Table 1) were compared to current NYSDEC Ambient Water Quality Class GA Standards and Guidance Values (April 2000 Revision). There were no detections above the Class GA standards in the groundwater samples.

4.2 Surface Water

Surface water results were compared to both NYSDEC Ambient Water Quality Class GA Groundwater Standards and Guidance Values and NYSDEC Class A Surface Water Standards and Guidance Values, where available, as these water bodies are tributaries to the Kensico Reservoir, which is a source of drinking water. Class A standards were available for surface water as a source of drinking water H(WS), human consumption of fish H(FC), fish propagation A(C), fish survival A(A), protection of wildlife (W), and aesthetics (E). Select standards rely on sample specific conditions (i.e., hardness

concentrations) and therefore were not included. Where appropriate, the most conservative standard or guidance value was used for comparison purposes.

TAL Metals and Chloride

The results of the TAL metals analysis revealed detectable concentrations of three non-RCRA metals (aluminum, iron, and manganese) and one RCRA metal (selenium) in the unfiltered downgradient/on-site surface water samples that exceeded their respective criteria (Table 2). Comparable, if not greater, concentrations of these same metals were detected in the upgradient/site background sample collected from SW-1, which would indicate a contribution from an upgradient/off-site source. Further, elevated levels (above either Class GA or Class standards or guidance values) of the metals chromium, cobalt, lead, mercury, and vanadium were also detected in SW-1, again indicating a contribution from an upgradient/off-site source. Results for the remaining metals analyzed were either non-detect at the respective analytical reporting limits, or less than the respective Class GA and/or Class A standards or guidance values.

Chloride was not detected above the Class GA standard in the surface water samples.

VOCs, SVOCs, and Cyanide

The VOC, toluene, was detected above the Class GA and Class A standard of 5 ug/l, respectively, in the upgradient/site background surface water sample, SW-1, at a concentration of 16 ug/l (Table 2). Toluene was also detected in the trip blank sample (Table 1) at a concentration of 10 ug/l. Because toluene was also detected in the trip blank sample, results are likely biased high. HDR|LMS is uncertain of the source of this contamination.

There were no detections above the Class GA standards of SVOCs or cyanide in the surface water samples (Table 2).

4.3 Sediment

The results of the sediment analysis (Table 3) were compared to the NYSDEC Technical Guidance for Screening Contaminated Sediments (January 1999 Revision).

TAL Metals and Chloride

The results of the TAL metals analysis indicate that the concentration of the RCRA metal lead exceeded the Severe Effect Level (SEL) in the sediment sample collected from SD-3. Concentrations of the non-RCRA metal copper exceeded the Lowest Effect Level (LEL) in SD-1, SD-3 and SD-4 and the non-RCRA metal manganese exceeded the LEL in SD-4 and SEL in SD-2. The concentration of copper detected in the upgradient/site background sediment location, SD-1, would indicate a contribution from an upgradient/off-site source.

According to the NYSDEC Technical Guidance for Screening Contaminated Sediments, a sediment is considered contaminated if either criterion is exceeded. If both criterion are exceeded, the sediment is considered to be severely impacted. If only the LEL criterion is exceeded, the impact is considered moderate.

Chloride was not detected in the sediment samples.

VOCs, SVOCs and Cyanide

The results indicate that there were no detections of VOCs, SVOCs or cyanide above the sediment criteria for the protection of human health.

5.0 CONCLUSIONS

Groundwater analytical results revealed elevated levels of the non-RCRA metals iron, manganese, and sodium and the RCRA metal selenium in the filtered results obtained at one or more of the downgradient well locations. An elevated sodium concentration was detected at the upgradient/site background well, LMW-2. The sodium concentrations detected at two of the downgradient wells (PC-1 and PC-3), however, were much higher. This would indicate a contribution from the landfill and perhaps some additional contribution from an upgradient/off-site source.

In comparison with concentrations detected in the downgradient wells, a higher selenium concentration was detected at the upgradient/site background well, LMW-2. This would indicate a contribution from an upgradient/off-site source and not necessarily a release from the landfill.

Concentrations of iron and manganese in the downgradient/on-site wells are greater than the concentrations detected in the upgradient/site background well, LMW-2, which would indicate some contribution from the landfill. Iron and manganese concentrations decrease in the downgradient/off-site monitoring well (PC-3) demonstrating that these metals are attenuating as they are migrating off-site. The remaining groundwater sample results were either non-detect or less than the respective NYSDEC Class GA standards or guidance values.

Surface water analytical results revealed elevated levels of the non-RCRA metals aluminum, iron, and manganese, and the RCRA metal selenium in the results obtained at one or more of the downgradient/on-site surface water sample locations. Comparable, if not greater, concentrations of these metals were detected in the sample results obtained from the upgradient/site background surface water location, SW-1. This would indicate a contribution from an upgradient/off-site source and not necessarily a release from the landfill. Elevated levels of several other metals including chromium, cobalt, lead, mercury, and vanadium and the VOC toluene were also detected in the sample collected from SW-1; again indicating a contribution from an upgradient/off-site source. The remaining surface water sample results were either non-detect or less than the respective NYSDEC Class GA and/or Class A standards or guidance values.

Sediment analytical results revealed elevated levels of the non-RCRA metals, copper and manganese, and the RCRA metal lead in one or more of the downgradient/on-site

sediment sample locations. A comparable copper concentration was detected in the upgradient/site background sediment sample location, SD-1. This would indicate a contribution from an upgradient/off-site source and not necessarily a release from the landfill. Concentrations of manganese in the downgradient/on-site sediment sample locations are greater than the concentration detected in the upgradient/site background sediment location, which would indicate some contribution from the landfill. Although the concentration of lead in the sample collected from the downgradient/on-site sediment sample location, SD-3, is elevated, SD-4 (downgradient from SD-3) is much lower indicating that lead is not significantly migrating off-site. The remaining sediment sample results were either non-detect or less than the respective NYSDEC sediment criteria.

ATTACHMENT A

GROUNDWATER MONITORING WELL LOGS

Date: 7/20/2007	METERS USED
Crew: BG/SQ	Temp.: YSI-556-01
Job No.: 137-063423-001	pH: YSI-556-01
Project: Harrison Subresidency Landfill	Cond.: YSI-556-01
Project Site: Harrison, New York	Turb.: Micro TPW

Well ID No.:	LMW-2	DTW Before Sampling:	12.51
Well Condition:	Good	Sample Date/Time:	7/20/07 at 1300
Well Depth/Diameter:	24.19'2"	Sampling Method:	Dedicated Bailer
Well Casing Type:	PVC	Sampling Depth(s):	
Screened Interval:	NA	DTW After Sampling:	
Casing Ht./Lock No.:		Chain-of-Custody No.(s):	
Reference Pt.:	TOC	Analytical Lab(s):	Veritech
Depth to Water (DTW):	11.98	Sampling Observations:	
Water Column Ht./Vol.:	12.21	Water clear	
Purge Est.:	11.3		
Purge Method(s):	Dedicated Bailer		
Purge Date/Time(s):			

SAMPLE CHEMISTRIES

	Temp. (°C)	pH	Sp. Cond	Turb.
Start	13.97	8.01	0.75	16.82
End				

Depth(s):
Rates (gpm):
Purged Volume: 3.8 gallons
DTW After Purging:
Yield Rate: L - M - H
Purge Observations:
Well dry after purging 3.8 gallons.
Water very turbid.

SAMPLE ANALYSES

Parameters	Inv. No.	Pres. Meth.	Filter
VOCs		Ice	N
SVOCs		Ice	N
Cn		Ice	N
Cl		Ice	N
TAL Metals (FILTERED ONLY*)			Y*

PURGE CHEMISTRIES

Vol.	Temp. (°C)	pH	Sp. Cond.	Turb.
0	12.41	7.16	0.787	99.5
3.8	12.38	7.78	0.77	747

* To be filtered and preserved by laboratory.

Air Temp: 70°
Weather Conditions: Fair

Crew Chief Signature _____

Date: _____

HDR|LMS

Well Sampling Log

Date: 7/20/2007
Crew: BG/SQ
Job No: 137-063423-001
Project: Harrison Subresidency Landfill
Project Site: Harrison, New York

Well ID No.: MW-4
Well Condition: Good
Well Depth/Diameter: 15.73'2"
Well Casing Type: PVC
Screened Interval: NA
Casing Ht./Lock No.:
Reference Pt.: TOC
Depth to Water (DTW): 5.2
Water Column Ht./Vol.: 10.53
Purge Est.: 9.8
Purge Method(s): Dedicated Bailer
Purge Date/Time(s):

Depth(s):
Rates (gpm):
Purged Volume: 9.8 gallons
DTW After Purging:
Yield Rate: L - M - H
Purge Observations:
Water slightly turbid then clear.

PURGE CHEMISTRIES

Vol.	Temp. (°C)	pH	Sp. Cond.	Turb.
0	16.2	6.91	1.108	25.6
3.3	15.56	6.97	1.126	49.38
6.6	16.24	7.04	1.115	23.73
9.8	16.91	7.13	1.079	12.28

METERS USED

Temp.: YSI-556-01
pH: YSI-556-01
Cond.: YSI-556-01
Turb.: Micro TPW

DTW Before Sampling: 5.7
Sample Date/Time: 7/20/07 at 1330
Sampling Method: Dedicated Bailer
Sampling Depth(s):
DTW After Sampling:
Chain-of-Custody No.(s):
Analytical Lab(s): Veritech
Sampling Observations: Water rust-colored and turbid.

SAMPLE CHEMISTRIES

	Temp. (°C)	pH	Sp. Cond.	Turb.
Start	19.91	7.92	1.081	74.9
End				

SAMPLE ANALYSES

Parameters	Inv. No.	Pres. Meth.	Filter
VOCs		Ice	N
SVOCs		Ice	N
Cn		Ice	N
Cl		Ice	N
TAL Metals (FILTERED ONLY*)			Y*

* To be filtered and preserved by laboratory.

Air Temp: 70°
Weather Conditions: Fair

Crew Chief Signature _____

Date: _____

HDR|LMS

Well Sampling Log

Date: 7/20/2007
Crew: BG/SQ
Job No: 137-063423-001
Project: Harrison Subresidency Landfill
Project Site: Harrison, New York

Well ID No.: PC-1
Well Condition: Good
Well Depth/Diameter: 16.82'2"
Well Casing Type: PVC
Screened Interval: NA
Casing Ht./Lock No.:
Reference Pt.: TOC
Depth to Water (DTW): 7.15
Water Column Ht./Vol.: 9.67
Purge Est.: 9
Purge Method(s): Peristaltic Pump
Purge Date/Time(s):

Depth(s):
Rates (gpm):
Purged Volume: 9 gallons
DTW After Purging:
Yield Rate: L - M - H
Purge Observations:
Water turbid then clear.

PURGE CHEMISTRIES

Vol.	Temp. (°C)	pH	Sp. Cond.	Turb.
0	18.64	7.17	0.808	104.5
3	18.5	7.41	0.71	7.53
6	18.4	7.37	0.681	6.76
9	18.8	7.37	0.677	6.54

METERS USED

Temp.: YSI-556-01
pH: YSI-556-01
Cond.: YSI-556-01
Turb.: Micro TPW

DTW Before Sampling: 7.18
Sample Date/Time: 7/20/07 at 1410
Sampling Method: Peristaltic Pump
Sampling Depth(s):
DTW After Sampling:
Chain-of-Custody No.(s):
Analytical Lab(s): Veritech
Sampling Observations: Water is clear.

SAMPLE CHEMISTRIES

	Temp. (°C)	pH	Sp. Cond.	Turb.
Start	19.36	7.75	0.688	3.26
End				

SAMPLE ANALYSES

Parameters	Inv. No.	Pres. Meth.	Filter
VOCs		Ice	N
SVOCs		Ice	N
Cn		Ice	N
Cl		Ice	Y*
TAL Metals (FILTERED ONLY*)			

* To be filtered and preserved by laboratory.

Air Temp: 70°
Weather Conditions: Fair

Crew Chief Signature _____

Date: _____

HDR|LMS

Well Sampling Log

Date: 7/20/2007
Crew: BG/SQ
Job No: 137-063423-001
Project: Harrison Subresidency Landfill
Project Site: Harrison, New York

Well ID No.: PC-2
Well Condition: Poor (bent)
Well Depth/Diameter: 11.49/2"
Well Casing Type: PVC
Screened Interval: NA
Casing Ht./Lock No.:
Reference Pt.: TOC
Depth to Water (DTW): 3.73
Water Column Ht./Vol.: 7.76
Purge Est.: 7.2
Purge Method(s): Dedicated Bailer
Purge Date/Time(s):

Depth(s):
Rates (gpm):
Purged Volume: 7.2 gallons
DTW After Purging:
Yield Rate: L - M - H
Purge Observations:
Water very turbid.

PURGE CHEMISTRIES

Vol.	Temp. (°C)	pH	Sp. Cond.	Turb.
0	19.88	7.29	0.922	93.67
2.4	16.23	7.12	0.936	205.6
4.8	17.23	7.03	0.914	231.9
7.2	15.54	6.97	0.938	288.7

METERS USED

Temp.: YSI-556-01
pH: YSI-556-01
Cond.: YSI-556-01
Turb.: Micro TPW

DTW Before Sampling: 3.67
Sample Date/Time: 7/20/07 at 1340
Sampling Method: Dedicated Bailer
Sampling Depth(s):
DTW After Sampling:
Chain-of-Custody No.(s):
Analytical Lab(s): Veritech
Sampling Observations: Water very turbid.

SAMPLE CHEMISTRIES

	Temp. (°C)	pH	Sp. Cond.	Turb.
Start	17.66	7.86	0.96	171.8
End				

SAMPLE ANALYSES

Parameters	Inv. No.	Pres. Meth.	Filter
VOCs		Ice	N
SVOCs		Ice	N
Cn		Ice	N
Cl		Ice	N
TAL Metals (FILTERED ONLY*)			Y*

* To be filtered and preserved by laboratory.

Air Temp: 70°
Weather Conditions: Fair

Crew Chief Signature _____

Date: _____

Date: 7/20/2007
Crew: BG/SQ
Job No: 137-063423-001
Project: Harrison Subresidency Landfill
Project Site: Harrison, New York

METERS USED

Temp.: YSI-556-01
pH: YSI-556-01
Cond.: YSI-556-01
Turb.: Micro TPW

Well ID No.: PC-3
Well Condition: Good
Well Depth/Diameter: 18.57'2"
Well Casing Type: PVC
Screened Interval: NA
Casing Ht./Lock No.:
Reference Pt.: TOC
Depth to Water (DTW): 9
Water Column Ht./Vol.: 9.57
Purge Est.: 9
Purge Method(s): Peristaltic Pump
Purge Date/Time(s):

DTW Before Sampling: 8.98
Sample Date/Time: 7/20/07 at 1550
Sampling Method: Peristaltic Pump
Sampling Depth(s):
DTW After Sampling:
Chain-of-Custody No.(s):
Analytical Lab(s): Veritech
Sampling Observations: Water clear

SAMPLE CHEMISTRIES

	Temp. (°C)	pH	Sp. Cond.	Turb.
Start	14.71	7.41	0.706	1.72
End				

SAMPLE ANALYSES

Parameters	Inv. No.	Pres. Meth.	Filter
VOCs		Ice	N
SVOCs		Ice	N
Cn		Ice	N
Cl		Ice	N
TAL Metals (FILTERED ONLY**)			Y*

Depth(s):
Rates (gpm):
Purged Volume: 9 gallons
DTW After Purging:
Yield Rate: L - M - H
Purge Observations:
 Water turbid during purge of second well volume then clear.

PURGE CHEMISTRIES

Vol.	Temp. (°C)	pH	Sp. Cond.	Turb.
0	17.36	7.89	1.466	10.88
3	15.6	7.53	0.675	207.2
6	15.03	7.45	0.701	14.55
9	14.97	7.45	0.716	5.69

* To be filtered and preserved by laboratory.

Air Temp: 70°
Weather Conditions: Fair

Crew Chief Signature _____

Date: _____

ATTACHMENT B

SURFACE WATER AND SEDIMENT LOGS

Date: 7/20/2007

Crew: BG/SQ

Site: Harrison Subresidency Landfill

FIELD DATA SHEET FOR SURFACE WATER

Meter(s) Used:

YSI - 556-01/Micro TPW

Sta. No.	Time	Sample Depth	Total Depth	Temp. °C	pH	Cond. umhos/cm	Turb. NTUs	Flow CFS	Sample Parameters	Comments
SW-1	1250	1"	1-2"	26.5	7.21	0.567	75.6	<1	TAL Metals (U), Cl, Cn, VOCs, SVOCS	Water Turbid
SW-2	1130	2"	2-3"	21.9	7.39	0.391	16.3	<1	TAL Metals (U), Cl, Cn, VOCs, SVOCS	Water Clear
SW-3	-	-	-	-	-	-	-	-	N/A	Stream Dry. No sample collected.
SW-4	1445	2"	2-3"	24.1	7.7	0.399	24.6	<1	TAL Metals (U), Cl, Cn, VOCs, SVOCS	Water Slightly Turbid

Weather: 70° Fair

FIELD DATA SHEET FOR SEDIMENT

Date: 7/20/2007
 Crew: BG/SQ
 Site: Harrison Subresidency Landfill

Sta. No.	Time	Sample Depth	Method	Texture	Color	Odor/ Staining	Sample Parameters	Comments
SD-1	1250	4-6"	Dedicated Spoon	Sandy/Silty	Drk to Med. Br.	None	TAL Metals, Cn, Cl, VOCs, SVOCs	
SD-2	1130	0-2"	Dedicated Spoon	Sandy/Silty	Drk to Med. Br.	None	TAL Metals, Cn, Cl, VOCs, SVOCs	
SD-3	1210	4-6"	Dedicated Spoon	Sandy/Silty	Drk to Med. Br.	None	TAL Metals, Cn, Cl, VOCs, SVOCs	Sream Dry
SD-4	1450	4-6"	Dedicated Spoon	Sandy/Silty	Drk to Med. Br.	None	TAL Metals, Cn, Cl, VOCs, SVOCs	
SD-5	1455	4-6"	Dedicated Spoon	Sandy/Silty	Drk to Med. Br.	None	TAL Metals, Cn, Cl, VOCs, SVOCs	Duplicate of SD-4

Weather: 70° Fair

ATTACHMENT C
GAS MONITORING LOGS

Date: 7/20/2007

Crew: BG/SQ

Site: Harrison Subpresidency Landfill

CGI: S/N 17045

PID: S/N 4539

FID: Foxboro OVA

AIR MONITORING FIELD DATA SHEET

Sample Point	Time	Inspector	% LEL CGI	PID Equiv.	FID PPM	CH4	Background		Observations/Notes
							PID	FID	
GV-1	1018	SQ	0	0	0		0		
GV-2	1020	SQ	0	0	0		0		
GV-3	1021	SQ	0	0	0		0		
GV-4	1022	SQ	0	0	0		0		
NP*	1024	SQ	0	0	0		0		
EP	1025	SQ	0	0	0		0		
SP	1026	SQ	0	0	0		0		
WP	1028	SQ	0	0	0		0		

*Note - "NP" indicates North Perimeter

Weather: 70°

Fair

ATTACHMENT D

LABORATORY ANALYTICAL DATA

Hampton-Clarke, Inc.

veritech laboratories



175 Route 46 West, Unit D
Fairfield, NJ 07004
(973) 244-9770
Federal ID: 222679402

Format: NYDOH-CatA

Project: Harrison Landfill

PO Number:

Client: HDR/LMS
One Blue Hill Plaza
P.O. Box 1509
Pearl River, NY 10965

Attn: Terry Schneider

Samples submitted on: 7/21/2007

AC31901-001
AC31901-002
AC31901-003
AC31901-004
AC31901-005
AC31901-006
AC31901-007
AC31901-008
AC31901-009
AC31901-010
AC31901-011
AC31901-012
AC31901-013
AC31901-014
AC31901-015
AC31901-016
AC31901-017
AC31901-018
AC31901-019

Date: 8/14/2007

HCI Project: 7072310

This report is a true report of results obtained from our tests of this material. In lieu of a formal contract document, the total aggregate liability of Veritech to all parties shall not exceed Veritech's total fee for analytical services rendered.

Jeri Rossi - Quality Assurance Director

Or

Stanley Gilewicz - Laboratory Director

CT #: PH-0671

MA #: NJ386

NJ #: 14622

NY #: 11408

PA #: 68-463

WV #: 353

USACE

0000

Form1 Inorganic Analysis Data Sheet

Sample ID: AC31901-007
 Client Id: LMW-2 F
 Matrix: AQUEOUS
 Level: LOW

% Solid: 0
 Units: UG/L
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-36-0	Antimony	3.9	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-38-2	Arsenic	4.9	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-39-3	Barium	0.45	110		108/02/07	8159	W8159C2m23		P	PEICP2
7440-41-7	Beryllium	0.24	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-43-9	Cadmium	0.78	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-70-2	Calcium	20	77000		108/02/07	8159	W8159C2m23		P	PEICP2
7440-47-3	Chromium	5.8	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-48-4	Cobalt	0.54	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-50-8	Copper	1.5	4.3		108/02/07	8159	W8159C2m23		P	PEICP2
7439-89-6	Iron	75	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7439-92-1	Lead	1.7	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7439-95-4	Magnesium	63	29000		108/02/07	8159	W8159C2m23		P	PEICP2
7439-96-5	Manganese	4.7	17		108/02/07	8159	W8159C2m23		P	PEICP2
7439-97-6	Mercury	0.21	ND		108/02/07	8159	8159SWMD19		CV	HGCV1
7440-02-0	Nickel	1.1	3.3		108/02/07	8159	W8159C2m23		P	PEICP2
7440-09-7	Potassium	52	3900		108/02/07	8159	W8159B2M123		P	PEICPRAD2
7782-49-2	Selenium	10	21		108/02/07	8159	W8159C2m23		P	PEICP2
7440-22-4	Silver	0.53	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-23-5	Sodium	130	39000		108/02/07	8159	W8159B2M123		P	PEICPRAD2
7440-28-0	Thallium	4.8	ND		108/02/07	8159	W8159C2m23		P	PEICP2
7440-62-2	Vanadium	1.7	3.9		108/02/07	8159	W8159C2m23		P	PEICP2
7440-66-6	Zinc	1.5	ND		108/02/07	8159	W8159C2m23		P	PEICP2

Comments: _____

Flag Codes:

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

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-006
Client Id: LMW-2 U
Data File: 2M22213.D
Analysis Date: 07/30/07 19:56
Date Rec/Extracted: 07/21/07-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	56-23-5	Carbon Tetrachloride	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	108-90-7	Chlorobenzene	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	67-66-3	Chloroform	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	74-87-3	Chloromethane	5.0	U
156-59-2	cis-1,2-Dichloroethene	5.0	U	107-06-2	1,2-Dichloroethane	5.0	U
10061-01-5	cis-1,3-Dichloropropene	5.0	U	78-87-5	1,2-Dichloropropane	5.0	U
124-48-1	Dibromochloromethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
1330-20-7	m&p-Xylenes	2.0	U	78-93-3	2-Butanone	5.0	U
75-09-2	Methylene Chloride	5.0	U	591-78-6	2-Hexanone	5.0	U
95-47-6	o-Xylene	1.0	U	108-10-1	4-Methyl-2-Pentanone	5.0	U
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	1.2	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
75-01-4	Vinyl Chloride	5.0	U				

Worksheet #: 54676

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 SEMIVOLATILE REPORT

Sample Number: AC31901-006

Client Id: LMW-2 U

Data File: 5M33231.D

Analysis Date: 07/27/07 20:22

Date Rec/Extracted: 07/21/07-07/25/07

Matrix: Aqueous

Initial Vol: 900ml

Final Vol: 1ml

Dilution: 1

Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	11	U	205-99-2	Benzo[b]fluoranthene	11	U
95-50-1	1,2-Dichlorobenzene	11	U	191-24-2	Benzo[g,h,i]perylene	11	U
541-73-1	1,3-Dichlorobenzene	11	U	207-08-9	Benzo[k]fluoranthene	11	U
106-46-7	1,4-Dichlorobenzene	11	U	111-91-1	bis(2-Chloroethoxy)methan	11	U
95-95-4	2,4,5-Trichlorophenol	11	U	111-44-4	bis(2-Chloroethyl)ether	11	U
88-06-2	2,4,6-Trichlorophenol	11	U	39638-32-9	bis(2-chloroisopropyl)ether	11	U
120-83-2	2,4-Dichlorophenol	11	U	117-81-7	bis(2-Ethylhexyl)phthalate	11	U
105-67-9	2,4-Dimethylphenol	11	U	85-68-7	Butylbenzylphthalate	11	U
51-28-5	2,4-Dinitrophenol	28	U	86-74-8	Carbazole	11	U
121-14-2	2,4-Dinitrotoluene	11	U	218-01-9	Chrysene	11	U
606-20-2	2,6-Dinitrotoluene	11	U	53-70-3	Dibenzo[a,h]anthracene	11	U
91-58-7	2-Chloronaphthalene	11	U	132-64-9	Dibenzofuran	11	U
95-57-8	2-Chlorophenol	11	U	84-66-2	Diethylphthalate	11	U
91-57-6	2-Methylnaphthalene	11	U	131-11-3	Dimethylphthalate	11	U
95-48-7	2-Methylphenol	11	U	84-74-2	Di-n-butylphthalate	11	U
88-74-4	2-Nitroaniline	11	U	117-84-0	Di-n-octylphthalate	11	U
88-75-5	2-Nitrophenol	11	U	206-44-0	Fluoranthene	11	U
106-44-5	3&4-Methylphenol	11	U	86-73-7	Fluorene	11	U
91-94-1	3,3'-Dichlorobenzidine	11	U	118-74-1	Hexachlorobenzene	11	U
99-09-2	3-Nitroaniline	11	U	87-68-3	Hexachlorobutadiene	11	U
534-52-1	4,6-Dinitro-2-methylphenol	28	U	77-47-4	Hexachlorocyclopentadiene	11	U
101-55-3	4-Bromophenyl-phenylether	11	U	67-72-1	Hexachloroethane	11	U
59-50-7	4-Chloro-3-methylphenol	11	U	193-39-5	Indeno[1,2,3-cd]pyrene	11	U
106-47-8	4-Chloroaniline	11	U	78-59-1	Isophorone	11	U
7005-72-3	4-Chlorophenyl-phenylether	11	U	91-20-3	Naphthalene	11	U
100-01-6	4-Nitroaniline	11	U	98-95-3	Nitrobenzene	11	U
100-02-7	4-Nitrophenol	28	U	621-64-7	N-Nitroso-di-n-propylamine	11	U
83-32-9	Acenaphthene	11	U	86-30-6	n-Nitrosodiphenylamine	11	U
208-96-8	Acenaphthylene	11	U	87-86-5	Pentachlorophenol	28	U
120-12-7	Anthracene	11	U	85-01-8	Phenanthrene	11	U
56-55-3	Benzo[a]anthracene	11	U	108-95-2	Phenol	11	U
50-32-8	Benzo[a]pyrene	11	U	129-00-0	Pyrene	11	U

Worksheet #: 54667

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: AC31901-009
 Client Id: MW-4 F
 Matrix: AQUEOUS
 Level: LOW

% Solid: 0
 Units: UG/L
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7440-36-0	Antimony	3.9	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7440-38-2	Arsenic	4.9	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7440-39-3	Barium	0.45	86		108/02/07	8159	W8159C2m24		P	PEICP2
7440-41-7	Beryllium	0.24	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7440-43-9	Cadmium	0.78	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7440-70-2	Calcium	20	51000		108/02/07	8159	W8159C2m24		P	PEICP2
7440-47-3	Chromium	5.8	7.8		108/02/07	8159	W8159C2m24		P	PEICP2
7440-48-4	Cobalt	0.54	33		108/02/07	8159	W8159C2m24		P	PEICP2
7440-50-8	Copper	1.5	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7439-89-6	Iron	75	22000		108/02/07	8159	W8159C2m24		P	PEICP2
7439-92-1	Lead	1.7	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7439-95-4	Magnesium	63	20000		108/02/07	8159	W8159C2m24		P	PEICP2
7439-96-5	Manganese	4.7	25000		108/02/07	8159	W8159C2m24		P	PEICP2
7439-97-6	Mercury	0.21	ND		108/02/07	8159	8159SWMD22		CV	HGCV1
7440-02-0	Nickel	1.1	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7440-09-7	Potassium	52	2900		108/02/07	8159	W8159B2MI24		P	PEICPRAD2
7782-49-2	Selenium	10	11		108/02/07	8159	W8159C2m24		P	PEICP2
7440-22-4	Silver	0.53	5.6		108/02/07	8159	W8159C2m24		P	PEICP2
7440-23-5	Sodium	130	35000		108/02/07	8159	W8159B2MI24		P	PEICPRAD2
7440-28-0	Thallium	4.8	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7440-62-2	Vanadium	1.7	ND		108/02/07	8159	W8159C2m24		P	PEICP2
7440-66-6	Zinc	1.5	ND		108/02/07	8159	W8159C2m24		P	PEICP2

Comments: _____

Flag Codes:

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

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Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-008
Client Id: MW-4 U
Data File: 2M22214.D
Analysis Date: 07/30/07 20:21
Date Rec/Extracted: 07/21/07-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	56-23-5	Carbon Tetrachloride	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	108-90-7	Chlorobenzene	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	67-66-3	Chloroform	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	74-87-3	Chloromethane	5.0	U
156-59-2	cis-1,2-Dichloroethene	5.0	U	107-06-2	1,2-Dichloroethane	5.0	U
10061-01-5	cis-1,3-Dichloropropene	5.0	U	78-87-5	1,2-Dichloropropane	5.0	U
124-48-1	Dibromochloromethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
1330-20-7	m&p-Xylenes	2.0	U	78-93-3	2-Butanone	5.0	U
75-09-2	Methylene Chloride	5.0	U	591-78-6	2-Hexanone	5.0	U
95-47-6	o-Xylene	1.0	U	108-10-1	4-Methyl-2-Pentanone	5.0	U
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	1.5	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
75-01-4	Vinyl Chloride	5.0	U				

Worksheet #: 54676

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 Inorganic Analysis Data Sheet

Sample ID: AC31901-013
 Client Id: PC-1 F
 Matrix: AQUEOUS
 Level: LOW

% Solid: 0
 Units: UG/L
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-36-0	Antimony	3.9	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-38-2	Arsenic	4.9	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-39-3	Barium	0.45	63		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-41-7	Beryllium	0.24	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-43-9	Cadmium	0.78	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-70-2	Calcium	20	49000		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-47-3	Chromium	5.8	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-48-4	Cobalt	0.54	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-50-8	Copper	1.5	2.4		108/02/07	8159	W8159C2m30	30	P	PEICP2
7439-89-6	Iron	75	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7439-92-1	Lead	1.7	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7439-95-4	Magnesium	63	5800		108/02/07	8159	W8159C2m30	30	P	PEICP2
7439-96-5	Manganese	4.7	820		108/02/07	8159	W8159C2m30	30	P	PEICP2
7439-97-6	Mercury	0.21	ND		108/02/07	8159	8159SWMD24		CV	HGCV1
7440-02-0	Nickel	1.1	1.3		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-09-7	Potassium	52	4200		108/02/07	8159	W8159B2MI30		P	PEICPRAD2
7782-49-2	Selenium	10	12		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-22-4	Silver	0.53	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-23-5	Sodium	130	77000		108/02/07	8159	W8159B2MI30		P	PEICPRAD2
7440-28-0	Thallium	4.8	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-62-2	Vanadium	1.7	1.8		108/02/07	8159	W8159C2m30	30	P	PEICP2
7440-66-6	Zinc	1.5	ND		108/02/07	8159	W8159C2m30	30	P	PEICP2

Comments: _____

Flag Codes:

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

7280

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-012
Client Id: PC-1 U
Data File: 2M22216.D
Analysis Date: 07/30/07 21:11
Date Rec/Extracted: 07/21/07-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	56-23-5	Carbon Tetrachloride	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	108-90-7	Chlorobenzene	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	67-66-3	Chloroform	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	74-87-3	Chloromethane	5.0	U
156-59-2	cis-1,2-Dichloroethene	5.0	U	107-06-2	1,2-Dichloroethane	5.0	U
10061-01-5	cis-1,3-Dichloropropene	5.0	U	78-87-5	1,2-Dichloropropane	5.0	U
124-48-1	Dibromochloromethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
1330-20-7	m&p-Xylenes	2.0	U	78-93-3	2-Butanone	5.0	U
75-09-2	Methylene Chloride	5.0	U	591-78-6	2-Hexanone	5.0	U
95-47-6	o-Xylene	1.0	U	108-10-1	4-Methyl-2-Pentanone	5.0	U
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	3.2	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
75-01-4	Vinyl Chloride	5.0	U				

Worksheet #: 54676

Total Target Concentration 3.2

*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 SEMIVOLATILE REPORT

Sample Number: AC31901-012
 Client Id: PC-1 U
 Data File: 5M33229.D
 Analysis Date: 07/27/07 19:37
 Date Rec/Extracted: 07/21/07-07/26/07

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

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	10	U	205-99-2	Benzo[b]fluoranthene	10	U
95-50-1	1,2-Dichlorobenzene	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
541-73-1	1,3-Dichlorobenzene	10	U	207-08-9	Benzo[k]fluoranthene	10	U
106-46-7	1,4-Dichlorobenzene	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
120-83-2	2,4-Dichlorophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
105-67-9	2,4-Dimethylphenol	10	U	85-68-7	Butylbenzylphthalate	10	U
51-28-5	2,4-Dinitrophenol	25	U	86-74-8	Carbazole	10	U
121-14-2	2,4-Dinitrotoluene	10	U	218-01-9	Chrysene	10	U
606-20-2	2,6-Dinitrotoluene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
91-58-7	2-Chloronaphthalene	10	U	132-64-9	Dibenzofuran	10	U
95-57-8	2-Chlorophenol	10	U	84-66-2	Diethylphthalate	10	U
91-57-6	2-Methylnaphthalene	10	U	131-11-3	Dimethylphthalate	10	U
95-48-7	2-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
88-74-4	2-Nitroaniline	10	U	117-84-0	Di-n-octylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	206-44-0	Fluoranthene	10	U
106-44-5	3&4-Methylphenol	10	U	86-73-7	Fluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	118-74-1	Hexachlorobenzene	10	U
99-09-2	3-Nitroaniline	10	U	87-68-3	Hexachlorobutadiene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	25	U	77-47-4	Hexachlorocyclopentadiene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
106-47-8	4-Chloroaniline	10	U	78-59-1	Isophorone	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	91-20-3	Naphthalene	10	U
100-01-6	4-Nitroaniline	10	U	98-95-3	Nitrobenzene	10	U
100-02-7	4-Nitrophenol	25	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
83-32-9	Acenaphthene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
208-96-8	Acenaphthylene	10	U	87-86-5	Pentachlorophenol	25	U
120-12-7	Anthracene	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 #: 54667

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.

01
01
01

Form1 Inorganic Analysis Data Sheet

Sample ID: AC31901-011
Client Id: PC-2 F
Matrix: AQUEOUS
Level: LOW

% Solid: 0
Units: UG/L
Date Rec: 7/23/2007

Lab Name: Veritech
Lab Code:
Contract:

Nras No:
Sdg No:
Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-36-0	Antimony	3.9	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-38-2	Arsenic	4.9	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-39-3	Barium	0.45	88	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-41-7	Beryllium	0.24	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-43-9	Cadmium	0.78	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-70-2	Calcium	20	69000	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-47-3	Chromium	5.8	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-48-4	Cobalt	0.54	2.5	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-50-8	Copper	1.5	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7439-89-6	Iron	75	13000	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7439-92-1	Lead	1.7	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7439-95-4	Magnesium	63	20000	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7439-96-5	Manganese	4.7	7700	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7439-97-6	Mercury	0.21	ND	1	08/02/07	8159	8159SWMD23	23	CV	HGCV1
7440-02-0	Nickel	1.1	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-09-7	Potassium	52	6300	1	08/02/07	8159	W8159B2M125	25	P	PEICPRAD2
7782-49-2	Selenium	10	13	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-22-4	Silver	0.53	2.0	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-23-5	Sodium	130	45000	1	08/02/07	8159	W8159B2M125	25	P	PEICPRAD2
7440-28-0	Thallium	4.8	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-62-2	Vanadium	1.7	2.0	1	08/02/07	8159	W8159C2m25	25	P	PEICP2
7440-66-6	Zinc	1.5	ND	1	08/02/07	8159	W8159C2m25	25	P	PEICP2

Comments: _____

Flag Codes:

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

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-010
Client Id: PC-2 U
Data File: 2M22215.D
Analysis Date: 07/30/07 20:46
Date Rec/Extracted: 07/21/07-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	56-23-5	Carbon Tetrachloride	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	108-90-7	Chlorobenzene	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	67-66-3	Chloroform	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	74-87-3	Chloromethane	5.0	U
156-59-2	cis-1,2-Dichloroethene	5.0	U	107-06-2	1,2-Dichloroethane	5.0	U
10061-01-5	cis-1,3-Dichloropropene	5.0	U	78-87-5	1,2-Dichloropropane	5.0	U
124-48-1	Dibromochloromethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
1330-20-7	m&p-Xylenes	2.0	U	78-93-3	2-Butanone	5.0	U
75-09-2	Methylene Chloride	5.0	U	591-78-6	2-Hexanone	5.0	U
95-47-6	o-Xylene	1.0	U	108-10-1	4-Methyl-2-Pentanone	5.0	U
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	3.5	75-27-4	Bromodichloromethane	5.0	U
156-60-5	trans-1,2-Dichloroethene	5.0	U	75-25-2	Bromoforn	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
75-01-4	Vinyl Chloride	5.0	U				

Worksheet #: 54676

Total Target Concentration 3.5

*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: AC31901-010
 Client Id: PC-2 U
 Data File: 5M33174.D
 Analysis Date: 07/26/07 17:43
 Date Rec/Extracted: 07/21/07-07/26/07

Matrix: Aqueous
 Initial Vol: 500ml
 Final Vol: 0.5ml
 Dilution: 1
 Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	10	U	205-99-2	Benzo[b]fluoranthene	10	U
95-50-1	1,2-Dichlorobenzene	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
541-73-1	1,3-Dichlorobenzene	10	U	207-08-9	Benzo[k]fluoranthene	10	U
106-46-7	1,4-Dichlorobenzene	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
120-83-2	2,4-Dichlorophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
105-67-9	2,4-Dimethylphenol	10	U	85-68-7	Butylbenzylphthalate	10	U
51-28-5	2,4-Dinitrophenol	25	U	86-74-8	Carbazole	10	U
121-14-2	2,4-Dinitrotoluene	10	U	218-01-9	Chrysene	10	U
606-20-2	2,6-Dinitrotoluene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
91-58-7	2-Chloronaphthalene	10	U	132-64-9	Dibenzofuran	10	U
95-57-8	2-Chlorophenol	10	U	84-66-2	Diethylphthalate	10	U
91-57-6	2-Methylnaphthalene	10	U	131-11-3	Dimethylphthalate	10	U
95-48-7	2-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
88-74-4	2-Nitroaniline	10	U	117-84-0	Di-n-octylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	206-44-0	Fluoranthene	10	U
106-44-5	3&4-Methylphenol	10	U	86-73-7	Fluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	118-74-1	Hexachlorobenzene	10	U
99-09-2	3-Nitroaniline	10	U	87-68-3	Hexachlorobutadiene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	25	U	77-47-4	Hexachlorocyclopentadiene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
106-47-8	4-Chloroaniline	10	U	78-59-1	Isophorone	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	91-20-3	Naphthalene	10	U
100-01-6	4-Nitroaniline	10	U	98-95-3	Nitrobenzene	10	U
100-02-7	4-Nitrophenol	25	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
83-32-9	Acenaphthene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
208-96-8	Acenaphthylene	10	U	87-86-5	Pentachlorophenol	25	U
120-12-7	Anthracene	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 #: 54667

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: AC31901-018
 Client Id: PC-3 F
 Matrix: AQUEOUS
 Level: LOW

% Solid: 0
 Units: UG/L
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

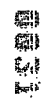
Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-36-0	Antimony	3.9	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-38-2	Arsenic	4.9	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-39-3	Barium	0.45	85		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-41-7	Beryllium	0.24	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-43-9	Cadmium	0.78	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-70-2	Calcium	20	41000		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-47-3	Chromium	5.8	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-48-4	Cobalt	0.54	0.94		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-50-8	Copper	1.5	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7439-89-6	Iron	75	2800		108/02/07	8159	W8159C2m31	31	P	PEICP2
7439-92-1	Lead	1.7	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7439-95-4	Magnesium	63	12000		108/02/07	8159	W8159C2m31	31	P	PEICP2
7439-96-5	Manganese	4.7	1800		108/02/07	8159	W8159C2m31	31	P	PEICP2
7439-97-6	Mercury	0.21	ND		108/02/07	8159	8159SWMD25	25	CV	HGCV1
7440-02-0	Nickel	1.1	1.9		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-09-7	Potassium	52	3800		108/02/07	8159	W8159B2M31	31	P	PEICPRAD2
7782-49-2	Selenium	10	12		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-22-4	Silver	0.53	0.61		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-23-5	Sodium	130	75000		108/02/07	8159	W8159B2M31	31	P	PEICPRAD2
7440-28-0	Thallium	4.8	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-62-2	Vanadium	1.7	1.7		108/02/07	8159	W8159C2m31	31	P	PEICP2
7440-66-6	Zinc	1.5	ND		108/02/07	8159	W8159C2m31	31	P	PEICP2

Comments: _____

Flag Codes:

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



Form1

ORGANICS VOLATILE REPORT

Sample Number: AC31901-017
 Client Id: PC-3 U
 Data File: 2M22218.D
 Analysis Date: 07/30/07 22:01
 Date Rec/Extracted: 07/21/07-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	56-23-5	Carbon Tetrachloride	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	108-90-7	Chlorobenzene	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	67-66-3	Chloroform	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	74-87-3	Chloromethane	5.0	U
156-59-2	cis-1,2-Dichloroethene	5.0	U	107-06-2	1,2-Dichloroethane	5.0	U
10061-01-5	cis-1,3-Dichloropropene	5.0	U	78-87-5	1,2-Dichloropropane	5.0	U
124-48-1	Dibromochloromethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
1330-20-7	m&p-Xylenes	2.0	U	78-93-3	2-Butanone	5.0	U
75-09-2	Methylene Chloride	5.0	U	591-78-6	2-Hexanone	5.0	U
95-47-6	o-Xylene	1.0	U	108-10-1	4-Methyl-2-Pentanone	5.0	U
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	1.6	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
75-01-4	Vinyl Chloride	5.0	U				

Worksheet #: 54676

Total Target Concentration 1.6

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: AC31901-017
 Client Id: PC-3 U
 Data File: 5M33227.D
 Analysis Date: 07/27/07 18:53
 Date Rec/Extracted: 07/21/07-07/26/07

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

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	10	U	205-99-2	Benzo[b]fluoranthene	10	U
95-50-1	1,2-Dichlorobenzene	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
541-73-1	1,3-Dichlorobenzene	10	U	207-08-9	Benzo[k]fluoranthene	10	U
106-46-7	1,4-Dichlorobenzene	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
120-83-2	2,4-Dichlorophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
105-67-9	2,4-Dimethylphenol	10	U	85-68-7	Butylbenzylphthalate	10	U
51-28-5	2,4-Dinitrophenol	25	U	86-74-8	Carbazole	10	U
121-14-2	2,4-Dinitrotoluene	10	U	218-01-9	Chrysene	10	U
606-20-2	2,6-Dinitrotoluene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
91-58-7	2-Chloronaphthalene	10	U	132-64-9	Dibenzofuran	10	U
95-57-8	2-Chlorophenol	10	U	84-66-2	Diethylphthalate	10	U
91-57-6	2-Methylnaphthalene	10	U	131-11-3	Dimethylphthalate	10	U
95-48-7	2-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
88-74-4	2-Nitroaniline	10	U	117-84-0	Di-n-octylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	206-44-0	Fluoranthene	10	U
106-44-5	3&4-Methylphenol	10	U	86-73-7	Fluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	118-74-1	Hexachlorobenzene	10	U
99-09-2	3-Nitroaniline	10	U	87-68-3	Hexachlorobutadiene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	25	U	77-47-4	Hexachlorocyclopentadiene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
106-47-8	4-Chloroaniline	10	U	78-59-1	Isophorone	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	91-20-3	Naphthalene	10	U
100-01-6	4-Nitroaniline	10	U	98-95-3	Nitrobenzene	10	U
100-02-7	4-Nitrophenol	25	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
83-32-9	Acenaphthene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
208-96-8	Acenaphthylene	10	U	87-86-5	Pentachlorophenol	25	U
120-12-7	Anthracene	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 #: 54667

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.

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Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-019
Client Id: TB
Data File: 3M37660.D
Analysis Date: 07/26/07 12:20
Date Rec/Extracted: 07/21/07-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	56-23-5	Carbon Tetrachloride	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	108-90-7	Chlorobenzene	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	67-66-3	Chloroform	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	74-87-3	Chloromethane	5.0	U
95-50-1	1,2-Dichlorobenzene	5.0	U	156-59-2	cis-1,2-Dichloroethene	5.0	U
107-06-2	1,2-Dichloroethane	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
78-87-5	1,2-Dichloropropane	5.0	U	124-48-1	Dibromochloromethane	5.0	U
541-73-1	1,3-Dichlorobenzene	5.0	U	100-41-4	Ethylbenzene	1.0	U
106-46-7	1,4-Dichlorobenzene	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
78-93-3	2-Butanone	5.0	U	75-09-2	Methylene Chloride	5.0	U
591-78-6	2-Hexanone	5.0	U	95-47-6	o-Xylene	1.0	U
108-10-1	4-Methyl-2-Pentanone	5.0	U	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	11
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	75-01-4	Vinyl Chloride	5.0	U

Worksheet #: 54676

Total Target Concentration 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 Inorganic Analysis Data Sheet

Sample ID: AC31901-005
 Client Id: SW-1
 Matrix: AQUEOUS
 Level: LOW

% Solid: 0
 Units: UG/L
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	47000		108/02/07	8159	W8159C2m22		P	PEICP2
7440-36-0	Antimony	3.9	ND		108/02/07	8159	W8159C2m22		P	PEICP2
7440-38-2	Arsenic	4.9	10		108/02/07	8159	W8159C2m22		P	PEICP2
7440-39-3	Barium	0.45	360		108/02/07	8159	W8159C2m22		P	PEICP2
7440-41-7	Beryllium	0.24	1.0		108/02/07	8159	W8159C2m22		P	PEICP2
7440-43-9	Cadmium	0.78	0.80		108/02/07	8159	W8159C2m22		P	PEICP2
7440-70-2	Calcium	20	71000		108/02/07	8159	W8159C2m22		P	PEICP2
7440-47-3	Chromium	5.8	73		108/02/07	8159	W8159C2m22		P	PEICP2
7440-48-4	Cobalt	0.54	21		108/02/07	8159	W8159C2m22		P	PEICP2
7440-50-8	Copper	1.5	97		108/02/07	8159	W8159C2m22		P	PEICP2
7439-89-6	Iron	75	78000		108/02/07	8159	W8159C2m22		P	PEICP2
7439-92-1	Lead	1.7	140		108/02/07	8159	W8159C2m22		P	PEICP2
7439-95-4	Magnesium	63	26000		108/02/07	8159	W8159C2m22		P	PEICP2
7439-96-5	Manganese	4.7	5400		108/02/07	8159	W8159C2m22		P	PEICP2
7439-97-6	Mercury	0.21	0.27		108/02/07	8159	8159SWMD18		CV	HGCV1
7440-02-0	Nickel	1.1	46		108/02/07	8159	W8159C2m22		P	PEICP2
7440-09-7	Potassium	52	11000		108/02/07	8159	W8159B2M122		P	PEICPRAD2
7782-49-2	Selenium	10	17		108/02/07	8159	W8159C2m22		P	PEICP2
7440-22-4	Silver	0.53	3.4		108/02/07	8159	W8159C2m22		P	PEICP2
7440-23-5	Sodium	130	45000		108/02/07	8159	W8159B2M122		P	PEICPRAD2
7440-28-0	Thallium	4.8	ND		108/02/07	8159	W8159C2m22		P	PEICP2
7440-62-2	Vanadium	1.7	110		108/02/07	8159	W8159C2m22		P	PEICP2
7440-66-6	Zinc	1.5	270		108/02/07	8159	W8159C2m22		P	PEICP2

Comments: _____

Flag Codes:

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

6233

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-005
Client Id: SW-1
Data File: 2M22212.D
Analysis Date: 07/30/07 19:31
Date Rec/Extracted: 07/21/07-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
79-00-5	1,1,2-Trichloroethane	5.0	U	74-87-3	Chloromethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	156-59-2	cis-1,2-Dichloroethene	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	10061-01-5	cis-1,3-Dichloropropene	5.0	U
107-06-2	1,2-Dichloroethane	5.0	U	124-48-1	Dibromochloromethane	5.0	U
78-87-5	1,2-Dichloropropane	5.0	U	100-41-4	Ethylbenzene	1.0	U
78-93-3	2-Butanone	5.0	U	1330-20-7	m&p-Xylenes	2.0	U
591-78-6	2-Hexanone	5.0	U	75-09-2	Methylene Chloride	5.0	U
108-10-1	4-Methyl-2-Pentanone	5.0	U	95-47-6	o-Xylene	1.0	U
67-64-1	Acetone	25	U	100-42-5	Styrene	5.0	U
71-43-2	Benzene	1.0	U	127-18-4	Tetrachloroethene	5.0	U
75-27-4	Bromodichloromethane	5.0	U	108-88-3	Toluene	1.0	16
75-25-2	Bromoform	5.0	U	156-60-5	trans-1,2-Dichloroethene	5.0	U
74-83-9	Bromomethane	5.0	U	10061-02-6	trans-1,3-Dichloropropene	5.0	U
75-15-0	Carbon Disulfide	5.0	U	79-01-6	Trichloroethene	5.0	U
56-23-5	Carbon Tetrachloride	5.0	U	75-01-4	Vinyl Chloride	5.0	U
108-90-7	Chlorobenzene	5.0	U				

Worksheet #: 54680

Total Target Concentration 16

*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: AC31901-005
 Client Id: SW-1
 Data File: 5M33232.D
 Analysis Date: 07/27/07 20:44
 Date Rec/Extracted: 07/21/07-07/25/07

Matrix: Aqueous
 Initial Vol: 960ml
 Final Vol: 1ml
 Dilution: 1
 Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	10	U	205-99-2	Benzo[b]fluoranthene	10	U
95-50-1	1,2-Dichlorobenzene	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
541-73-1	1,3-Dichlorobenzene	10	U	207-08-9	Benzo[k]fluoranthene	10	U
106-46-7	1,4-Dichlorobenzene	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
120-83-2	2,4-Dichlorophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
105-67-9	2,4-Dimethylphenol	10	U	85-68-7	Butylbenzylphthalate	10	U
51-28-5	2,4-Dinitrophenol	26	U	86-74-8	Carbazole	10	U
121-14-2	2,4-Dinitrotoluene	10	U	218-01-9	Chrysene	10	U
606-20-2	2,6-Dinitrotoluene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
91-58-7	2-Chloronaphthalene	10	U	132-64-9	Dibenzofuran	10	U
95-57-8	2-Chlorophenol	10	U	84-66-2	Diethylphthalate	10	U
91-57-6	2-Methylnaphthalene	10	U	131-11-3	Dimethylphthalate	10	U
95-48-7	2-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
88-74-4	2-Nitroaniline	10	U	117-84-0	Di-n-octylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	206-44-0	Fluoranthene	10	U
106-44-5	3&4-Methylphenol	10	U	86-73-7	Fluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	118-74-1	Hexachlorobenzene	10	U
99-09-2	3-Nitroaniline	10	U	87-68-3	Hexachlorobutadiene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	26	U	77-47-4	Hexachlorocyclopentadiene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
106-47-8	4-Chloroaniline	10	U	78-59-1	Isophorone	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	91-20-3	Naphthalene	10	U
100-01-6	4-Nitroaniline	10	U	98-95-3	Nitrobenzene	10	U
100-02-7	4-Nitrophenol	26	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
83-32-9	Acenaphthene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
208-96-8	Acenaphthylene	10	U	87-86-5	Pentachlorophenol	26	U
120-12-7	Anthracene	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 #: 54667

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: AC31901-001
 Client Id: SW-2
 Matrix: AQUEOUS
 Level: LOW

% Solid: 0
 Units: UG/L
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	93		108/02/07	8159	W8159C2m20		P	PEICP2
7440-36-0	Antimony	3.9	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-38-2	Arsenic	4.9	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-39-3	Barium	0.45	26		108/02/07	8159	W8159C2m20		P	PEICP2
7440-41-7	Beryllium	0.24	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-43-9	Cadmium	0.78	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-70-2	Calcium	20	44000		108/02/07	8159	W8159C2m20		P	PEICP2
7440-47-3	Chromium	5.8	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-48-4	Cobalt	0.54	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-50-8	Copper	1.5	22		108/02/07	8159	W8159C2m20		P	PEICP2
7439-89-6	Iron	75	450		108/02/07	8159	W8159C2m20		P	PEICP2
7439-92-1	Lead	1.7	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7439-95-4	Magnesium	63	13000		108/02/07	8159	W8159C2m20		P	PEICP2
7439-96-5	Manganese	4.7	520		108/02/07	8159	W8159C2m20		P	PEICP2
7439-97-6	Mercury	0.21	ND		108/02/07	8159	8159SWMD17		CV	HGCV1
7440-02-0	Nickel	1.1	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-09-7	Potassium	52	3500		108/02/07	8159	W8159B2M120		P	PEICPRAD2
7782-49-2	Selenium	10	14		108/02/07	8159	W8159C2m20		P	PEICP2
7440-22-4	Silver	0.53	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-23-5	Sodium	130	11000		108/02/07	8159	W8159B2M120		P	PEICPRAD2
7440-28-0	Thallium	4.8	ND		108/02/07	8159	W8159C2m20		P	PEICP2
7440-62-2	Vanadium	1.7	2.4		108/02/07	8159	W8159C2m20		P	PEICP2
7440-66-6	Zinc	1.5	ND		108/02/07	8159	W8159C2m20		P	PEICP2

Comments: _____

Flag Codes:

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

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-001
Client Id: SW-2
Data File: 2M22211.D
Analysis Date: 07/30/07 19:06
Date Rec/Extracted: 07/21/07-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	56-23-5	Carbon Tetrachloride	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	108-90-7	Chlorobenzene	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	67-66-3	Chloroform	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	74-87-3	Chloromethane	5.0	U
156-59-2	cis-1,2-Dichloroethene	5.0	U	107-06-2	1,2-Dichloroethane	5.0	U
10061-01-5	cis-1,3-Dichloropropene	5.0	U	78-87-5	1,2-Dichloropropane	5.0	U
124-48-1	Dibromochloromethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
1330-20-7	m&p-Xylenes	2.0	U	78-93-3	2-Butanone	5.0	U
75-09-2	Methylene Chloride	5.0	U	591-78-6	2-Hexanone	5.0	U
95-47-6	o-Xylene	1.0	U	108-10-1	4-Methyl-2-Pentanone	5.0	U
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	2.8	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
75-01-4	Vinyl Chloride	5.0	U				

Worksheet #: 54676

Total Target Concentration 2.8

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.

Form 1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC31901-001
 Client Id: SW-2
 Data File: 5M33153.D
 Analysis Date: 07/26/07 09:24
 Date Rec/Extracted: 07/21/07-07/25/07

Matrix: Aqueous
 Initial Vol: 500ml
 Final Vol: 0.5ml
 Dilution: 1
 Solids: 0

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	10	U	205-99-2	Benzo[b]fluoranthene	10	U
95-50-1	1,2-Dichlorobenzene	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
541-73-1	1,3-Dichlorobenzene	10	U	207-08-9	Benzo[k]fluoranthene	10	U
106-46-7	1,4-Dichlorobenzene	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
120-83-2	2,4-Dichlorophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
105-67-9	2,4-Dimethylphenol	10	U	85-68-7	Butylbenzylphthalate	10	U
51-28-5	2,4-Dinitrophenol	25	U	86-74-8	Carbazole	10	U
121-14-2	2,4-Dinitrotoluene	10	U	218-01-9	Chrysene	10	U
606-20-2	2,6-Dinitrotoluene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
91-58-7	2-Chloronaphthalene	10	U	132-64-9	Dibenzofuran	10	U
95-57-8	2-Chlorophenol	10	U	84-66-2	Diethylphthalate	10	U
91-57-6	2-Methylnaphthalene	10	U	131-11-3	Dimethylphthalate	10	U
95-48-7	2-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
88-74-4	2-Nitroaniline	10	U	117-84-0	Di-n-octylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	206-44-0	Fluoranthene	10	U
106-44-5	3&4-Methylphenol	10	U	86-73-7	Fluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	118-74-1	Hexachlorobenzene	10	U
99-09-2	3-Nitroaniline	10	U	87-68-3	Hexachlorobutadiene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	25	U	77-47-4	Hexachlorocyclopentadiene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
106-47-8	4-Chloroaniline	10	U	78-59-1	Isophorone	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	91-20-3	Naphthalene	10	U
100-01-6	4-Nitroaniline	10	U	98-95-3	Nitrobenzene	10	U
100-02-7	4-Nitrophenol	25	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
83-32-9	Acenaphthene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
208-96-8	Acenaphthylene	10	U	87-86-5	Pentachlorophenol	25	U
120-12-7	Anthracene	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 #: 54667

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: AC31901-014
 Client Id: SW-4
 Matrix: AQUEOUS
 Level: LOW

% Solid: 0
 Units: UG/L
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	2600		108/02/07	8159	W8159C2m13		P	PEICP2
7440-36-0	Antimony	3.9	ND		108/02/07	8159	W8159C2m13		P	PEICP2
7440-38-2	Arsenic	4.9	ND		108/02/07	8159	W8159C2m13		P	PEICP2
7440-39-3	Barium	0.45	94		108/02/07	8159	W8159C2m13		P	PEICP2
7440-41-7	Beryllium	0.24	ND		108/02/07	8159	W8159C2m13		P	PEICP2
7440-43-9	Cadmium	0.78	ND		108/02/07	8159	W8159C2m13		P	PEICP2
7440-70-2	Calcium	20	47000		108/02/07	8159	W8159C2m13		P	PEICP2
7440-47-3	Chromium	5.8	6.9		108/02/07	8159	W8159C2m13		P	PEICP2
7440-48-4	Cobalt	0.54	2.5		108/02/07	8159	W8159C2m13		P	PEICP2
7440-50-8	Copper	1.5	35		108/02/07	8159	W8159C2m13		P	PEICP2
7439-89-6	Iron	75	8200		108/02/07	8159	W8159C2m13		P	PEICP2
7439-92-1	Lead	1.7	15		108/02/07	8159	W8159C2m13		P	PEICP2
7439-95-4	Magnesium	63	14000		108/02/07	8159	W8159C2m13		P	PEICP2
7439-96-5	Manganese	4.7	2600		108/02/07	8159	W8159C2m13		P	PEICP2
7439-97-6	Mercury	0.21	ND		108/02/07	8159	8159SWMD13		CV	HGCV1
7440-02-0	Nickel	1.1	5.9		108/02/07	8159	W8159C2m13		P	PEICP2
7440-09-7	Potassium	52	3800		108/02/07	8159	W8159B2M13		P	PEICPRAD2
7782-49-2	Selenium	10	17		108/02/07	8159	W8159C2m13		P	PEICP2
7440-22-4	Silver	0.53	0.78		108/02/07	8159	W8159C2m13		P	PEICP2
7440-23-5	Sodium	130	12000		108/02/07	8159	W8159B2M13		P	PEICPRAD2
7440-28-0	Thallium	4.8	ND		108/02/07	8159	W8159C2m13		P	PEICP2
7440-62-2	Vanadium	1.7	9.7		108/02/07	8159	W8159C2m13		P	PEICP2
7440-66-6	Zinc	1.5	38		108/02/07	8159	W8159C2m13		P	PEICP2

Comments: _____

Flag Codes:

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

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-014
 Client Id: SW-4
 Data File: 2M22217.D
 Analysis Date: 07/30/07 21:36
 Date Rec/Extracted: 07/21/07-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	56-23-5	Carbon Tetrachloride	5.0	U
79-34-5	1,1,2,2-Tetrachloroethane	5.0	U	108-90-7	Chlorobenzene	5.0	U
79-00-5	1,1,2-Trichloroethane	5.0	U	75-00-3	Chloroethane	5.0	U
75-34-3	1,1-Dichloroethane	5.0	U	67-66-3	Chloroform	5.0	U
75-35-4	1,1-Dichloroethene	5.0	U	74-87-3	Chloromethane	5.0	U
156-59-2	cis-1,2-Dichloroethene	5.0	U	107-06-2	1,2-Dichloroethane	5.0	U
10061-01-5	cis-1,3-Dichloropropene	5.0	U	78-87-5	1,2-Dichloropropane	5.0	U
124-48-1	Dibromochloromethane	5.0	U	100-41-4	Ethylbenzene	1.0	U
1330-20-7	m&p-Xylenes	2.0	U	78-93-3	2-Butanone	5.0	U
75-09-2	Methylene Chloride	5.0	U	591-78-6	2-Hexanone	5.0	U
95-47-6	o-Xylene	1.0	U	108-10-1	4-Methyl-2-Pentanone	5.0	U
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	1.8	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
75-01-4	Vinyl Chloride	5.0	U				

Worksheet #: 54676

Total Target Concentration 1.8

*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 SEMIVOLATILE REPORT

Sample Number: AC31901-014
 Client Id: SW-4
 Data File: 5M33228.D
 Analysis Date: 07/27/07 19:15
 Date Rec/Extracted: 07/21/07-07/26/07

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

Units: ug/L

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	10	U	205-99-2	Benzo[h]fluoranthene	10	U
95-50-1	1,2-Dichlorobenzene	10	U	191-24-2	Benzo[g,h,i]perylene	10	U
541-73-1	1,3-Dichlorobenzene	10	U	207-08-9	Benzo[k]fluoranthene	10	U
106-46-7	1,4-Dichlorobenzene	10	U	111-91-1	bis(2-Chloroethoxy)methan	10	U
95-95-4	2,4,5-Trichlorophenol	10	U	111-44-4	bis(2-Chloroethyl)ether	10	U
88-06-2	2,4,6-Trichlorophenol	10	U	39638-32-9	bis(2-chloroisopropyl)ether	10	U
120-83-2	2,4-Dichlorophenol	10	U	117-81-7	bis(2-Ethylhexyl)phthalate	10	U
105-67-9	2,4-Dimethylphenol	10	U	85-68-7	Butylbenzylphthalate	10	U
51-28-5	2,4-Dinitrophenol	25	U	86-74-8	Carbazole	10	U
121-14-2	2,4-Dinitrotoluene	10	U	218-01-9	Chrysene	10	U
606-20-2	2,6-Dinitrotoluene	10	U	53-70-3	Dibenzo[a,h]anthracene	10	U
91-58-7	2-Chloronaphthalene	10	U	132-64-9	Dibenzofuran	10	U
95-57-8	2-Chlorophenol	10	U	84-66-2	Diethylphthalate	10	U
91-57-6	2-Methylnaphthalene	10	U	131-11-3	Dimethylphthalate	10	U
95-48-7	2-Methylphenol	10	U	84-74-2	Di-n-butylphthalate	10	U
88-74-4	2-Nitroaniline	10	U	117-84-0	Di-n-octylphthalate	10	U
88-75-5	2-Nitrophenol	10	U	206-44-0	Fluoranthene	10	U
106-44-5	3&4-Methylphenol	10	U	86-73-7	Fluorene	10	U
91-94-1	3,3'-Dichlorobenzidine	10	U	118-74-1	Hexachlorobenzene	10	U
99-09-2	3-Nitroaniline	10	U	87-68-3	Hexachlorobutadiene	10	U
534-52-1	4,6-Dinitro-2-methylphenol	25	U	77-47-4	Hexachlorocyclopentadiene	10	U
101-55-3	4-Bromophenyl-phenylether	10	U	67-72-1	Hexachloroethane	10	U
59-50-7	4-Chloro-3-methylphenol	10	U	193-39-5	Indeno[1,2,3-cd]pyrene	10	U
106-47-8	4-Chloroaniline	10	U	78-59-1	Isophorone	10	U
7005-72-3	4-Chlorophenyl-phenylether	10	U	91-20-3	Naphthalene	10	U
100-01-6	4-Nitroaniline	10	U	98-95-3	Nitrobenzene	10	U
100-02-7	4-Nitrophenol	25	U	621-64-7	N-Nitroso-di-n-propylamine	10	U
83-32-9	Acenaphthene	10	U	86-30-6	n-Nitrosodiphenylamine	10	U
208-96-8	Acenaphthylene	10	U	87-86-5	Pentachlorophenol	25	U
120-12-7	Anthracene	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 #: 54667

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.

Form 1 Inorganic Analysis Data Sheet

Sample ID: AC31901-004
 Client Id: SD-1
 Matrix: SOIL
 Level: LOW

% Solid: 57
 Units: MG/KG
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	17	5700	100	07/25/07	8156	38156BMDI22	22	P	PEICPRAD1
7440-36-0	Antimony	0.98	ND	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-38-2	Arsenic	0.42	1.5	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-39-3	Barium	0.045	34	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-41-7	Beryllium	0.015	0.15	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-43-9	Cadmium	0.033	0.13	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-70-2	Calcium	41	8800	100	07/25/07	8156	38156BMD122	22	P	PEICPRAD1
7440-47-3	Chromium	0.61	12	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-48-4	Cobalt	0.071	3.5	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-50-8	Copper	0.55	17	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7439-89-6	Iron	11	9900	100	07/25/07	8156	38156BMD122	22	P	PEICPRAD1
7439-92-1	Lead	0.89	18	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7439-95-4	Magnesium	12	5900	100	07/25/07	8156	38156BMD122	22	P	PEICPRAD1
7439-96-5	Manganese	2.1	130	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7439-97-6	Mercury	0.019	ND	167	07/25/07	8156	38156SMD118	18	CV	HGCV1
7440-02-0	Nickel	0.27	7.4	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-09-7	Potassium	33	710	100	07/25/07	8156	38156BMD122	22	P	PEICPRAD1
7782-49-2	Selenium	0.85	ND	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-22-4	Silver	0.10	ND	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-23-5	Sodium	26	78	100	07/25/07	8156	38156BMD122	22	P	PEICPRAD1
7440-28-0	Thallium	0.81	ND	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-62-2	Vanadium	0.044	17	100	07/24/07	8156	38156AMD122	22	P	PEICP1
7440-66-6	Zinc	0.66	35	100	07/24/07	8156	38156AMD122	22	P	PEICP1

Comments: _____

Flag Codes:

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

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-004
Client Id: SD-1
Data File: 1M23379.D
Analysis Date: 07/24/07 21:55
Date Rec/Extracted: 07/21/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 57

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0088	U	56-23-5	Carbon Tetrachloride	0.0088	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0088	U	108-90-7	Chlorobenzene	0.0088	U
79-00-5	1,1,2-Trichloroethane	0.0088	U	75-00-3	Chloroethane	0.0088	U
75-34-3	1,1-Dichloroethane	0.0088	U	67-66-3	Chloroform	0.0088	U
75-35-4	1,1-Dichloroethene	0.0088	U	74-87-3	Chloromethane	0.0088	U
95-50-1	1,2-Dichlorobenzene	0.0088	U	156-59-2	cis-1,2-Dichloroethene	0.0088	U
107-06-2	1,2-Dichloroethane	0.0088	U	10061-01-5	cis-1,3-Dichloropropene	0.0088	U
78-87-5	1,2-Dichloropropane	0.0088	U	124-48-1	Dibromochloromethane	0.0088	U
541-73-1	1,3-Dichlorobenzene	0.0088	U	100-41-4	Ethylbenzene	0.0088	U
106-46-7	1,4-Dichlorobenzene	0.0088	U	1330-20-7	m&p-Xylenes	0.018	U
78-93-3	2-Butanone	0.018	U	75-09-2	Methylene Chloride	0.018	0.055 B
591-78-6	2-Hexanone	0.018	U	95-47-6	o-Xylene	0.0088	U
108-10-1	4-Methyl-2-Pentanone	0.0088	U	100-42-5	Styrene	0.0088	U
67-64-1	Acetone	0.088	0.13	127-18-4	Tetrachloroethene	0.0088	U
71-43-2	Benzene	0.0088	U	108-88-3	Toluene	0.0088	U
75-27-4	Bromodichloromethane	0.0088	U	156-60-5	trans-1,2-Dichloroethene	0.0088	U
75-25-2	Bromoform	0.0088	U	10061-02-6	trans-1,3-Dichloropropene	0.0088	U
74-83-9	Bromomethane	0.0088	U	79-01-6	Trichloroethene	0.0088	U
75-15-0	Carbon Disulfide	0.0088	0.0070 J	75-01-4	Vinyl Chloride	0.0088	U

Worksheet #: 54676

Total Target Concentration 0.192

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: AC31901-004
 Client Id: SD-1
 Data File: 9M04810.D
 Analysis Date: 07/30/07 18:38
 Date Rec/Extracted: 07/21/07-07/24/07

Matrix: Soil
 Initial Vol: 30g
 Final Vol: 1ml
 Dilution: 1
 Solids: 57

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	0.58	U	205-99-2	Benzo[b]fluoranthene	0.58	U
191-24-2	Benzo[g,h,i]perylene	0.58	U	207-08-9	Benzo[k]fluoranthene	0.58	U
111-91-1	bis(2-Chloroethoxy)methan	0.58	U	95-95-4	2,4,5-Trichlorophenol	0.58	U
111-44-4	bis(2-Chloroethyl)ether	0.58	U	88-06-2	2,4,6-Trichlorophenol	0.58	U
39638-32-9	bis(2-chloroisopropyl)ether	0.58	U	120-83-2	2,4-Dichlorophenol	0.58	U
117-81-7	bis(2-Ethylhexyl)phthalate	0.58	U	105-67-9	2,4-Dimethylphenol	0.58	U
85-68-7	Butylbenzylphthalate	0.58	U	51-28-5	2,4-Dinitrophenol	1.5	U
86-74-8	Carbazole	0.58	U	121-14-2	2,4-Dinitrotoluene	0.58	U
218-01-9	Chrysene	0.58	U	606-20-2	2,6-Dinitrotoluene	0.58	U
53-70-3	Dibenzo[a,h]anthracene	0.58	U	91-58-7	2-Chloronaphthalene	0.58	U
132-64-9	Dibenzofuran	0.58	U	95-57-8	2-Chlorophenol	0.58	U
84-66-2	Diethylphthalate	0.58	U	91-57-6	2-Methylnaphthalene	0.58	U
131-11-3	Dimethylphthalate	0.58	U	95-48-7	2-Methylphenol	0.58	U
84-74-2	Di-n-butylphthalate	0.58	0.071 JB	88-74-4	2-Nitroaniline	0.58	U
117-84-0	Di-n-octylphthalate	0.58	U	88-75-5	2-Nitrophenol	0.58	U
206-44-0	Fluoranthene	0.58	U	106-44-5	3&4-Methylphenol	0.58	U
86-73-7	Fluorene	0.58	U	91-94-1	3,3'-Dichlorobenzidine	0.58	U
118-74-1	Hexachlorobenzene	0.58	U	99-09-2	3-Nitroaniline	0.58	U
87-68-3	Hexachlorobutadiene	0.58	U	534-52-1	4,6-Dinitro-2-methylphenol	1.5	U
77-47-4	Hexachlorocyclopentadiene	1.5	U	101-55-3	4-Bromophenyl-phenylether	0.58	U
67-72-1	Hexachloroethane	0.58	U	59-50-7	4-Chloro-3-methylphenol	0.58	U
193-39-5	Indeno[1,2,3-cd]pyrene	0.58	U	106-47-8	4-Chloroaniline	0.58	U
78-59-1	Isophorone	0.58	U	7005-72-3	4-Chlorophenyl-phenylether	0.58	U
91-20-3	Naphthalene	0.58	U	100-01-6	4-Nitroaniline	0.58	U
98-95-3	Nitrobenzene	0.58	U	100-02-7	4-Nitrophenol	0.58	U
621-64-7	N-Nitroso-di-n-propylamine	0.58	U	83-32-9	Acenaphthene	0.58	U
86-30-6	n-Nitrosodiphenylamine	0.58	U	208-96-8	Acenaphthylene	0.58	U
87-86-5	Pentachlorophenol	2.9	U	120-12-7	Anthracene	0.58	U
85-01-8	Phenanthrene	0.58	U	56-55-3	Benzo[a]anthracene	0.58	U
108-95-2	Phenol	0.58	U	50-32-8	Benzo[a]pyrene	0.58	U
129-00-0	Pyrene	0.58	U				

Worksheet #: 54667

Total Target Concentration 0.071

*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: AC31901-002
Client Id: SD-2
Matrix: SOIL
Level: LOW

% Solid: 79
Units: MG/KG
Date Rec: 7/23/2007

Lab Name: Veritech
Lab Code:
Contract:

Nras No:
Sdg No:
Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	12	3700	100	07/25/07	8156	8156BMDL13	13	P	PEICPRAD1
7440-36-0	Antimony	0.71	ND	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-38-2	Arsenic	0.30	2.0	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-39-3	Barium	0.032	52	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-41-7	Beryllium	0.011	0.042	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-43-9	Cadmium	0.024	0.18	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-70-2	Calcium	29	42000	100	07/25/07	8156	8156BMDL13	13	P	PEICPRAD1
7440-47-3	Chromium	0.44	7.4	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-48-4	Cobalt	0.051	3.6	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-50-8	Copper	0.40	11	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7439-89-6	Iron	8.2	11000	100	07/25/07	8156	8156BMDL13	13	P	PEICPRAD1
7439-92-1	Lead	0.64	11	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7439-95-4	Magnesium	9.0	25000	100	07/25/07	8156	8156BMDL13	13	P	PEICPRAD1
7439-96-5	Manganese	1.5	1600	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7439-97-6	Mercury	0.014	ND	167	07/25/07	8156	8156SMD13	13	CV	HGCV1
7440-02-0	Nickel	0.19	8.1	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-09-7	Potassium	24	920	100	07/25/07	8156	8156BMDL13	13	P	PEICPRAD1
7782-49-2	Selenium	0.61	ND	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-22-4	Silver	0.075	0.16	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-23-5	Sodium	19	53	100	07/25/07	8156	8156BMDL13	13	P	PEICPRAD1
7440-28-0	Thallium	0.58	ND	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-62-2	Vanadium	0.032	14	100	07/24/07	8156	8156AMD13	13	P	PEICP1
7440-66-6	Zinc	0.47	33	100	07/24/07	8156	8156AMD13	13	P	PEICP1

Comments: _____

Flag Codes:

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

CT #: PH-0671
 MA #: NJ386
 NJ #: 14622

NY #: 11408
 PA #: 68-463

Report Of Analysis

veritech laboratories

To: HDR/LMS

7 Coates Drive
 Suite 2
 Goshen

NY 10924

Attention: T. Schneider
 Project: Harrison Landfill

Date Collected: 10/29/2007
 Date Submitted: 10/30/2007
 Date Reported: 11/13/2007

Lab#: AC33918-001

SampleID: SD-2

DRAFT

TestGroup	Analyte	DF	Units	MDL/PQL/RL	RESULT
AC33918-001	SD-2				
	Volatile Organics (no search) 8260				
	1,1,1-Trichloroethane	1	mg/Kg	0.029	ND
	1,1,2,2-Tetrachloroethane	1	mg/Kg	0.029	ND
	1,1,2-trichloro-1,2,2-trifluoroethane	1	mg/Kg	0.029	ND
	1,1,2-Trichloroethane	1	mg/Kg	0.029	ND
	1,1-Dichloroethane	1	mg/Kg	0.029	ND
	1,1-Dichloroethene	1	mg/Kg	0.029	ND
	1,2,3-Trichloropropane	1	mg/Kg	0.029	ND
	1,2,4-Trimethylbenzene	1	mg/Kg	0.0059	ND
	1,2-Dibromo-3-chloropropane	1	mg/Kg	0.029	ND
	1,2-Dibromoethane	1	mg/Kg	0.029	ND
	1,2-Dichlorobenzene	1	mg/Kg	0.029	ND
	1,2-Dichloroethane	1	mg/Kg	0.029	ND
	1,2-Dichloropropane	1	mg/Kg	0.029	ND
	1,3,5-Trimethylbenzene	1	mg/Kg	0.0059	ND
	1,3-Dichlorobenzene	1	mg/Kg	0.029	ND
	1,3-Dichloropropane	1	mg/Kg	0.029	ND
	1,4-Dichlorobenzene	1	mg/Kg	0.029	ND
	1,4-Dioxane	1	mg/Kg	1.5	ND
	2-Butanone	1	mg/Kg	0.029	ND
	2-Chloroethylvinylether	1	mg/Kg	0.029	ND
	2-Hexanone	1	mg/Kg	0.029	ND
	4-Isopropyltoluene	1	mg/Kg	0.0059	ND
	4-Methyl-2-Pentanone	1	mg/Kg	0.029	ND
	Acetone	1	mg/Kg	0.15	ND
	Benzene	1	mg/Kg	0.0059	ND
	Bromodichloromethane	1	mg/Kg	0.029	ND
	Bromoform	1	mg/Kg	0.029	ND
	Bromomethane	1	mg/Kg	0.029	ND
	Carbon disulfide	1	mg/Kg	0.029	0.0064J
	Carbon tetrachloride	1	mg/Kg	0.029	ND
	Chlorobenzene	1	mg/Kg	0.029	ND
	Chloroethane	1	mg/Kg	0.029	ND
	Chloroform	1	mg/Kg	0.029	ND
	Chloromethane	1	mg/Kg	0.029	ND
	cis-1,2-Dichloroethene	1	mg/Kg	0.029	ND
	cis-1,3-Dichloropropene	1	mg/Kg	0.029	ND
	Cyclohexane	1	mg/Kg	0.029	ND
	Dibromochloromethane	1	mg/Kg	0.029	ND
	Dichlorodifluoromethane	1	mg/Kg	0.029	ND
	Ethylbenzene	1	mg/Kg	0.0059	ND
	Isopropylbenzene	1	mg/Kg	0.0059	ND
	m&p-Xylenes	1	mg/Kg	0.012	ND
	Methyl Acetate	1	mg/Kg	0.029	ND
	Methylcyclohexane	1	mg/Kg	0.029	ND

ND = Not Detected

Lab#: AC33918-001

SampleID: SD-2

DRAFT

TestGroup	Analyte	DF	Units	MDL/PQL/RL	RESULT
	Methylene chloride	1	mg/Kg	0.029	0.060B
	Methyl-t-butyl ether	1	mg/Kg	0.0059	ND
	n-Butylbenzene	1	mg/Kg	0.0059	ND
	n-Propylbenzene	1	mg/Kg	0.0059	ND
	o-Xylene	1	mg/Kg	0.0059	ND
	sec-Butylbenzene	1	mg/Kg	0.0059	ND
	Styrene	1	mg/Kg	0.029	ND
	t-Butyl Alcohol	1	mg/Kg	0.15	ND
	t-Butylbenzene	1	mg/Kg	0.0059	ND
	Tetrachloroethene	1	mg/Kg	0.029	ND
	Toluene	1	mg/Kg	0.0059	ND
	trans-1,2-Dichloroethene	1	mg/Kg	0.029	ND
	trans-1,3-dichloropropene	1	mg/Kg	0.029	ND
	Trichloroethene	1	mg/Kg	0.029	ND
	Trichlorofluoromethane	1	mg/Kg	0.029	ND
	Vinyl chloride	1	mg/Kg	0.029	ND

RL Definitions: SW846 Organics reported to PQL
 SW846 Inorganics reported to PQL

Clean Water Act Organics reported to MDL
 Clean Water Act Inorganics reported to PQL

CLP Organics reported to CRDL
 CLP Inorganics reported to CRDL

This report is a true report of results obtained from our tests of this material. In lieu of a formal contract document, the total aggregate liability of Veritech to all parties shall not exceed Veritech's total fee for analytical services rendered.

Or

 Jeri Rossi - Quality Assurance Director

 Stanley Gilewicz - Laboratory Director

ND = Not Detected

 Veritech Report Of Analysis
 175 Route 46 West, Unit D, Fairfield, NJ 07004

Veritech Project: 7103008

Page 2 of 2

Form1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC31901-002
 Client Id: SD-2
 Data File: 9M04878.D
 Analysis Date: 08/03/07 09:46
 Date Rec/Extracted: 07/21/07-07/24/07

Matrix: Soil
 Initial Vol: 30g
 Final Vol: 1ml
 Dilution: 1
 Solids: 79

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	0.42	U	205-99-2	Benzo[b]fluoranthene	0.42	U
191-24-2	Benzo[g,h,i]perylene	0.42	U	207-08-9	Benzo[k]fluoranthene	0.42	U
111-91-1	bis(2-Chloroethoxy)methan	0.42	U	95-95-4	2,4,5-Trichlorophenol	0.42	U
111-44-4	bis(2-Chloroethyl)ether	0.42	U	88-06-2	2,4,6-Trichlorophenol	0.42	U
39638-32-9	bis(2-chloroisopropyl)ether	0.42	U	120-83-2	2,4-Dichlorophenol	0.42	U
117-81-7	bis(2-Ethylhexyl)phthalate	0.42	U	105-67-9	2,4-Dimethylphenol	0.42	U
85-68-7	Butylbenzylphthalate	0.42	U	51-28-5	2,4-Dinitrophenol	1.1	U
86-74-8	Carbazole	0.42	U	121-14-2	2,4-Dinitrotoluene	0.42	U
218-01-9	Chrysene	0.42	U	606-20-2	2,6-Dinitrotoluene	0.42	U
53-70-3	Dibenzo[a,h]anthracene	0.42	U	91-58-7	2-Chloronaphthalene	0.42	U
132-64-9	Dibenzofuran	0.42	U	95-57-8	2-Chlorophenol	0.42	U
84-66-2	Diethylphthalate	0.42	U	91-57-6	2-Methylnaphthalene	0.42	U
131-11-3	Dimethylphthalate	0.42	U	95-48-7	2-Methylphenol	0.42	U
84-74-2	Di-n-butylphthalate	0.42	U	88-74-4	2-Nitroaniline	0.42	U
117-84-0	Di-n-octylphthalate	0.42	U	88-75-5	2-Nitrophenol	0.42	U
206-44-0	Fluoranthene	0.42	U	106-44-5	3&4-Methylphenol	0.42	U
86-73-7	Fluorene	0.42	U	91-94-1	3,3'-Dichlorobenzidine	0.42	U
118-74-1	Hexachlorobenzene	0.42	U	99-09-2	3-Nitroaniline	0.42	U
87-68-3	Hexachlorobutadiene	0.42	U	534-52-1	4,6-Dinitro-2-methylphenol	1.1	U
77-47-4	Hexachlorocyclopentadiene	0.42	U	101-55-3	4-Bromophenyl-phenylether	0.42	U
67-72-1	Hexachloroethane	0.42	U	59-50-7	4-Chloro-3-methylphenol	0.42	U
193-39-5	Indeno[1,2,3-cd]pyrene	0.42	U	106-47-8	4-Chloroaniline	0.42	U
78-59-1	Isophorone	0.42	U	7005-72-3	4-Chlorophenyl-phenylether	0.42	U
91-20-3	Naphthalene	0.42	U	100-01-6	4-Nitroaniline	0.42	U
98-95-3	Nitrobenzene	0.42	U	100-02-7	4-Nitrophenol	0.42	U
621-64-7	N-Nitroso-di-n-propylamine	0.42	U	83-32-9	Acenaphthene	0.42	U
86-30-6	n-Nitrosodiphenylamine	0.42	U	208-96-8	Acenaphthylene	0.42	U
87-86-5	Pentachlorophenol	1.1	U	120-12-7	Anthracene	0.42	U
85-01-8	Phenanthrene	0.42	U	56-55-3	Benzo[a]anthracene	0.42	U
108-95-2	Phenol	0.42	U	50-32-8	Benzo[a]pyrene	0.42	U
129-00-0	Pyrene	0.42	0.051 J				

Worksheet #: 54667

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 Inorganic Analysis Data Sheet

Sample ID: AC31901-003
 Client Id: SD-3
 Matrix: SOIL
 Level: LOW

% Solid: 55
 Units: MG/KG
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	18	8800	100	07/25/07	8156	38156BMDL20	20	P	PEICPRAD1
7440-36-0	Antimony	1.0	ND	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-38-2	Arsenic	0.43	3.0	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-39-3	Barium	0.046	54	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-41-7	Beryllium	0.016	0.37	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-43-9	Cadmium	0.034	0.55	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-70-2	Calcium	42	23000	100	07/25/07	8156	38156BMDL20	20	P	PEICPRAD1
7440-47-3	Chromium	0.63	15	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-48-4	Cobalt	0.073	5.4	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-50-8	Copper	0.57	22	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7439-89-6	Iron	12	13000	100	07/25/07	8156	38156BMDL20	20	P	PEICPRAD1
7439-92-1	Lead	0.92	120	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7439-95-4	Magnesium	13	12000	100	07/25/07	8156	38156BMDL20	20	P	PEICPRAD1
7439-96-5	Manganese	2.2	200	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7439-97-6	Mercury	0.020	ND	167	07/25/07	8156	38156SMD117	17	CV	HGCV1
7440-02-0	Nickel	0.28	12	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-09-7	Potassium	34	880	100	07/25/07	8156	38156BMDL20	20	P	PEICPRAD1
7782-49-2	Selenium	0.88	ND	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-22-4	Silver	0.11	ND	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-23-5	Sodium	27	180	100	07/25/07	8156	38156BMDL20	20	P	PEICPRAD1
7440-28-0	Thallium	0.83	ND	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-62-2	Vanadium	0.045	25	100	07/24/07	8156	38156AMD120	20	P	PEICP1
7440-66-6	Zinc	0.68	81	100	07/24/07	8156	38156AMD120	20	P	PEICP1

Comments: _____

Flag Codes:

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

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-003
 Client Id: SD-3
 Data File: 1M23378.D
 Analysis Date: 07/24/07 21:28
 Date Rec/Extracted: 07/21/07-NA

Matrix: Soil
 Initial Vol: 5g
 Final Vol: NA
 Dilution: 1
 Solids: 55

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0091	U	56-23-5	Carbon Tetrachloride	0.0091	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0091	U	108-90-7	Chlorobenzene	0.0091	U
79-00-5	1,1,2-Trichloroethane	0.0091	U	75-00-3	Chloroethane	0.0091	U
75-34-3	1,1-Dichloroethane	0.0091	U	67-66-3	Chloroform	0.0091	U
75-35-4	1,1-Dichloroethene	0.0091	U	74-87-3	Chloromethane	0.0091	U
95-50-1	1,2-Dichlorobenzene	0.0091	U	156-59-2	cis-1,2-Dichloroethene	0.0091	U
107-06-2	1,2-Dichloroethane	0.0091	U	10061-01-5	cis-1,3-Dichloropropene	0.0091	U
78-87-5	1,2-Dichloropropane	0.0091	U	124-48-1	Dibromochloromethane	0.0091	U
541-73-1	1,3-Dichlorobenzene	0.0091	U	100-41-4	Ethylbenzene	0.0091	U
106-46-7	1,4-Dichlorobenzene	0.0091	U	1330-20-7	m&p-Xylenes	0.018	U
78-93-3	2-Butanone	0.018	U	75-09-2	Methylene Chloride	0.018	0.041 B
591-78-6	2-Hexanone	0.018	U	95-47-6	o-Xylene	0.0091	U
108-10-1	4-Methyl-2-Pentanone	0.0091	U	100-42-5	Styrene	0.0091	U
67-64-1	Acetone	0.091	0.13	127-18-4	Tetrachloroethene	0.0091	U
71-43-2	Benzene	0.0091	U	108-88-3	Toluene	0.0091	U
75-27-4	Bromodichloromethane	0.0091	U	156-60-5	trans-1,2-Dichloroethene	0.0091	U
75-25-2	Bromoform	0.0091	U	10061-02-6	trans-1,3-Dichloropropene	0.0091	U
74-83-9	Bromomethane	0.0091	U	79-01-6	Trichloroethene	0.0091	U
75-15-0	Carbon Disulfide	0.0091	0.0020 J	75-01-4	Vinyl Chloride	0.0091	U

Worksheet #: 54676

Total Target Concentration 0.173

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: AC31901-003
 Client Id: SD-3
 Data File: 9M04809.D
 Analysis Date: 07/30/07 18:15
 Date Rec/Extracted: 07/21/07-07/24/07

Matrix: Soil
 Initial Vol: 30g
 Final Vol: 1ml
 Dilution: 1
 Solids: 55

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	0.61	U	205-99-2	Benzo[b]fluoranthene	0.61	0.14 J
191-24-2	Benzo[g,h,i]perylene	0.61	0.082 J	207-08-9	Benzo[k]fluoranthene	0.61	0.080 J
111-91-1	bis(2-Chloroethoxy)methan	0.61	U	95-95-4	2,4,5-Trichlorophenol	0.61	U
111-44-4	bis(2-Chloroethyl)ether	0.61	U	88-06-2	2,4,6-Trichlorophenol	0.61	U
39638-32-9	bis(2-chloroisopropyl)ether	0.61	U	120-83-2	2,4-Dichlorophenol	0.61	U
117-81-7	bis(2-Ethylhexyl)phthalate	0.61	U	105-67-9	2,4-Dimethylphenol	0.61	U
85-68-7	Butylbenzylphthalate	0.61	U	51-28-5	2,4-Dinitrophenol	1.5	U
86-74-8	Carbazole	0.61	U	121-14-2	2,4-Dinitrotoluene	0.61	U
218-01-9	Chrysene	0.61	0.16 J	606-20-2	2,6-Dinitrotoluene	0.61	U
53-70-3	Dibenzo[a,h]anthracene	0.61	U	91-58-7	2-Chloronaphthalene	0.61	U
132-64-9	Dibenzofuran	0.61	U	95-57-8	2-Chlorophenol	0.61	U
84-66-2	Diethylphthalate	0.61	U	91-57-6	2-Methylnaphthalene	0.61	U
131-11-3	Dimethylphthalate	0.61	U	95-48-7	2-Methylphenol	0.61	U
84-74-2	Di-n-butylphthalate	0.61	U	88-74-4	2-Nitroaniline	0.61	U
117-84-0	Di-n-octylphthalate	0.61	U	88-75-5	2-Nitrophenol	0.61	U
206-44-0	Fluoranthene	0.61	0.34 J	106-44-5	3&4-Methylphenol	0.61	U
86-73-7	Fluorene	0.61	U	91-94-1	3,3'-Dichlorobenzidine	0.61	U
118-74-1	Hexachlorobenzene	0.61	U	99-09-2	3-Nitroaniline	0.61	U
87-68-3	Hexachlorobutadiene	0.61	U	534-52-1	4,6-Dinitro-2-methylphenol	1.5	U
77-47-4	Hexachlorocyclopentadiene	1.5	U	101-55-3	4-Bromophenyl-phenylether	0.61	U
67-72-1	Hexachloroethane	0.61	U	59-50-7	4-Chloro-3-methylphenol	0.61	U
193-39-5	Indeno[1,2,3-cd]pyrene	0.61	0.078 J	106-47-8	4-Chloroaniline	0.61	U
78-59-1	Isophorone	0.61	U	7005-72-3	4-Chlorophenyl-phenylether	0.61	U
91-20-3	Naphthalene	0.61	U	100-01-6	4-Nitroaniline	0.61	U
98-95-3	Nitrobenzene	0.61	U	100-02-7	4-Nitrophenol	0.61	U
621-64-7	N-Nitroso-di-n-propylamine	0.61	U	83-32-9	Acenaphthene	0.61	U
86-30-6	n-Nitrosodiphenylamine	0.61	U	208-96-8	Acenaphthylene	0.61	U
87-86-5	Pentachlorophenol	3.0	U	120-12-7	Anthracene	0.61	0.072 J
85-01-8	Phenanthrene	0.61	0.31 J	56-55-3	Benzo[a]anthracene	0.61	0.18 J
108-95-2	Phenol	0.61	U	50-32-8	Benzo[a]pyrene	0.61	0.12 J
129-00-0	Pyrene	0.61	0.43 J				

Worksheet #: 54667

Total Target Concentration 1.992

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: AC31901-015
Client Id: SD-4
Matrix: SOIL
Level: LOW

% Solid: 71
Units: MG/KG
Date Rec: 7/23/2007

Lab Name: Veritech
Lab Code:
Contract:

Nras No:
Sdg No:
Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	14	3600	100	07/25/07	8156	S8156BMDL	23	P	PEICPRAD1
7440-36-0	Antimony	0.78	ND	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-38-2	Arsenic	0.34	2.7	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-39-3	Barium	0.036	36	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-41-7	Beryllium	0.012	0.088	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-43-9	Cadmium	0.026	0.24	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-70-2	Calcium	33	9400	100	07/25/07	8156	S8156BMD	23	P	PEICPRAD1
7440-47-3	Chromium	0.49	8.2	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-48-4	Cobalt	0.057	3.5	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-50-8	Copper	0.44	18	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7439-89-6	Iron	9.1	12000	100	07/25/07	8156	S8156BMD	23	P	PEICPRAD1
7439-92-1	Lead	0.71	19	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7439-95-4	Magnesium	10	6700	100	07/25/07	8156	S8156BMD	23	P	PEICPRAD1
7439-96-5	Manganese	1.7	590	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7439-97-6	Mercury	0.016	ND	167	07/25/07	8156	S8156SMD	19	CV	HGCV1
7440-02-0	Nickel	0.22	9.1	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-09-7	Potassium	26	560	100	07/25/07	8156	S8156BMD	23	P	PEICPRAD1
7782-49-2	Selenium	0.68	ND	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-22-4	Silver	0.083	0.088	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-23-5	Sodium	21	51	100	07/25/07	8156	S8156BMD	23	P	PEICPRAD1
7440-28-0	Thallium	0.65	ND	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-62-2	Vanadium	0.035	11	100	07/24/07	8156	S8156AMD	23	P	PEICP1
7440-66-6	Zinc	0.53	56	100	07/24/07	8156	S8156AMD	23	P	PEICP1

Comments: _____

Flag Codes:

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

Form1

ORGANICS VOLATILE REPORT

Sample Number: AC31901-015
 Client Id: SD-4
 Data File: 1M23380.D
 Analysis Date: 07/24/07 22:22
 Date Rec/Extracted: 07/21/07-NA

Matrix: Soil
 Initial Vol: 5g
 Final Vol: NA
 Dilution: 1
 Solids: 71

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0070	U	56-23-5	Carbon Tetrachloride	0.0070	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0070	U	108-90-7	Chlorobenzene	0.0070	U
79-00-5	1,1,2-Trichloroethane	0.0070	U	75-00-3	Chloroethane	0.0070	U
75-34-3	1,1-Dichloroethane	0.0070	U	67-66-3	Chloroform	0.0070	U
75-35-4	1,1-Dichloroethene	0.0070	U	74-87-3	Chloromethane	0.0070	U
95-50-1	1,2-Dichlorobenzene	0.0070	U	156-59-2	cis-1,2-Dichloroethene	0.0070	U
107-06-2	1,2-Dichloroethane	0.0070	U	10061-01-5	cis-1,3-Dichloropropene	0.0070	U
78-87-5	1,2-Dichloropropane	0.0070	U	124-48-1	Dibromochloromethane	0.0070	U
541-73-1	1,3-Dichlorobenzene	0.0070	U	100-41-4	Ethylbenzene	0.0070	U
106-46-7	1,4-Dichlorobenzene	0.0070	U	1330-20-7	m&p-Xylenes	0.014	U
78-93-3	2-Butanone	0.014	U	75-09-2	Methylene Chloride	0.014	0.038 B
591-78-6	2-Hexanone	0.014	U	95-47-6	o-Xylene	0.0070	U
108-10-1	4-Methyl-2-Pentanone	0.0070	U	100-42-5	Styrene	0.0070	U
67-64-1	Acetone	0.070	0.078	127-18-4	Tetrachloroethene	0.0070	U
71-43-2	Benzene	0.0070	U	108-88-3	Toluene	0.0070	U
75-27-4	Bromodichloromethane	0.0070	U	156-60-5	trans-1,2-Dichloroethene	0.0070	U
75-25-2	Bromoform	0.0070	U	10061-02-6	trans-1,3-Dichloropropene	0.0070	U
74-83-9	Bromomethane	0.0070	U	79-01-6	Trichloroethene	0.0070	U
75-15-0	Carbon Disulfide	0.0070	0.0032 J	75-01-4	Vinyl Chloride	0.0070	U

Worksheet #: 54676

Total Target Concentration 0.1192

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.

Form 1

ORGANICS SEMIVOLATILE REPORT

Sample Number: AC31901-015
 Client Id: SD-4
 Data File: 9M04768.D
 Analysis Date: 07/27/07 22:08
 Date Rec/Extracted: 07/21/07-07/24/07

Matrix: Soil
 Initial Vol: 30g
 Final Vol: 1ml
 Dilution: 1
 Solids: 71

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	0.47	U	205-99-2	Benzo[b]fluoranthene	0.47	U
191-24-2	Benzo[g,h,i]perylene	0.47	U	207-08-9	Benzo[k]fluoranthene	0.47	U
111-91-1	bis(2-Chloroethoxy)methan	0.47	U	95-95-4	2,4,5-Trichlorophenol	0.47	U
111-44-4	bis(2-Chloroethyl)ether	0.47	U	88-06-2	2,4,6-Trichlorophenol	0.47	U
39638-32-9	bis(2-chloroisopropyl)ether	0.47	U	120-83-2	2,4-Dichlorophenol	0.47	U
117-81-7	bis(2-Ethylhexyl)phthalate	0.47	U	105-67-9	2,4-Dimethylphenol	0.47	U
85-68-7	Butylbenzylphthalate	0.47	U	51-28-5	2,4-Dinitrophenol	1.2	U
86-74-8	Carbazole	0.47	U	121-14-2	2,4-Dinitrotoluene	0.47	U
218-01-9	Chrysene	0.47	U	606-20-2	2,6-Dinitrotoluene	0.47	U
53-70-3	Dibenzo[a,h]anthracene	0.47	U	91-58-7	2-Chloronaphthalene	0.47	U
132-64-9	Dibenzofuran	0.47	U	95-57-8	2-Chlorophenol	0.47	U
84-66-2	Diethylphthalate	0.47	U	91-57-6	2-Methylnaphthalene	0.47	U
131-11-3	Dimethylphthalate	0.47	U	95-48-7	2-Methylphenol	0.47	U
84-74-2	Di-n-butylphthalate	0.47	U	88-74-4	2-Nitroaniline	0.47	U
117-84-0	Di-n-octylphthalate	0.47	U	88-75-5	2-Nitrophenol	0.47	U
206-44-0	Fluoranthene	0.47	U	106-44-5	3&4-Methylphenol	0.47	U
86-73-7	Fluorene	0.47	U	91-94-1	3,3'-Dichlorobenzidine	0.47	U
118-74-1	Hexachlorobenzene	0.47	U	99-09-2	3-Nitroaniline	0.47	U
87-68-3	Hexachlorobutadiene	0.47	U	534-52-1	4,6-Dinitro-2-methylphenol	1.2	U
77-47-4	Hexachlorocyclopentadiene	1.2	U	101-55-3	4-Bromophenyl-phenylether	0.47	U
67-72-1	Hexachloroethane	0.47	U	59-50-7	4-Chloro-3-methylphenol	0.47	U
193-39-5	Indeno[1,2,3-cd]pyrene	0.47	U	106-47-8	4-Chloroaniline	0.47	U
78-59-1	Isophorone	0.47	U	7005-72-3	4-Chlorophenyl-phenylether	0.47	U
91-20-3	Naphthalene	0.47	U	100-01-6	4-Nitroaniline	0.47	U
98-95-3	Nitrobenzene	0.47	U	100-02-7	4-Nitrophenol	0.47	U
621-64-7	N-Nitroso-di-n-propylamine	0.47	U	83-32-9	Acenaphthene	0.47	U
86-30-6	n-Nitrosodiphenylamine	0.47	U	208-96-8	Acenaphthylene	0.47	U
87-86-5	Pentachlorophenol	2.3	U	120-12-7	Anthracene	0.47	U
85-01-8	Phenanthrene	0.47	U	56-55-3	Benzo[a]anthracene	0.47	U
108-95-2	Phenol	0.47	U	50-32-8	Benzo[a]pyrene	0.47	U
129-00-0	Pyrene	0.47	U				

Worksheet #: 54667

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: AC31901-016
 Client Id: SD-5
 Matrix: SOIL
 Level: LOW

% Solid: 69
 Units: MG/KG
 Date Rec: 7/23/2007

Lab Name: Veritech
 Lab Code:
 Contract:

Nras No:
 Sdg No:
 Case No:

Cas No.	Analyte	RL	Conc	Dil Fact	Analysis Date:	Prep Batch	File:	Seq Num:	M	Instr
7429-90-5	Aluminum	14	3700	100	07/25/07	8156	38156BMDI	24	P	PEICPRAD1
7440-36-0	Antimony	0.81	ND	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-38-2	Arsenic	0.35	1.3	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-39-3	Barium	0.037	47	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-41-7	Beryllium	0.012	0.057	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-43-9	Cadmium	0.027	0.27	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-70-2	Calcium	34	11000	100	07/25/07	8156	38156BMD	24	P	PEICPRAD1
7440-47-3	Chromium	0.50	8.6	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-48-4	Cobalt	0.059	3.1	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-50-8	Copper	0.45	8.7	100	07/24/07	8156	38156AMD	24	P	PEICP1
7439-89-6	Iron	9.3	10000	100	07/25/07	8156	38156BMD	24	P	PEICPRAD1
7439-92-1	Lead	0.73	16	100	07/24/07	8156	38156AMD	24	P	PEICP1
7439-95-4	Magnesium	10	7100	100	07/25/07	8156	38156BMD	24	P	PEICPRAD1
7439-96-5	Manganese	1.7	940	100	07/24/07	8156	38156AMD	24	P	PEICP1
7439-97-6	Mercury	0.016	ND	167	07/25/07	8156	38156SMD	22	CV	HGCV1
7440-02-0	Nickel	0.22	8.6	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-09-7	Potassium	27	760	100	07/25/07	8156	38156BMD	24	P	PEICPRAD1
7782-49-2	Selenium	0.70	ND	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-22-4	Silver	0.086	0.12	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-23-5	Sodium	21	58	100	07/25/07	8156	38156BMD	24	P	PEICPRAD1
7440-28-0	Thallium	0.67	ND	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-62-2	Vanadium	0.036	12	100	07/24/07	8156	38156AMD	24	P	PEICP1
7440-66-6	Zinc	0.54	58	100	07/24/07	8156	38156AMD	24	P	PEICP1

Comments: _____

Flag Codes:

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

Form1
ORGANICS VOLATILE REPORT

Sample Number: AC31901-016
Client Id: SD-5
Data File: 1M23381.D
Analysis Date: 07/24/07 22:49
Date Rec/Extracted: 07/21/07-NA

Matrix: Soil
Initial Vol: 5g
Final Vol: NA
Dilution: 1
Solids: 69

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
71-55-6	1,1,1-Trichloroethane	0.0072	U	56-23-5	Carbon Tetrachloride	0.0072	U
79-34-5	1,1,2,2-Tetrachloroethane	0.0072	U	108-90-7	Chlorobenzene	0.0072	U
79-00-5	1,1,2-Trichloroethane	0.0072	U	75-00-3	Chloroethane	0.0072	U
75-34-3	1,1-Dichloroethane	0.0072	U	67-66-3	Chloroform	0.0072	U
75-35-4	1,1-Dichloroethene	0.0072	U	74-87-3	Chloromethane	0.0072	U
95-50-1	1,2-Dichlorobenzene	0.0072	U	156-59-2	cis-1,2-Dichloroethene	0.0072	U
107-06-2	1,2-Dichloroethane	0.0072	U	10061-01-5	cis-1,3-Dichloropropene	0.0072	U
78-87-5	1,2-Dichloropropane	0.0072	U	124-48-1	Dibromochloromethane	0.0072	U
541-73-1	1,3-Dichlorobenzene	0.0072	U	100-41-4	Ethylbenzene	0.0072	U
106-46-7	1,4-Dichlorobenzene	0.0072	U	1330-20-7	m&p-Xylenes	0.014	U
78-93-3	2-Butanone	0.014	U	75-09-2	Methylene Chloride	0.014	0.036 B
591-78-6	2-Hexanone	0.014	U	95-47-6	o-Xylene	0.0072	U
108-10-1	4-Methyl-2-Pentanone	0.0072	U	100-42-5	Styrene	0.0072	U
67-64-1	Acetone	0.072	0.14	127-18-4	Tetrachloroethene	0.0072	U
71-43-2	Benzene	0.0072	U	108-88-3	Toluene	0.0072	U
75-27-4	Bromodichloromethane	0.0072	U	156-60-5	trans-1,2-Dichloroethene	0.0072	U
75-25-2	Bromoform	0.0072	U	10061-02-6	trans-1,3-Dichloropropene	0.0072	U
74-83-9	Bromomethane	0.0072	U	79-01-6	Trichloroethene	0.0072	U
75-15-0	Carbon Disulfide	0.0072	0.0052 J	75-01-4	Vinyl Chloride	0.0072	U

Worksheet #: 54676

Total Target Concentration 0.1812

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: AC31901-016

Client Id: SD-5

Data File: 9M04767.D

Analysis Date: 07/27/07 21:45

Date Rec/Extracted: 07/21/07-07/24/07

Matrix: Soil

Initial Vol: 30g

Final Vol: 1ml

Dilution: 1

Solids: 69

Units: mg/Kg

Cas #	Compound	RL	Conc	Cas #	Compound	RL	Conc
120-82-1	1,2,4-Trichlorobenzene	0.48	U	205-99-2	Benzo[b]fluoranthene	0.48	U
191-24-2	Benzo[g,h,i]perylene	0.48	U	207-08-9	Benzo[k]fluoranthene	0.48	U
111-91-1	bis(2-Chloroethoxy)methan	0.48	U	95-95-4	2,4,5-Trichlorophenol	0.48	U
111-44-4	bis(2-Chloroethyl)ether	0.48	U	88-06-2	2,4,6-Trichlorophenol	0.48	U
39638-32-9	bis(2-chloroisopropyl)ether	0.48	U	120-83-2	2,4-Dichlorophenol	0.48	U
117-81-7	bis(2-Ethylhexyl)phthalate	0.48	U	105-67-9	2,4-Dimethylphenol	0.48	U
85-68-7	Butylbenzylphthalate	0.48	U	51-28-5	2,4-Dinitrophenol	1.2	U
86-74-8	Carbazole	0.48	U	121-14-2	2,4-Dinitrotoluene	0.48	U
218-01-9	Chrysene	0.48	U	606-20-2	2,6-Dinitrotoluene	0.48	U
53-70-3	Dibenzo[a,h]anthracene	0.48	U	91-58-7	2-Chloronaphthalene	0.48	U
132-64-9	Dibenzofuran	0.48	U	95-57-8	2-Chlorophenol	0.48	U
84-66-2	Diethylphthalate	0.48	U	91-57-6	2-Methylnaphthalene	0.48	U
131-11-3	Dimethylphthalate	0.48	U	95-48-7	2-Methylphenol	0.48	U
84-74-2	Di-n-butylphthalate	0.48	U	88-74-4	2-Nitroaniline	0.48	U
117-84-0	Di-n-octylphthalate	0.48	U	88-75-5	2-Nitrophenol	0.48	U
206-44-0	Fluoranthene	0.48	0.051 J	106-44-5	3&4-Methylphenol	0.48	U
86-73-7	Fluorene	0.48	U	91-94-1	3,3'-Dichlorobenzidine	0.48	U
118-74-1	Hexachlorobenzene	0.48	U	99-09-2	3-Nitroaniline	0.48	U
87-68-3	Hexachlorobutadiene	0.48	U	534-52-1	4,6-Dinitro-2-methylphenol	1.2	U
77-47-4	Hexachlorocyclopentadiene	1.2	U	101-55-3	4-Bromophenyl-phenylether	0.48	U
67-72-1	Hexachloroethane	0.48	U	59-50-7	4-Chloro-3-methylphenol	0.48	U
193-39-5	Indeno[1,2,3-cd]pyrene	0.48	U	106-47-8	4-Chloroaniline	0.48	U
78-59-1	Isophorone	0.48	U	7005-72-3	4-Chlorophenyl-phenylether	0.48	U
91-20-3	Naphthalene	0.48	U	100-01-6	4-Nitroaniline	0.48	U
98-95-3	Nitrobenzene	0.48	U	100-02-7	4-Nitrophenol	0.48	U
621-64-7	N-Nitroso-di-n-propylamine	0.48	U	83-32-9	Acenaphthene	0.48	U
86-30-6	n-Nitrosodiphenylamine	0.48	U	208-96-8	Acenaphthylene	0.48	U
87-86-5	Pentachlorophenol	2.4	U	120-12-7	Anthracene	0.48	U
85-01-8	Phenanthrene	0.48	U	56-55-3	Benzo[a]anthracene	0.48	U
108-95-2	Phenol	0.48	U	50-32-8	Benzo[a]pyrene	0.48	U
129-00-0	Pyrene	0.48	0.070 J				

Worksheet #: 54667

Total Target Concentration 0.121

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.

VERITECH Wet Chem Form1 Analysis Summary



Lab#: AC31901-001 Matrix Aqueous Client SampleID: SW-2	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICW	1	6.1	mg/L	1.5	07/24/07	07/25/07
Cyanide	CN-W-9010	1	0.014	mg/l	0.01	07/27/07	07/27/07

Lab#: AC31901-002 Matrix Soil Client SampleID: SD-2	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICS	1	ND	mg/kg	190	07/24/07	07/26/07
Cyanide	CN-SOIL	1	ND	mg/kg	0.32	08/01/07	08/01/07

Lab#: AC31901-003 Matrix Soil Client SampleID: SD-3	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICS	1	ND	mg/kg	270	07/24/07	07/26/07
Cyanide	CN-SOIL	1	0.47	mg/kg	0.45	08/01/07	08/01/07

Lab#: AC31901-004 Matrix Soil Client SampleID: SD-1	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICS	1	ND	mg/kg	260	07/24/07	07/26/07
Cyanide	CN-SOIL	1	ND	mg/kg	0.44	08/01/07	08/01/07

Lab#: AC31901-005 Matrix Aqueous Client SampleID: SW-1	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICW	1	15	mg/L	1.5	07/24/07	07/25/07
Cyanide	CN-W-9010	1	ND	mg/l	0.01	07/27/07	07/27/07

Lab#: AC31901-006 Matrix Aqueous Client SampleID: LMW-2 U	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICW	1	9.1	mg/L	1.5	07/24/07	07/25/07
Cyanide	CN-W-9010	1	ND	mg/l	0.01	07/27/07	07/27/07

Lab#: AC31901-008 Matrix Aqueous Client SampleID: MW-4 U	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
--	--

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICW	1	30	mg/L	1.5	07/24/07	07/25/07
Cyanide	CN-W-9010	1	ND	mg/l	0.01	07/27/07	07/27/07

Lab#: AC31901-010 Matrix Aqueous Client SampleID: PC-2 U	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICW	1	20	mg/L	1.5	07/24/07	07/25/07
Cyanide	CN-W-9010	1	ND	mg/l	0.01	07/27/07	07/27/07

VERITECH Wet Chem Form1 Analysis Summary

Lab#: AC31901-012 Matrix: Aqueous Client SampleID: PC-1 U	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICW	1	52	mg/L	1.5	07/24/07	07/25/07
Cyanide	CN-W-9010	1	ND	mg/l	0.01	07/27/07	07/27/07

Lab#: AC31901-014 Matrix: Aqueous Client SampleID: SW-4	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
---	--

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICW	1	6.4	mg/L	1.5	07/24/07	07/25/07
Cyanide	CN-W-9010	1	ND	mg/l	0.01	07/27/07	07/27/07

Lab#: AC31901-015 Matrix: Soil Client SampleID: SD-4	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
--	--

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICS	1	ND	mg/kg	210	07/24/07	07/26/07
Cyanide	CN-SOIL	1	ND	mg/kg	0.35	08/01/07	08/01/07

Lab#: AC31901-016 Matrix: Soil Client SampleID: SD-5	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
--	--

Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICS	1	ND	mg/kg	220	07/24/07	07/26/07
Cyanide	CN-SOIL	1	ND	mg/kg	0.36	08/01/07	08/01/07

Lab#: AC31901-017 Matrix: Aqueous Client SampleID: PC-3 U	Project Number: 7072310 Received Date: 7/21/2007 Collect Date: 7/20/2007
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Analysis	TestGroup	Dilution:	Result	Units:	PQL:	Prep Date:	Analysis Date:
Chloride	CHLORIDE-ICW	1	84	mg/L	1.5	07/24/07	07/25/07
Cyanide	CN-W-9010	1	0.012	mg/l	0.01	07/27/07	07/27/07

CT #: PH-0671
 MA #: NJ386
 NJ #: 14622

NY #: 11408
 PA #: 68-463

Report Of Analysis

veritech laboratories

To: HDR/LMS

7 Coates Drive
 Suite 2
 Goshen

NY 10924

Attention: T.Schneider
 Project: Harrison Landfill

Date Collected: 10/29/2007
 Date Submitted: 10/30/2007
 Date Reported: 11/13/2007

Lab#: AC33918-001	SampleID: SD-2	DRAFT
TestGroup	Analyte	DF Units MDL/PQL/RL RESULT

AC33918-001 **SD-2**

Volatile Organics (no search) 8260

1,1,1-Trichloroethane	1	mg/Kg	0.029	ND
1,1,2,2-Tetrachloroethane	1	mg/Kg	0.029	ND
1,1,2-trichloro-1,2,2-trifluoroethane	1	mg/Kg	0.029	ND
1,1,2-Trichloroethane	1	mg/Kg	0.029	ND
1,1-Dichloroethane	1	mg/Kg	0.029	ND
1,1-Dichloroethene	1	mg/Kg	0.029	ND
1,2,3-Trichloropropane	1	mg/Kg	0.029	ND
1,2,4-Trimethylbenzene	1	mg/Kg	0.0059	ND
1,2-Dibromo-3-chloropropane	1	mg/Kg	0.029	ND
1,2-Dibromoethane	1	mg/Kg	0.029	ND
1,2-Dichlorobenzene	1	mg/Kg	0.029	ND
1,2-Dichloroethane	1	mg/Kg	0.029	ND
1,2-Dichloropropane	1	mg/Kg	0.029	ND
1,3,5-Trimethylbenzene	1	mg/Kg	0.0059	ND
1,3-Dichlorobenzene	1	mg/Kg	0.029	ND
1,3-Dichloropropane	1	mg/Kg	0.029	ND
1,4-Dichlorobenzene	1	mg/Kg	0.029	ND
1,4-Dioxane	1	mg/Kg	1.5	ND
2-Butanone	1	mg/Kg	0.029	ND
2-Chloroethylvinylether	1	mg/Kg	0.029	ND
2-Hexanone	1	mg/Kg	0.029	ND
4-Isopropyltoluene	1	mg/Kg	0.0059	ND
4-Methyl-2-Pentanone	1	mg/Kg	0.029	ND
Acetone	1	mg/Kg	0.15	ND
Benzene	1	mg/Kg	0.0059	ND
Bromodichloromethane	1	mg/Kg	0.029	ND
Bromoform	1	mg/Kg	0.029	ND
Bromomethane	1	mg/Kg	0.029	ND
Carbon disulfide	1	mg/Kg	0.029	0.0064J
Carbon tetrachloride	1	mg/Kg	0.029	ND
Chlorobenzene	1	mg/Kg	0.029	ND
Chloroethane	1	mg/Kg	0.029	ND
Chloroform	1	mg/Kg	0.029	ND
Chloromethane	1	mg/Kg	0.029	ND
cis-1,2-Dichloroethene	1	mg/Kg	0.029	ND
cis-1,3-Dichloropropene	1	mg/Kg	0.029	ND
Cyclohexane	1	mg/Kg	0.029	ND
Dibromochloromethane	1	mg/Kg	0.029	ND
Dichlorodifluoromethane	1	mg/Kg	0.029	ND
Ethylbenzene	1	mg/Kg	0.0059	ND
Isopropylbenzene	1	mg/Kg	0.0059	ND
m&p-Xylenes	1	mg/Kg	0.012	ND
Methyl Acetate	1	mg/Kg	0.029	ND
Methylcyclohexane	1	mg/Kg	0.029	ND

ND = Not Detected

TestGroup	Analyte	DF	Units	MDL/PQL/RL	RESULT
	Methylene chloride	1	mg/Kg	0.029	0.060B
	Methyl-t-butyl ether	1	mg/Kg	0.0059	ND
	n-Butylbenzene	1	mg/Kg	0.0059	ND
	n-Propylbenzene	1	mg/Kg	0.0059	ND
	o-Xylene	1	mg/Kg	0.0059	ND
	sec-Butylbenzene	1	mg/Kg	0.0059	ND
	Styrene	1	mg/Kg	0.029	ND
	t-Butyl Alcohol	1	mg/Kg	0.15	ND
	t-Butylbenzene	1	mg/Kg	0.0059	ND
	Tetrachloroethene	1	mg/Kg	0.029	ND
	Toluene	1	mg/Kg	0.0059	ND
	trans-1,2-Dichloroethene	1	mg/Kg	0.029	ND
	trans-1,3-dichloropropene	1	mg/Kg	0.029	ND
	Trichloroethene	1	mg/Kg	0.029	ND
	Trichlorofluoromethane	1	mg/Kg	0.029	ND
	Vinyl chloride	1	mg/Kg	0.029	ND

RL Definitions: SW846 Organics reported to PQL
SW846 Inorganics reported to PQL

Clean Water Act Organics reported to MDL
Clean Water Act Inorganics reported to PQL

CLP Organics reported to CRDL
CLP Inorganics reported to CRDL

This report is a true report of results obtained from our tests of this material. In lieu of a formal contract document, the total aggregate liability of Veritech to all parties shall not exceed Veritech's total fee for analytical services rendered.

Or

Jeri Rossi - Quality Assurance Director

Stanley Gilewicz - Laboratory Director

