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October 23, 2006

Mr. John Lovejoy  
Bureau of Environmental Protection  
Nassau County Department of Health  
240 Old Country Rd.  
Mineola, New York 11501

**Re: Sanitary Leaching Pool Closure Work Plan  
Former Penetrex Processing Company  
Site No. 1-30-034  
1 Shore Road, Glenwood Landing, New York**

Dear Mr. Lovejoy:

P.W. Grosser Consulting, Inc. (PWGC) has prepared the following Sanitary Leaching Pool Closure Work Plan to document activities regarding the above-referenced site's leaching structures, and to detail procedures for their proper closure.

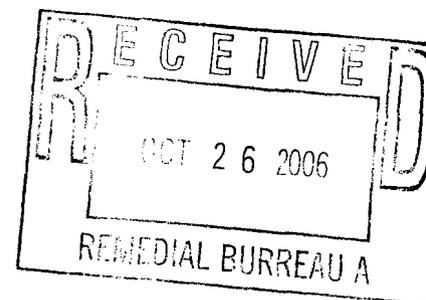
**Background**

The subject site is comprised of a one-acre parcel developed with a two-story commercial building, asphalt parking, communications tower, and other ancillary improvements, and is located at 1 Shore Road, Glenwood Landing, Town of North Hempstead, New York. The property is identified in Nassau County Tax maps as Section 20 – Block K – Lots 10 through 12. It is bound on its western boundary by Shore Road and on its eastern boundary by West Street. The site is located generally north of Scudders Lane and is situated near and adjoining several major oil storage facilities, coastal terminals, and municipal power stations near Hempstead Harbor. A site vicinity map is included as **Figure 1**.

The site is a Class II Inactive Hazardous Waste Site under NYSDEC oversight. Investigation activities continue at the site in an attempt to down-grade and/or delist the site from the NYSDEC registry.

In May 2003, a storm drain and sanitary leaching pool remediation was performed. Storm drains DW-1, DW-2, and DW-3, and sanitary leaching pools DW-4, DW-5, DW-6, DW-7, and the structures which comprise sanitary system A-1 were remediated and/or closed. UIC structures are indicated on **Figure 2**. Remedial Activities are documented in the "Storm Drain and Sanitary Leaching Pool Remediation and Closure Report," September 2003, previously submitted by PWGC under separate cover.

The United States Environmental Protection Agency (USEPA) acknowledged the satisfactory remediation or closure of the Class V UIC structures in a November 2003 letter. This letter is included as **Appendix A**.



**ACEC**

### **Geophysical Investigation**

PWGC conducted a geophysical investigation at the site on June 23, 2006 as part of a subsurface investigation deemed appropriate by the New York State Department of Environmental Protection (NYSDEC). The investigation included the use of a metal detector and ground penetrating radar (GPR) by NAEVA Geophysics, Inc. to locate subsurface structures in the parking area of the site. During the investigation, several "metal detector anomalies" were detected.

### **Leaching Pool Excavation and Sampling**

On August 24, 2006, excavation of the metal detector anomalies was performed at the site to identify them. Excavation activities were coordinated and overseen by PWGC. Excavation services were provided by Eastern Environmental Solutions. A representative from the NYSDEC was present to witness the activities.

Three of the anomalies uncovered were previously-undetected leaching structures. Upon discovery, these leaching structures were labeled DW-8, DW-9, and DW-10.

The concrete cover of leaching structure DW-8 was uncovered approximately 1.5 feet below grade surface. It was constructed of eight-foot diameter pre-cast concrete leaching rings. A four-inch diameter pipe entered the structure from the direction of DW-2, which had been abandoned. The depth to the bottom sediment was measured at 11 feet below grade surface. There was approximately four inches of water, believed to be groundwater as this corresponds to groundwater measurements in nearby monitoring wells.

The steel cover of DW-9 was uncovered approximately five inches below grade surface. It had been abandoned and was full of sand to a depth of five feet below grade. There was no visible piping.

The concrete cover of leaching structure DW-10 was uncovered approximately 2 feet below grade surface. It was constructed of eight-foot diameter pre-cast concrete leaching rings. As with DW-8, a four-inch diameter pipe entered the structure from the direction of the abandoned DW-2. The depth to the bottom sediment was measured at 13 feet below grade surface. There was approximately two feet of water, corresponding to the site's depth of groundwater.

Sediment samples were collected from the floors of structures DW-8 and DW-10 utilizing a stainless steel hand auger. The samples were placed in pre-cleaned laboratory supplied glassware and placed in a cooler packed with ice for transport to American Analytical Laboratories in Farmingdale, New York, an ELAP-certified laboratory.

The sampling of DW-9 occurred on August 30, 2006, as part of a NYSDEC soil boring investigation at the site. A Geoprobe™ was used to sample through the center of the leaching structure. Soils were collected continuously from the surface of the fill material to a depth of 20 feet below grade surface. Soils were characterized by a PWGC field hydrogeologist. The interface between the original bottom sand of the structure and the fill material added during the structure's previous abandonment was apparent due to the distinct difference in the color and grain size of the sands above and below. This interface was determined to be 13 feet below grade surface. The soil collected directly below the interface was collected in pre-cleaned laboratory supplied glassware and placed in a cooler packed with ice for transport to American Analytical Laboratories. The sample was labeled SB-8, 13' – 15' as it was collected from the eighth soil boring of the subsurface investigation.

Samples were submitted to the laboratory for analysis of VOCs by EPA method 8260, SVOCs by EPA method 8270, and the eight RCRA Metals in accordance with the NCDH procedures and protocols.

### **Analytical Results**

Analytical results were compared to Recommended Soil Cleanup Objectives (RSCOs) specified in the NYSDEC Technical and Administrative Guidance Memo (TAGM) #4046, to determine if remediation of the UIC structures is

warranted. Analytical data for samples DW-8 and DW-10 is shown on **Tables 1, 2, and 3**. Analytical data for sample SB-8, 13' – 15' (DW-9) is shown on **Tables 4, 5, and 6**. Laboratory data sheets are included as **Appendix B**.

The results indicate that concentrations of VOCs and metals detected in the three samples are within Recommended Soil Cleanup Objectives.

Concentrations of SVOCs detected in the three samples were within Cleanup Objectives with the exception of benzo(b)fluoranthene and benzo(a)pyrene in samples DW-8 and DW-10.

### **Recommended Closure Procedures**

Minor detections of VOCs reflect low level impact remaining at the site related to the original known contamination sources.

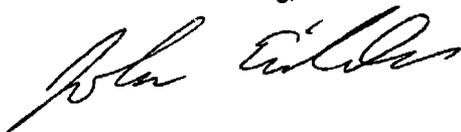
Although concentrations of two SVOCs exceeded recommended soil cleanup objectives, it is apparent that the structures did not receive improper discharge. SVOCs have not been detected in groundwater samples collected during the NYSDEC subsurface investigation of the site. For this reason, SVOCs had been previously eliminated as a concern by the NYSDEC.

PWGC recommends that the three UIC structures (DW-8, DW-9, and DW-10) be permanently abandoned to grade with clean, inert fill material, and re-paved over with asphalt. The covers and metal casings will be removed prior to abandonment. Continued groundwater sampling will be conducted as part of the NYSDEC investigation.

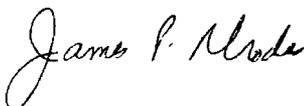
A United States Environmental Protection Agency (USEPA) Inventory of Injection Wells (EPA Form 7520-16) had been previously submitted in October 2003. A second Form 7520-16 will be submitted documenting the additional structures discovered at the site.

Please contact us if you have any questions or require additional information.

Very truly yours,  
**P.W. Grosser Consulting, Inc.**



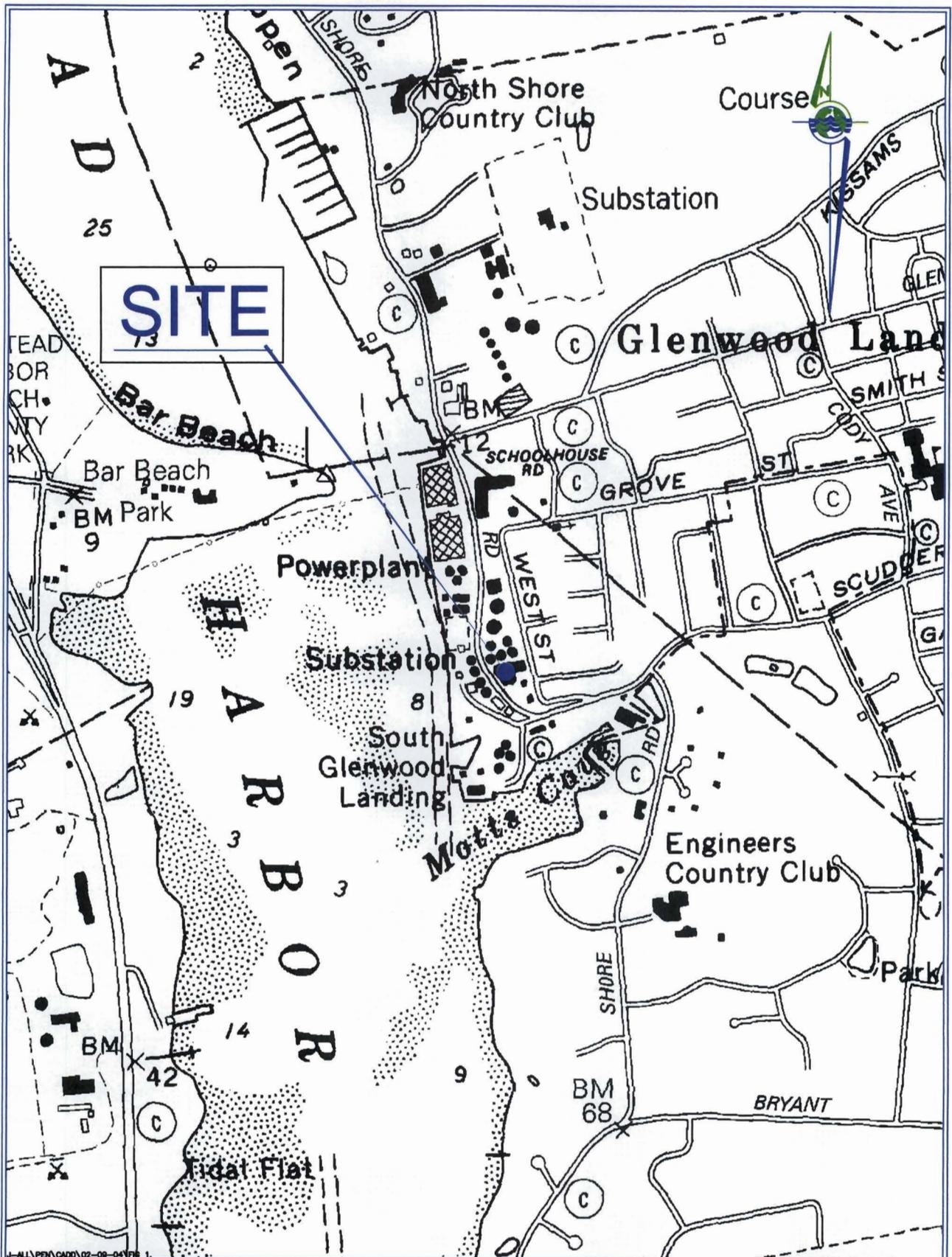
John Eichler  
Field Hydrogeologist



James Rhodes, CPG  
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Cc: G. Bobersky, NYSDEC  
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J. Nealon, NYSDOH  
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**Figures**



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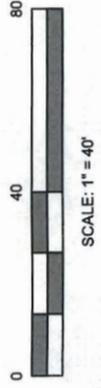
### Site Location Map

Project	PEN0001	Figure No.	1
Designed by	JPR		
Checked by	JPR		
Date			02/09/04



**LEGEND**

- ⊕ MONITORING WELL
- DRY WELL / LEACHING STRUCTURE



**PROPOSED SOIL BORING LOCATIONS**  
**FORMER PENETREX PROCESSING NYSDEC**  
 I.D. No. 130034

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Project:	PEN0001	Designed By:	JPR	Figure No.:	2
CAD Operator:	LLG	Approved By:	JPR	Date:	10/03/06

SOURCE: YEC, INC., SURVEY MAP 10, JULY 1992

**Tables**

**TABLE 1**

Soil Sample Analytical Results for Volatile Organic Compounds by EPA Method 8260  
One Shore Road, Glenwood Landing, New York

<b>Compound</b>	<b>NYSDEC Standard<sup>(1)</sup></b>	<b>DW-8</b>	<b>DW-10</b>	<b>Trip Blank</b>
Dichlorodifluoromethane	10,000*	< 6.2	< 6.4	< 1.0
1,1,1,2-Tetrachloroethane	10,000*	< 6.2	< 6.4	< 1.0
1,1,1-Trichloroethane	700	< 6.2	< 6.4	< 1.0
1,1,2,2-Tetrachloroethane	400	< 6.2	< 6.4	< 1.0
1,1,2-Trichloroethane	10,000*	< 6.2	< 6.4	< 1.0
1,1-Dichloroethane	300	< 6.2	< 6.4	< 1.0
1,1-Dichloroethene	300	< 6.2	< 6.4	< 1.0
1,1-Dichloropropene	10,000*	< 6.2	< 6.4	< 1.0
1,2,3-Trichlorobenzene	8,300	< 6.2	< 6.4	< 1.0
1,2,3-Trichloropropane	10,000*	< 6.2	< 6.4	< 1.0
1,2,4-Trichlorobenzene	8,300	< 6.2	< 6.4	< 1.0
1,2,4-Trimethylbenzene	NS	< 6.2	< 6.4	< 1.0
1,2-Dichlorobenzene	7,900	< 6.2	< 6.4	< 1.0
1,2-Dichloroethane	20 or MDL	< 6.2	< 6.4	< 1.0
1,2-Dichloropropane	10,000*	< 6.2	< 6.4	< 1.0
1,3,5-Trimethylbenzene	3,300	< 6.2	< 6.4	< 1.0
1,3-Dichlorobenzene	1,600	< 6.2	< 6.4	< 1.0
1,3-Dichloropropane	10,000*	< 6.2	< 6.4	< 1.0
1,4-Dichlorobenzene	8,500	< 6.2	< 6.4	< 1.0
2,2-Dichloropropane	10,000*	< 6.2	< 6.4	< 1.0
2-Chlorotoluene	10,000*	< 6.2	< 6.4	< 1.0
4-Chlorotoluene	10,000*	< 6.2	< 6.4	< 1.0
Benzene	60	< 6.2	< 6.4	< 1.0
Bromobenzene	10,000*	< 6.2	< 6.4	< 1.0
Bromochloromethane	10,000*	< 6.2	< 6.4	< 1.0
Bromodichloromethane	10,000*	< 6.2	< 6.4	< 1.0
Bromoform	10,000*	< 6.2	< 6.4	< 1.0
Bromomethane	10,000*	< 6.2	< 6.4	< 1.0
Carbon Tetrachloride	800	< 6.2	< 6.4	< 1.0
Chlorobenzene	1,100	< 6.2	< 6.4	< 1.0
Chloroethane	200	< 6.2	< 6.4	< 1.0
Chloroform	400	< 6.2	< 6.4	< 1.0
Chloromethane	10,000*	< 6.2	< 6.4	< 1.0
cis-1,2-Dichloroethene	200	7.7	67	< 1.0
cis-1,3-Dichloropropene	NS	< 6.2	< 6.4	< 1.0
Dibromochloromethane	NS	< 6.2	< 6.4	< 1.0
Dibromomethane	10,000*	< 6.2	< 6.4	< 1.0
Ethyl benzene	5,500	< 6.2	< 6.4	< 1.0
Hexachlorobutadiene	10,000*	< 6.2	< 6.4	< 1.0
Isopropylbenzene	2,300	< 6.2	< 6.4	< 1.0
m + p Xylene	1,200	< 12	< 13	< 2.0
Methyl tertiary butyl ether	1,200	< 6.2	< 6.4	< 1.0
Methylene Chloride	50 or MDL	37 B	26 B	16 B
n-Butylbenzene	NS	< 6.2	< 6.4	< 1.0

n-Propylbenzene	3,700	< 6.2	< 6.4	< 1.0
o Xylene	1,200	< 6.2	< 6.4	< 1.0
4-Isopropyltoluene	NS	< 6.2	< 6.4	< 1.0
sec-Butylbenzene	NS	< 6.2	< 6.4	< 1.0
Styrene	10,000*	< 6.2	< 6.4	< 1.0
tert-Butylbenzene	10,000*	< 6.2	< 6.4	< 1.0
Tetrachloroethene	1,300	52	240	< 1.0
Toluene	1,500	< 6.2	< 6.4	< 1.0
trans-1,2-Dichloroethene	200	< 6.2	< 6.4	< 1.0
trans-1,3-Dichloropropene	NS	< 6.2	< 6.4	< 1.0
Trichloroethene	500	11	34	< 1.0
Trichlorofluoromethane	10,000*	< 6.2	< 6.4	< 1.0
Vinyl chloride	200	< 6.2	< 6.4	< 1.0
2-propanol	NS	< 62	< 64	< 50
1,2-Dibromoethane	NS	< 6.2	< 6.4	< 1.0
2-Butanone	300	< 6.2	< 6.4	< 1.0
2-Chloroethyl vinyl ether	NS	< 6.2	< 6.4	< 1.0
2-Hexanone	NS	< 6.2	< 6.4	< 1.0
4-Methyl-2-pentanone	1,000	< 6.2	< 6.4	< 1.0
Acetone	200	< 6.2	< 6.4	< 1.0
Acrolein	NS	< 31	< 32	< 1.0
Acrylonitrile	NS	< 6.2	< 6.4	< 1.0
Carbon Disulfide	2,700	< 6.2	< 6.4	< 1.0
Chlorodifluoromethane	NS	< 6.2	< 6.4	< 1.0
Diisopropyl ether	NS	< 6.2	< 6.4	< 1.0
Ethanol	NS	< 31	< 32	< 1.0
Ethyl acetate	NS	< 6.2	< 6.4	< 1.0
Freon-114	NS	< 6.2	< 6.4	< 1.0
Isopropyl acetate	NS	< 6.2	< 6.4	< 1.0
n-Amyl acetate	NS	< 6.2	< 6.4	< 1.0
Naphthalene	NS	< 6.2	< 6.4	< 1.0
n-Butyl acetate	NS	< 6.2	< 6.4	< 1.0
n-Propyl acetate	NS	< 6.2	< 6.4	< 1.0
p-Diethylbenzene	NS	< 6.2	< 6.4	< 1.0
p-ethyltoluene	NS	< 6.2	< 6.4	< 1.0
t-Butyl alcohol	NS	< 6.2	< 6.4	< 1.0
Vinyl acetate	NS	< 6.2	< 6.4	< 1.0
Freon-113	6,000	< 6.2	< 6.4	< 1.0
1,2,4,5-tetramethylbenzene	NS	< 6.2	< 6.4	< 1.0
1,2-dibromo-3-chloropropane	NS	< 6.2	< 6.4	< 1.0

**Notes:**

1 - NYSDEC Recommended Soil Cleanup Objectives (RSCO), Technical and Administrative Guidance Memo (TAGM) 4046, 12/00.

NS - No RSCO

Bolded text denotes RSCO Exceedance.

\* - No specific RSCO established, RSCO of 10,000 ug/kg for total VOCs is used.

All units are ug/kg.

B - Compound detected in laboratory blank.

**TABLE 2**

Soil Sample Analytical Results for Semi-volatile Organic Compounds by EPA Method 8270  
One Shore Road, Glenwood Landing, New York

Compound	NYSDEC Standard <sup>(1)</sup>	DW-8 (8/24/06)	DW-10 (8/24/06)
Phenol	30 or MDL	< 140	< 150
Bis(2-chloroethyl)ether	50,000*	< 140	<150
1,3-Dichlorobenzene	1,000	< 140	<150
1,4-Dichlorobenzene	1,800	< 140	<150
1,2-Dichlorobenzene	1,100	< 140	<150
Benzyl alcohol	NS	< 140	<150
2-Methylphenol	100 or MDL	< 140	<150
N-Nitrosodi-n-propylamine	50,000*	< 140	<150
Hexachloroethane	50,000*	< 140	<150
4-Methylphenol	900	< 140	<150
2-Chlorophenol	800	< 140	<150
Nitrobenzene	50 or MDL	< 140	<150
Bis(2-chloroethoxy)methane	50,000*	< 140	<150
1,2,4-Trichlorobenzene	8,300	< 140	<150
Isophorone	1,000	< 140	<150
2,4-Dimethylphenol	NS	< 140	<150
Hexachlorobutadiene	50,000*	< 140	<150
Naphthalene	13,000	< 140	<150
2,4-Dichlorophenol	400	< 140	<150
4-Chloroaniline	3,100	< 140	<150
2,4,6-Trichlorophenol	NS	< 140	<150
2,4,5-Trichlorophenol	100	< 140	<150
Hexachlorocyclopentadiene	50,000*	< 140	<150
2-Methylnaphthalene	36,400	< 140	<150
2-Nitroaniline	430 or MDL	< 140	<150
2-Chloronaphthalene	50,000*	< 140	<150
4-Chloro-3-methylphenol	240 or MDL	< 140	<150
2,6-Dinitrotoluene	500	< 140	<150
2-Nitrophenol	330 or MDL	< 140	<150
3-Nitroaniline	500 or MDL	< 140	<150
Dimethylphthalate	2,300	< 140	<150
2,4-Dinitrophenol	200 or MDL	< 140	<150
Acenaphthylene	41,000	< 140	<150
2,4-Dinitrotoluene	50,000*	< 140	<150
Acenaphthene	50,000*	< 140	<150
Dibenzofuran	6,200	< 140	<150
4-Nitrophenol	100 or MDL	< 140	<150
Fluorene	50,000*	< 140	<150
4-Nitroaniline	NS	< 140	<150
4-Bromophenyl phenyl ether	50,000*	< 140	<150
Hexachlorobenzene	400	< 140	<150
Diethylphthalate	7,100	< 140	<150
4-Chlorophenyl phenyl ether	NS	< 140	<150
Pentachlorophenol	1,000 or MDL	< 140	<150
n-Nitrosodiphenylamine	NS	< 140	<150
4,6-Dintro-2-methylphenol	NS	< 140	<150
Phenanthrene	50,000*	280	340

Anthracene	50,000*	76 J	<150
Carbazole	NS	< 140	<150
Di-n-butylphthalate	50,000*	< 140	<150
Fluoroanthene	50,000*	390	500
Pyrene	50,000*	430	480
Butylbenzylphthalate	50,000*	< 140	<150
Benzo(a)anthracene	224 or MDL	180	190
Chrysene	400	200	230
3,3-Dichlorobenzidine	1,400	< 140	<150
Bis(2-ethylhexyl)phthalate	50,000*	< 140	<150
Di-n-octylphthalate	50,000*	< 140	<150
Benzo(b)fluoranthene	61 or MDL	<b>250</b>	<b>250</b>
Benzo(k)fluoranthene	610 or MDL	87 J	99 J
Benzo (a) pyrene	61 or MDL	<b>210</b>	<b>180</b>
Indeno(1,2,3-cd)pyrene	3,200	120 J	120 J
Dibenzo(a,h)anthracene	14.3 or MDL	< 140	<150
Benzo(g,h,i)perylene	50,000*	130 J	120 J
3+4 Methylphenol	20 or MDL	<140	<150
Aniline	100	<140	<150
Azobenzene	NS	<140	<150
Benzidene	NS	<140	<150
Benzoic Acid	NS	<140	<150
Bis(2-chloroisopropyl) ether	NS	<140	<150
N-Nitrosodimethylamine	NS	<140	<150
Pyridine	NS	<140	<150

**Notes:**

1 - NYSDEC Recommended Soil Cleanup Objectives (RSCO), Technical and Administrative Guidance Memo (TAGM) 4046, 12/00.

NS - No RSCO established for this compound.

Bolded text denotes RSCO Exceedance.

\* - No specific RSCO established, RSCO of 50,000 µg/kg for total SVOCs is used.

All units are µg/kg.

J - Result is less than the reporting limit, but greater than or equal to the method detection limit.

MDL - Method Detection Limit

**TABLE 3**

Soil Sample Analytical Results for Total Metals by EPA Method 6010  
One Shore Road, Glenwood Landing, New York

<b>Compound</b>	<b>NYSDEC Standard<sup>(1)</sup></b>	<b>Eastern USA Background<sup>(2)</sup></b>	<b>DW-8</b>	<b>DW-10</b>
Arsenic	7.5 or SB	3-12	3.42	1.06
Barium	300 or SB	15-600	35.9	7.32
Cadmium	10	0.1-1	0.651	0.643
Chromium	50	1.5-40	10.5	5.96
Lead	400	200-500	94.3	62.4
Mercury	0.1 or SB	0.001-0.2	0.174	0.0664
Selenium	2 or SB	0.1-3.9	< 0.592	0.664
Silver	SB	N/A	< 0.474	<0.497

**Notes:**

1 - NYSDEC Recommended Soil Cleanup Objectives, Technical and Administrative Guidance Memo (TAGM) 4046, 12/00.

All units are mg/kg.

SB - Site Background.

N/A - Not Available.

**TABLE 4**  
Soil Sample Analytical Results for Volatile Organic Compounds by EPA Method 8260  
One Shore Road, Glenwood Landing, New York

Compound	NYSDEC Standard	SB-8 13'-15' (DW-9)	Trip Blank 8/30/2006	Field Blank 8/30/2006
Dichlorodifluoromethane	10,000*	< 6.1	< 1.0	< 1.0
1,1,1,2-Tetrachloroethane	10,000*	< 6.1	< 1.0	< 1.0
1,1,1-Trichloroethane	700	< 6.1	< 1.0	< 1.0
1,1,2,2-Tetrachloroethane	400	< 6.1	< 1.0	< 1.0
1,1,2-Trichloroethane	10,000*	< 6.1	< 1.0	< 1.0
1,1-Dichloroethane	300	< 6.1	< 1.0	< 1.0
1,1-Dichloroethene	300	< 6.1	< 1.0	< 1.0
1,1-Dichloropropene	10,000*	< 6.1	< 1.0	< 1.0
1,2,3-Trichlorobenzene	8,300	< 6.1	< 1.0	< 1.0
1,2,3-Trichloropropane	10,000*	< 6.1	< 1.0	< 1.0
1,2,4-Trichlorobenzene	8,300	< 6.1	< 1.0	< 1.0
1,2,4-Trimethylbenzene	NS	< 6.1	< 1.0	< 1.0
1,2-Dichlorobenzene	7,900	< 6.1	< 1.0	< 1.0
1,2-Dichloroethane	20 or MDL	< 6.1	< 1.0	< 1.0
1,2-Dichloropropane	10,000*	< 6.1	< 1.0	< 1.0
1,3,5-Trimethylbenzene	3,300	< 6.1	< 1.0	< 1.0
1,3-Dichlorobenzene	1,600	< 6.1	< 1.0	< 1.0
1,3-Dichloropropane	10,000*	< 6.1	< 1.0	< 1.0
1,4-Dichlorobenzene	8,500	< 6.1	< 1.0	< 1.0
2,2-Dichloropropane	10,000*	< 6.1	< 1.0	< 1.0
2-Chlorotoluene	10,000*	< 6.1	< 1.0	< 1.0
4-Chlorotoluene	10,000*	< 6.1	< 1.0	< 1.0
Benzene	60	< 6.1	< 1.0	< 1.0
Bromobenzene	10,000*	< 6.1	< 1.0	< 1.0
Bromochloromethane	10,000*	< 6.1	< 1.0	< 1.0
Bromodichloromethane	10,000*	< 6.1	< 1.0	< 1.0
Bromoform	10,000*	< 6.1	< 1.0	< 1.0
Bromomethane	10,000*	< 6.1	< 1.0	< 1.0
Carbon Tetrachloride	800	< 6.1	< 1.0	< 1.0
Chlorobenzene	1,100	< 6.1	< 1.0	< 1.0
Chloroethane	200	< 6.1	< 1.0	< 1.0
Chloroform	400	< 6.1	< 1.0	< 1.0
Chloromethane	10,000*	< 6.1	< 1.0	< 1.0
cis-1,2-Dichloroethene	200	< 6.1	< 1.0	< 1.0
cis-1,3-Dichloropropene	NS	< 6.1	< 1.0	< 1.0
Dibromochloromethane	NS	< 6.1	< 1.0	< 1.0
Dibromomethane	10,000*	< 6.1	< 1.0	< 1.0
Ethyl benzene	5,500	< 6.1	< 1.0	< 1.0
Hexachlorobutadiene	10,000*	< 6.1	< 1.0	< 1.0
Isopropylbenzene	2,300	< 6.1	< 1.0	< 1.0
m + p Xylene	1,200	< 12	< 2.0	< 2.0
Methyl tertiary butyl ether	1,200	< 6.1	< 1.0	< 1.0
Methylene Chloride	50 or MDL	11 B	5.1 B	5.2 B

n-Butylbenzene	NS	< 6.1	< 1.0	< 1.0
n-Propylbenzene	3,700	< 6.1	< 1.0	< 1.0
o Xylene	1,200	< 6.1	< 1.0	< 1.0
4-Isopropyltoluene	NS	< 6.1	< 1.0	< 1.0
sec-Butylbenzene	NS	< 6.1	< 1.0	< 1.0
Styrene	10,000*	< 6.1	< 1.0	< 1.0
tert-Butylbenzene	10,000*	< 6.1	< 1.0	< 1.0
Tetrachloroethene	1,300	< 6.1	< 1.0	< 1.0
Toluene	1,500	26	< 1.0	< 1.0
trans-1,2-Dichloroethene	200	< 6.1	< 1.0	< 1.0
trans-1,3-Dichloropropene	NS	< 6.1	< 1.0	< 1.0
Trichloroethene	500	< 6.1	< 1.0	< 1.0
Trichlorofluoromethane	10,000*	< 6.1	< 1.0	< 1.0
Vinyl chloride	200	< 6.1	< 1.0	< 1.0
2-propanol	NS	< 61	< 50	< 50
1,2-Dibromoethane	NS	< 6.1	< 1.0	< 1.0
2-Butanone	300	< 6.1	< 1.0	< 1.0
2-Chloroethyl vinyl ether	NS	< 6.1	< 1.0	< 1.0
2-Hexanone	NS	< 6.1	< 1.0	< 1.0
4-Methyl-2-pentanone	1,000	< 6.1	< 1.0	< 1.0
Acetone	200	< 6.1	< 1.0	< 1.0
Acrolein	NS	< 31	< 1.0	< 1.0
Acrylonitrile	NS	< 6.1	< 1.0	< 1.0
Carbon Disulfide	2,700	< 6.1	< 1.0	< 1.0
Chlorodifluoromethane	NS	< 6.1	< 1.0	< 1.0
Diisopropyl ether	NS	< 6.1	< 1.0	< 1.0
Ethanol	NS	< 31	< 1.0	< 1.0
Ethyl acetate	NS	< 6.1	< 1.0	< 1.0
Freon-114	NS	< 6.1	< 1.0	< 1.0
Isopropyl acetate	NS	< 6.1	< 1.0	< 1.0
n-Amyl acetate	NS	< 6.1	< 1.0	< 1.0
Naphthalene	NS	< 6.1	< 1.0	< 1.0
n-Butyl acetate	NS	< 6.1	< 1.0	< 1.0
n-Propyl acetate	NS	< 6.1	< 1.0	< 1.0
p-Diethylbenzene	NS	< 6.1	< 1.0	< 1.0
p-ethyltoluene	NS	< 6.1	< 1.0	< 1.0
t-Butyl alcohol	NS	< 6.1	< 1.0	< 1.0
Vinyl acetate	NS	< 6.1	< 1.0	< 1.0
Freon-113	6,000	< 6.1	< 1.0	< 1.0
1,2,4,5-tetramethylbenzene	NS	< 6.1	< 1.0	< 1.0
1,2-dibromo-3-chloropropane	NS	< 6.1	< 1.0	< 1.0

**Notes:**

1 - NYSDEC Recommended Soil Cleanup Objectives (RSCO), Technical and Administrative Guidance Memo (TAGM) 4046, 12/00.

NS-No RSCO established for this compound.

\* - No specific RSCO established, RSCO of 10,000 ug/kg for total VOCs is used.

All units are µg/kg.

B - Compound detected in laboratory blank.

**TABLE 5**  
 Soil Sample Analytical Results  
 Semi-volatile Organic Compounds by EPA Method 8270  
 One Shore Road, Glenwood Landing, New York

Compound	NYSDEC Standard <sup>(1)</sup>	SB-8, 13'-15' (DW-9)
Phenol	30 or MDL	< 150
Bis(2-chloroethyl)ether	50,000*	< 150
1,3-Dichlorobenzene	1,000	< 150
1,4-Dichlorobenzene	1,800	< 150
1,2-Dichlorobenzene	1,100	< 150
Benzyl alcohol	NS	< 150
2-Methylphenol	100 or MDL	< 150
N-Nitrosodi-n-propylamine	50,000*	< 150
Hexachloroethane	50,000*	< 150
4-Methylphenol	900	< 150
2-Chlorophenol	800	< 150
Nitrobenzene	50 or MDL	< 150
Bis(2-chloroethoxy)methane	50,000*	< 150
1,2,4-Trichlorobenzene	8,300	< 150
Isophorone	1,000	< 150
2,4-Dimethylphenol	NS	< 150
Hexachlorobutadiene	50,000*	< 150
Naphthalene	13,000	< 150
2,4-Dichlorophenol	400	< 150
4-Chloroaniline	3,100	< 150
2,4,6-Trichlorophenol	NS	< 150
2,4,5-Trichlorophenol	100	< 150
Hexachlorocyclopentadiene	50,000*	< 150
2-Methylnaphthalene	36,400	< 150
2-Nitroaniline	430 or MDL	< 150
2-Chloronaphthalene	50,000*	< 150
4-Chloro-3-methylphenol	240 or MDL	< 150
2,6-Dinitrotoluene	500	< 150
2-Nitrophenol	330 or MDL	< 150
3-Nitroaniline	500 or MDL	< 150
Dimethylphthalate	2,300	< 150
2,4-Dinitrophenol	200 or MDL	< 150
Acenaphthylene	41,000	< 150
2,4-Dinitrotoluene	50,000*	< 150
Acenaphthene	50,000*	< 150
Dibenzofuran	6,200	< 150
4-Nitrophenol	100 or MDL	< 150
Fluorene	50,000*	< 150
4-Nitroaniline	NS	< 150
4-Bromophenyl phenyl ether	50,000*	< 150
Hexachlorobenzene	400	< 150
Diethylphthalate	7,100	< 150

4-Chlorophenyl phenyl ether	NS	< 150
Pentachlorophenol	1,000 or MDL	< 150
n-Nitrosodiphenylamine	NS	< 150
4,6-Dintro-2-methylphenol	NS	< 150
Phenanthrene	50,000*	< 150
Anthracene	50,000*	< 150
Carbazole	NS	< 150
Di-n-butylphthalate	50,000*	< 150
Fluoroanthene	50,000*	< 150
Pyrene	50,000*	< 150
Butylbenzylphthalate	50,000*	< 150
Benzo(a)anthracene	224 or MDL	< 150
Chrysene	400	< 150
3,3-Dichlorobenzidine	1,400	< 150
Bis(2-ethylhexyl)phthalate	50,000*	< 150
Di-n-octylphthalate	50,000*	< 150
Benzo(b)fluoranthene	61 or MDL	< 150
Benzo(k)fluoranthene	610 or MDL	< 150
Benzo (a) pyrene	61 or MDL	< 150
Indeno(1,2,3-cd)pyrene	3,200	< 150
Dibenzo(a,h)anthracene	14.3 or MDL	< 150
Benzo(g,h,i)perylene	50,000*	< 150
3+4 Methylphenol	20 or MDL	< 150
Aniline	100	< 150
Azobenzene	NS	< 150
Benzidene	NS	< 150
Benzoic Acid	NS	< 150
Bis(2-chloroisopropyl) ether	NS	< 150
N-Nitrosodimethylamine	NS	< 150
Pyridine	NS	< 150

**Notes:**

1 - NYSDEC Recommended Soil Cleanup Objectives (RSCO), Technical and Administrative Guidance Memo (TAGM) 4046, 12/00.

NS - No RSCO established for this compound.

\* - No specific RSCO established, RSCO of 50,000 µg/kg for total SVOCs is used.

All units are µg/kg.

MDL - Method Detection Limit.

**TABLE 6**

Soil Sample Analytical Results for Total Metals by EPA Method 6010  
One Shore Road, Glenwood Landing, New York

<b>Compound</b>	<b>NYSDEC RSCO (1)</b>	<b>Eastern USA Background</b>	<b>SB-8, 13'-15' (DW-9)</b>
Arsenic	7.5 or SB	.3-12	0.601
Barium	300 or SB	15-600	3.06
Cadmium	10	0.1-1	0.835
Chromium	50	1.5-40	1.13
Lead	400	200-500	38.6
Mercury	0.1 or SB	0.001-0.2	< 0.0116
Selenium	2 or SB	0.1-3.9	< 0.601
Silver	SB	N/A	< 0.481

**Notes:**

1 - NYSDEC Recommended Soil Cleanup Objectives, Technical and Administrative  
Guidance Memo (TAGM) 4046, 12/00.

All units are mg/kg.

SB - Site Background.

**Appendix A**  
**USEPA Closure Letter**



ENVIRONMENTAL PROTECTION AGENCY  
 REGION 2  
 290 BROADWAY  
 NEW YORK, NY 10007-1866

NOV 12 2003

**CERTIFIED MAIL - RETURN RECEIPT REQUESTED**

**Article Number: 7002 2030 0006 5360 7344**

**Property Owner:**

Larry Weinberger  
 Glenwood Realty, LLC  
 P.O. Box 1356  
 Roslyn Heights, NY 11577

**Facility:**

Former Penetrex Site  
 1 Shore Road  
 Glenwood Landing, NY 11547  
 NYSDEC Site # 1-30-034

Re: Class V Underground Injection Wells and File Closure  
 Class V Underground Injection Wells and Authorization by Rule

Dear Mr. Weinberger:

The Ground Water Compliance Section of the U.S. Environmental Protection Agency (EPA) has reviewed the Storm Drain and Sanitary Leaching Pool Remediation and Closure Report prepared by P.W. Grosser Consulting, Inc. (dated September 9, 2003), concerning the closure and removal from service of drywells DW-2 and DW-3. DW-2 and DW-3 have been closed and sealed to the satisfaction of this office, and the EPA will now close its Underground Injection Control (UIC) file on drywells DW-2 and DW-3.

The Storm Drain and Sanitary Leaching Pool Remediation and Closure Report also concerned the remediations of five Class V injection wells. These were the remediation of the western sanitary septic system (seepage pit, A-1, and the distribution box), the active stormwater drainage well DW-1, the inactive storm water drainage well DW-4 converted to a residential secondary overflow leaching pool, and the active cesspool DW-5. These injection wells have been remediated to the satisfaction of this office.

The continued use of these as Class V wells, in accordance with applicable EPA regulations, is approved. The injection wells are considered active Class V UIC wells. There may be additional regulatory requirements when future regulations for Class V wells become promulgated. The Class V injection wells are authorized by rule pursuant to Title 40 of the Code of Federal Regulations (C.F.R.), §144.24.

In accordance with 40 C.F.R., § § 144.84(b)(2), 144.85, and 144.88 (Table), existing large capacity cesspools must be closed by April 5, 2005. You must notify the UIC program director (below) of your intent to close the cesspool at least 30 days prior to closure.

Internet Address (URL) • <http://www.epa.gov>

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Should any conditions change (such as return of the residential well DW-4 to a commercial well, injectate composition, accidental spills into the system, sealing the drywells, injection of cooling water more than 90° Fahrenheit, construction of additional wells, etc.), you are **required** to notify this office, specifically:

**Charles Hillenbrand, Chief  
Ground Water Compliance Section, 2DECA-WCB, 20<sup>th</sup> Floor  
U.S. Environmental Protection Agency  
290 Broadway New York, NY 10007-1866**

If you have any questions, you may contact Margaret Halley of my staff at (212) 637-3092.

Sincerely,



**Charles Hillenbrand, Ph.D., Chief  
Ground Water Compliance Section**

cc: Bill Spitz, NYSDEC, Region 1

Walter Parish, NYSDEC, Region 1

John Lovejoy, NCHD-Bureau of Water Supply Protection

Joe Jones  
NYSDEC Central Office  
625 Broadway  
Albany NY 12207-2942

Dave Yudelson, Esq.  
Sive, Paget & Riesel PC  
460 Park Avenue  
New York, NY 10022

Bruce Mackay  
P. W. Grosser Consulting, Inc.  
Suite 7  
630 Johnson Avenue  
Bohemia NY 11716-2618

**Appendix B**  
**Laboratory Data Sheets**



NYSDOH 11418  
NJDEP NY050  
CTDOH PH-0205  
PADEP 68-00573

Friday, September 01, 2006

Jim Rhodes  
P.W. Grosser Consulting  
630 Johnson Avenue  
Suite 7  
Bohemia, NY 11716

TEL: (631) 589-6353  
FAX (631) 589-8705

RE: PEN-Glenwood Landing

Order No.: 0608273

Dear Jim Rhodes:

American Analytical Laboratories, LLC. received 5 sample(s) on 8/24/2006 for the analyses presented in the following report.

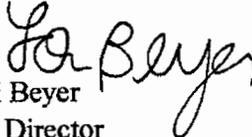
Samples were analyzed in accordance with the test procedures documented on the chain of custody and detailed throughout the text of this report.

The limits provided in the data package are analytical reporting limits and not Federal or Local mandated values to which the sample results should be compared.

There were no problems with the analyses and all data for associated QC met laboratory specifications. If there are any exceptions a Case Narrative is provided in the report.

If you have any questions regarding these tests results, please do not hesitate to call (631) 454-6100 or email me directly at [lbeyer@american-analytical.com](mailto:lbeyer@american-analytical.com).

Sincerely,

  
Lori Beyer  
Lab Director

---

**CLIENT:** P.W. Grosser Consulting  
**Project:** PEN-Glenwood Landing  
**Lab Order:** 0608273

**Work Order Sample Summary**

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<b>Lab Sample ID</b>	<b>Client Sample ID</b>	<b>Tag Number</b>	<b>Date Collected</b>	<b>Date Received</b>
0608273-01A	DW-8	10005	8/24/2006 11:25:00 AM	8/24/2006
0608273-02A	DW-10	10005	8/24/2006 12:50:00 PM	8/24/2006
0608273-03A	MDA-10	10005	8/24/2006 10:30:00 AM	8/24/2006
0608273-04A	Trip Blank	10005	8/24/2006	8/24/2006
0608273-05A	Field Blank	10005	8/24/2006 11:10:00 AM	8/24/2006



**AMERICAN ANALYTICAL LABORATORIES, LLC**

**56 TOLEDO STREET**

**FARMINGDALE, NEW YORK 11735**

**TELEPHONE: (631) 454-6100 FAX: (631) 454-8027**

**DATA REPORTING QUALIFIERS**

For reporting results, the following "Results Qualifiers" are used:

<b>Value</b>	If the result is greater than or equal to the detection limit, report the value
<b>U</b>	Indicates the compound was analyzed for but was not detected. Report the minimum detection limit for the sample with the U, i.e. "10U". This is not necessarily the instrument detection limit attainable for this particular sample based on any concentration or dilution that may have been required.
<b>J</b>	Indicates an estimated value. The flag is used: <ol style="list-style-type: none"><li>(1) When estimating a concentration for a tentatively identified compound (library search hits, where a 1:1 response is assumed.)</li><li>(2) When the mass spectral data indicated the identification, however the result was less than the specified detection limit greater than zero. If the detection limit was 10ug/L and a concentration of 3ug/L was calculated report as 3J. This flag is used when similar situations arise on any organic parameter i.e. Pesticide, PCBs and others.</li></ol>
<b>B</b>	Indicates the analyte was found in the blank as well as the sample report "10B".
<b>E</b>	Indicates the analytes concentration exceeds the calibrated range of the instrument for that specific analysis.
<b>D</b>	This flag identifies all compounds identified in an analysis at a secondary dilution factor.
<b>P</b>	This flag is used for Pesticide / PCB target analyte when there is >25% difference for detected concentrations between the two GC Columns. The higher of the two values is reported on Form I and flagged with a "P".
<b>N</b>	This flag indicates presumptive evidence of a compound. This is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It applies to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the flag is not used.
<b>H</b>	Indicates sample was received and/or analyzed outside of The method allowable holding time

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	DW-8
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 11:25:00 AM
<b>Lab ID:</b>	0608273-01A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>MERCURY</b>						
		<b>SW7471B</b>				<b>Analyst: WN</b>
Mercury	0.174	0.0114		mg/Kg-dry	1	8/29/2006 3:34:17 PM
<b>PERCENT MOISTURE</b>						
		<b>D2216</b>				<b>Analyst: PA</b>
Percent Moisture	16.9	0		wt%	1	8/28/2006
<b>NCDH METALS</b>						
		<b>SW6010B</b>				<b>Analyst: JP</b>
Arsenic	3.42	0.592		mg/Kg-dry	1	8/29/2006 10:53:25 AM
Barium	35.9	0.474		mg/Kg-dry	1	8/29/2006 10:53:25 AM
Cadmium	0.651	0.237		mg/Kg-dry	1	8/29/2006 10:53:25 AM
Chromium	10.5	0.474		mg/Kg-dry	1	8/29/2006 10:53:25 AM
Lead	94.3	0.355		mg/Kg-dry	1	8/29/2006 10:53:25 AM
Selenium	U	0.592		mg/Kg-dry	1	8/29/2006 10:53:25 AM
Silver	U	0.474		mg/Kg-dry	1	8/29/2006 10:53:25 AM
<b>SEMIVOLATILE SW-846 METHOD 8270</b>						
		<b>SW8270D</b>		<b>SW3550A</b>		<b>Analyst: RN</b>
1,2,4-Trichlorobenzene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
1,2-Dichlorobenzene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
1,3-Dichlorobenzene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
1,4-Dichlorobenzene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2,4,5-Trichlorophenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2,4,6-Trichlorophenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2,4-Dichlorophenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2,4-Dimethylphenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2,4-Dinitrophenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2,4-Dinitrotoluene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2,6-Dinitrotoluene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2-Chloronaphthalene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2-Chlorophenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2-Methylnaphthalene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2-Methylphenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2-Nitroaniline	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
2-Nitrophenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
3,3'-Dichlorobenzidine	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
3+4-Methylphenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
3-Nitroaniline	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
4,6-Dinitro-2-methylphenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
4-Bromophenyl phenyl ether	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
4-Chloro-3-methylphenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
4-Chloroaniline	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
4-Chlorophenyl phenyl ether	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
4-Nitroaniline	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	DW-8
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 11:25:00 AM
<b>Lab ID:</b>	0608273-01A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>SEMIVOLATILE SW-846 METHOD 8270</b>		<b>SW8270D</b>		<b>SW3550A</b>		<b>Analyst: RN</b>
4-Nitrophenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Acenaphthene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Acenaphthylene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Aniline	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Anthracene	76	140	J	µg/Kg-dry	1	8/29/2006 11:52:00 PM
Azobenzene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Benzdine	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Benzo(a)anthracene	180	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Benzo(a)pyrene	210	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Benzo(b)fluoranthene	250	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Benzo(g,h,i)perylene	130	140	J	µg/Kg-dry	1	8/29/2006 11:52:00 PM
Benzo(k)fluoranthene	87	140	J	µg/Kg-dry	1	8/29/2006 11:52:00 PM
Benzoic acid	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Benzyl alcohol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Bis(2-chloroethoxy)methane	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Bis(2-chloroethyl)ether	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Bis(2-chloroisopropyl)ether	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Bis(2-ethylhexyl)phthalate	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Butyl benzyl phthalate	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Carbazole	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Chrysene	200	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Dibenzo(a,h)anthracene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Dibenzofuran	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Diethyl phthalate	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Dimethyl phthalate	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Di-n-butyl phthalate	110	140	J	µg/Kg-dry	1	8/29/2006 11:52:00 PM
Di-n-octyl phthalate	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Fluoranthene	390	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Fluorene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Hexachlorobenzene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Hexachlorobutadiene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Hexachlorocyclopentadiene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Hexachloroethane	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Indeno(1,2,3-c,d)pyrene	120	140	J	µg/Kg-dry	1	8/29/2006 11:52:00 PM
Isophorone	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Naphthalene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Nitrobenzene	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
N-Nitrosodimethylamine	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
N-Nitrosodi-n-propylamine	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	DW-8
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 11:25:00 AM
<b>Lab ID:</b>	0608273-01A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>SEMIVOLATILE SW-846 METHOD 8270</b>		<b>SW8270D</b>		<b>SW3550A</b>		Analyst: RN
N-Nitrosodiphenylamine	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Pentachlorophenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Phenanthrene	280	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Phenol	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Pyrene	430	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
Pyridine	U	140		µg/Kg-dry	1	8/29/2006 11:52:00 PM
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: LDS
1,1,1,2-Tetrachloroethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,1,1-Trichloroethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,1,2,2-Tetrachloroethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,1,2-Trichloro-1,2,2-trifluoroethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,1,2-Trichloroethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,1-Dichloroethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,1-Dichloroethene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,1-Dichloropropene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2,3-Trichlorobenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2,3-Trichloropropane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2,4,5-Tetramethylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2,4-Trichlorobenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2,4-Trimethylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2-Dibromo-3-chloropropane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2-Dibromoethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2-Dichlorobenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2-Dichloroethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,2-Dichloropropane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,3,5-Trimethylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,3-Dichlorobenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,3-dichloropropane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
1,4-Dichlorobenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
2,2-Dichloropropane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
2-Butanone	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
2-Chloroethyl vinyl ether	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
2-Chlorotoluene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
2-Hexanone	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
2-Propanol	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
4-Chlorotoluene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
4-Isopropyltoluene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
4-Methyl-2-pentanone	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Acetone	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

**CLIENT:** P.W. Grosser Consulting  
**Lab Order:** 0608273  
**Project:** PEN-Glenwood Landing  
**Lab ID:** 0608273-01A

**Client Sample ID:** DW-8  
**Tag Number:** 10005  
**Collection Date:** 8/24/2006 11:25:00 AM  
**Matrix:** SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>	<b>SW5030A</b>			Analyst: LDS
Acrolein	U	31		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Acrylonitrile	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Benzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Bromobenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Bromochloromethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Bromodichloromethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Bromoform	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Bromomethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Carbon disulfide	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Carbon tetrachloride	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Chlorobenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Chlorodifluoromethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Chloroethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Chloroform	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Chloromethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
cis-1,2-Dichloroethene	7.7	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
cis-1,3-Dichloropropene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Dibromochloromethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Dibromomethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Dichlorodifluoromethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Diisopropyl ether	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Ethanol	U	31		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Ethyl acetate	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Ethylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Freon-114	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Hexachlorobutadiene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Isopropyl acetate	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Isopropylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
m,p-Xylene	U	12		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Methyl tert-butyl ether	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Methylene chloride	37	6.2	B	µg/Kg-dry	1	8/29/2006 8:09:00 AM
n-Amyl acetate	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Naphthalene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
n-Butyl acetate	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
n-Butylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
n-Propyl acetate	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
n-Propylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
o-Xylene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
p-Diethylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM

**Qualifiers:** B Analyte detected in the associated Method Blank  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
U Indicates the compound was analyzed for but not detected

E Value above quantitation range  
J Analyte detected below quantitation limits  
S Spike Recovery outside accepted recovery limits  
X Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	DW-8
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 11:25:00 AM
<b>Lab ID:</b>	0608273-01A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
p-Ethyltoluene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
sec-Butylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Styrene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
t-Butyl alcohol	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
tert-Butylbenzene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Tetrachloroethene	52	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Toluene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
trans-1,2-Dichloroethene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
trans-1,3-Dichloropropene	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Trichloroethene	11	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Trichlorofluoromethane	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Vinyl acetate	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM
Vinyl chloride	U	6.2		µg/Kg-dry	1	8/29/2006 8:09:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	DW-10
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 12:50:00 PM
<b>Lab ID:</b>	0608273-02A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>MERCURY</b>		<b>SW7471B</b>		<b>SW7471B</b>		Analyst: <b>WN</b>
Mercury	0.0664	0.0112		mg/Kg-dry	1	8/29/2006 3:36:25 PM
<b>PERCENT MOISTURE</b>		<b>D2216</b>				Analyst: <b>PA</b>
Percent Moisture	20.7	0		wt%	1	8/28/2006
<b>NCDH METALS</b>		<b>SW6010B</b>		<b>SW3050A</b>		Analyst: <b>JP</b>
Arsenic	1.06	0.621		mg/Kg-dry	1	8/29/2006 10:56:07 AM
Barium	7.32	0.497		mg/Kg-dry	1	8/29/2006 10:56:07 AM
Cadmium	0.643	0.248		mg/Kg-dry	1	8/29/2006 10:56:07 AM
Chromium	5.96	0.497		mg/Kg-dry	1	8/29/2006 10:56:07 AM
Lead	62.4	0.373		mg/Kg-dry	1	8/29/2006 10:56:07 AM
Selenium	0.664	0.621		mg/Kg-dry	1	8/29/2006 10:56:07 AM
Silver	U	0.497		mg/Kg-dry	1	8/29/2006 10:56:07 AM
<b>SEMIVOLATILE SW-846 METHOD 8270</b>		<b>SW8270D</b>		<b>SW3550A</b>		Analyst: <b>RN</b>
1,2,4-Trichlorobenzene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
1,2-Dichlorobenzene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
1,3-Dichlorobenzene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
1,4-Dichlorobenzene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2,4,5-Trichlorophenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2,4,6-Trichlorophenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2,4-Dichlorophenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2,4-Dimethylphenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2,4-Dinitrophenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2,4-Dinitrotoluene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2,6-Dinitrotoluene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2-Chloronaphthalene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2-Chlorophenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2-Methylnaphthalene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2-Methylphenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2-Nitroaniline	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
2-Nitrophenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
3,3'-Dichlorobenzidine	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
3+4-Methylphenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
3-Nitroaniline	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
4,6-Dinitro-2-methylphenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
4-Bromophenyl phenyl ether	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
4-Chloro-3-methylphenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
4-Chloroaniline	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
4-Chlorophenyl phenyl ether	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
4-Nitroaniline	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

**CLIENT:** P.W. Grosser Consulting  
**Lab Order:** 0608273  
**Project:** PEN-Glenwood Landing  
**Lab ID:** 0608273-02A

**Client Sample ID:** DW-10  
**Tag Number:** 10005  
**Collection Date:** 8/24/2006 12:50:00 PM  
**Matrix:** SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>SEMIVOLATILE SW-846 METHOD 8270</b>		<b>SW8270D</b>		<b>SW3550A</b>		<b>Analyst: RN</b>
4-Nitrophenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Acenaphthene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Acenaphthylene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Aniline	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Anthracene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Azobenzene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Benzidine	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Benzo(a)anthracene	190	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Benzo(a)pyrene	180	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Benzo(b)fluoranthene	250	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Benzo(g,h,i)perylene	120	150	J	µg/Kg-dry	1	8/30/2006 12:17:00 AM
Benzo(k)fluoranthene	99	150	J	µg/Kg-dry	1	8/30/2006 12:17:00 AM
Benzoic acid	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Benzyl alcohol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Bis(2-chloroethoxy)methane	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Bis(2-chloroethyl)ether	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Bis(2-chloroisopropyl)ether	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Bis(2-ethylhexyl)phthalate	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Butyl benzyl phthalate	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Carbazole	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Chrysene	230	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Dibenzo(a,h)anthracene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Dibenzofuran	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Diethyl phthalate	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Dimethyl phthalate	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Di-n-butyl phthalate	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Di-n-octyl phthalate	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Fluoranthene	500	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Fluorene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Hexachlorobenzene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Hexachlorobutadiene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Hexachlorocyclopentadiene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Hexachloroethane	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Indeno(1,2,3-c,d)pyrene	120	150	J	µg/Kg-dry	1	8/30/2006 12:17:00 AM
Isophorone	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Naphthalene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Nitrobenzene	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
N-Nitrosodimethylamine	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
N-Nitrosodi-n-propylamine	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P. W. Grosser Consulting	<b>Client Sample ID:</b>	DW-10
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 12:50:00 PM
<b>Lab ID:</b>	0608273-02A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>SEMIVOLATILE SW-846 METHOD 8270</b>		<b>SW8270D</b>		<b>SW3550A</b>		<b>Analyst: RN</b>
N-Nitrosodiphenylamine	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Pentachlorophenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Phenanthrene	340	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Phenol	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Pyrene	480	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
Pyridine	U	150		µg/Kg-dry	1	8/30/2006 12:17:00 AM
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
1,1,1,2-Tetrachloroethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,1,1-Trichloroethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,1,2,2-Tetrachloroethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,1,2-Trichloro-1,2,2-trifluoroethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,1,2-Trichloroethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,1-Dichloroethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,1-Dichloroethene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,1-Dichloropropene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2,3-Trichlorobenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2,3-Trichloropropane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2,4,5-Tetramethylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2,4-Trichlorobenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2,4-Trimethylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2-Dibromo-3-chloropropane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2-Dibromoethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2-Dichlorobenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2-Dichloroethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,2-Dichloropropane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,3,5-Trimethylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,3-Dichlorobenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,3-dichloropropane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
1,4-Dichlorobenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
2,2-Dichloropropane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
2-Butanone	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
2-Chloroethyl vinyl ether	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
2-Chlorotoluene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
2-Hexanone	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
2-Propanol	U	64		µg/Kg-dry	1	8/29/2006 8:46:00 AM
4-Chlorotoluene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
4-Isopropyltoluene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
4-Methyl-2-pentanone	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Acetone	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	DW-10
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 12:50:00 PM
<b>Lab ID:</b>	0608273-02A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
Acrolein	U	32		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Acrylonitrile	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Benzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Bromobenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Bromochloromethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Bromodichloromethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Bromofom	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Bromomethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Carbon disulfide	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Carbon tetrachloride	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Chlorobenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Chlorodifluoromethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Chloroethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Chloroform	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Chloromethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
cis-1,2-Dichloroethene	67	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
cis-1,3-Dichloropropene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Dibromochloromethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Dibromomethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Dichlorodifluoromethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Diisopropyl ether	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Ethanol	U	32		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Ethyl acetate	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Ethylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Freon-114	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Hexachlorobutadiene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Isopropyl acetate	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Isopropylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
m,p-Xylene	U	13		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Methyl tert-butyl ether	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Methylene chloride	26	6.4	B	µg/Kg-dry	1	8/29/2006 8:46:00 AM
n-Amyl acetate	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Naphthalene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
n-Butyl acetate	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
n-Butylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
n-Propyl acetate	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
n-Propylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
o-Xylene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
p-Diethylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	DW-10
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 12:50:00 PM
<b>Lab ID:</b>	0608273-02A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
p-Ethyltoluene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
sec-Butylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Styrene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
t-Butyl alcohol	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
tert-Butylbenzene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Tetrachloroethene	240	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Toluene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
trans-1,2-Dichloroethene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
trans-1,3-Dichloropropene	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Trichloroethene	34	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Trichlorofluoromethane	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Vinyl acetate	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM
Vinyl chloride	U	6.4		µg/Kg-dry	1	8/29/2006 8:46:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	MDA-10
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 10:30:00 AM
<b>Lab ID:</b>	0608273-03A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>PERCENT MOISTURE</b>		<b>D2216</b>				Analyst: PA
Percent Moisture	20.2	0		wt%	1	8/28/2006
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: LDS
1,1,1,2-Tetrachloroethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,1,1-Trichloroethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,1,2,2-Tetrachloroethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,1,2-Trichloro-1,2,2-trifluoroethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,1,2-Trichloroethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,1-Dichloroethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,1-Dichloroethene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,1-Dichloropropene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2,3-Trichlorobenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2,3-Trichloropropane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2,4,5-Tetramethylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2,4-Trichlorobenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2,4-Trimethylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2-Dibromo-3-chloropropane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2-Dibromoethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2-Dichlorobenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2-Dichloroethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,2-Dichloropropane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,3,5-Trimethylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,3-Dichlorobenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,3-dichloropropane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
1,4-Dichlorobenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
2,2-Dichloropropane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
2-Butanone	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
2-Chloroethyl vinyl ether	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
2-Chlorotoluene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
2-Hexanone	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
2-Propanol	U	65		µg/Kg-dry	1	8/29/2006 9:23:00 AM
4-Chlorotoluene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
4-Isopropyltoluene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
4-Methyl-2-pentanone	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Acetone	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Acrolein	U	32		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Acrylonitrile	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Benzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Bromobenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Bromochloromethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

**CLIENT:** P.W. Grosser Consulting  
**Lab Order:** 0608273  
**Project:** PEN-Glenwood Landing  
**Lab ID:** 0608273-03A

**Client Sample ID:** MDA-10  
**Tag Number:** 10005  
**Collection Date:** 8/24/2006 10:30:00 AM  
**Matrix:** SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
Bromodichloromethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Bromoform	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Bromomethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Carbon disulfide	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Carbon tetrachloride	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Chlorobenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Chlorodifluoromethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Chloroethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Chloroform	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Chloromethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
cis-1,2-Dichloroethene	920	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
cis-1,3-Dichloropropene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Dibromochloromethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Dibromomethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Dichlorodifluoromethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Diisopropyl ether	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Ethanol	U	32		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Ethyl acetate	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Ethylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Freon-114	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Hexachlorobutadiene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Isopropyl acetate	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Isopropylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
m,p-Xylene	U	13		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Methyl tert-butyl ether	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Methylene chloride	21	6.5	B	µg/Kg-dry	1	8/29/2006 9:23:00 AM
n-Amyl acetate	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Naphthalene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
n-Butyl acetate	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
n-Butylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
n-Propyl acetate	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
n-Propylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
o-Xylene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
p-Diethylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
p-Ethyltoluene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
sec-Butylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Styrene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
t-Butyl alcohol	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
tert-Butylbenzene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	MDA-10
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 10:30:00 AM
<b>Lab ID:</b>	0608273-03A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>						
		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: LDS
Tetrachloroethene	410	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Toluene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
trans-1,2-Dichloroethene	55	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
trans-1,3-Dichloropropene	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Trichloroethene	56	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Trichlorofluoromethane	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Vinyl acetate	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM
Vinyl chloride	U	6.5		µg/Kg-dry	1	8/29/2006 9:23:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Trip Blank
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006
<b>Lab ID:</b>	0608273-04A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
1,1,1,2-Tetrachloroethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,1,1-Trichloroethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,1,2,2-Tetrachloroethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,1,2-Trichloro-1,2,2-trifluoroethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,1,2-Trichloroethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,1-Dichloroethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,1-Dichloroethene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,1-Dichloropropene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2,3-Trichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2,3-Trichloropropane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2,4,5-Tetramethylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2,4-Trichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2,4-Trimethylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2-Dibromo-3-chloropropane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2-Dibromoethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2-Dichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2-Dichloroethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,2-Dichloropropane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,3,5-Trimethylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,3-Dichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,3-dichloropropane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
1,4-Dichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
2,2-Dichloropropane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
2-Butanone	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
2-Chloroethyl vinyl ether	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
2-Chlorotoluene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
2-Hexanone	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
2-Propanol	U	50		µg/L	1	8/29/2006 6:16:00 AM
4-Chlorotoluene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
4-Isopropyltoluene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
4-Methyl-2-pentanone	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Acetone	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Acrolein	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Acrylonitrile	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Benzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Bromobenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Bromochloromethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Bromodichloromethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Bromoform	U	1.0		µg/L	1	8/29/2006 6:16:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

**CLIENT:** P.W. Grosser Consulting  
**Lab Order:** 0608273  
**Project:** PEN-Glenwood Landing  
**Lab ID:** 0608273-04A

**Client Sample ID:** Trip Blank  
**Tag Number:** 10005  
**Collection Date:** 8/24/2006  
**Matrix:** LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
Bromomethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Carbon disulfide	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Carbon tetrachloride	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Chlorobenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Chlorodifluoromethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Chloroethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Chloroform	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Chloromethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
cis-1,2-Dichloroethene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
cis-1,3-Dichloropropene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Dibromochloromethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Dibromomethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Dichlorodifluoromethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Diisopropyl ether	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Ethanol	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Ethyl acetate	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Ethylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Freon-114	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Hexachlorobutadiene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Isopropyl acetate	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Isopropylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
m,p-Xylene	U	2.0		µg/L	1	8/29/2006 6:16:00 AM
Methyl tert-butyl ether	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Methylene chloride	16	1.0	B	µg/L	1	8/29/2006 6:16:00 AM
n-Amyl acetate	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Naphthalene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
n-Butyl acetate	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
n-Butylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
n-Propyl acetate	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
n-Propylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
o-Xylene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
p-Diethylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
p-Ethyltoluene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
sec-Butylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Styrene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
t-Butyl alcohol	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
tert-Butylbenzene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Tetrachloroethene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Toluene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM

**Qualifiers:**  
 B Analyte detected in the associated Method Blank  
 H Holding times for preparation or analysis exceeded  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

E Value above quantitation range  
 J Analyte detected below quantitation limits  
 S Spike Recovery outside accepted recovery limits  
 X Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Trip Blank
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006
<b>Lab ID:</b>	0608273-04A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>						
		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: LDS
trans-1,2-Dichloroethene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
trans-1,3-Dichloropropene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Trichloroethene	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Trichlorofluoromethane	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Vinyl acetate	U	1.0		µg/L	1	8/29/2006 6:16:00 AM
Vinyl chloride	U	1.0		µg/L	1	8/29/2006 6:16:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Field Blank
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 11:10:00 AM
<b>Lab ID:</b>	0608273-05A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: LDS
1,1,1,2-Tetrachloroethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,1,1-Trichloroethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,1,2,2-Tetrachloroethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,1,2-Trichloro-1,2,2-trifluoroethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,1,2-Trichloroethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,1-Dichloroethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,1-Dichloroethene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,1-Dichloropropene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2,3-Trichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2,3-Trichloropropane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2,4,5-Tetramethylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2,4-Trichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2,4-Trimethylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2-Dibromo-3-chloropropane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2-Dibromoethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2-Dichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2-Dichloroethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,2-Dichloropropane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,3,5-Trimethylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,3-Dichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,3-dichloropropane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
1,4-Dichlorobenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
2,2-Dichloropropane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
2-Butanone	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
2-Chloroethyl vinyl ether	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
2-Chlorotoluene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
2-Hexanone	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
2-Propanol	U	50		µg/L	1	8/29/2006 6:53:00 AM
4-Chlorotoluene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
4-Isopropyltoluene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
4-Methyl-2-pentanone	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Acetone	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Acrolein	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Acrylonitrile	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Benzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Bromobenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Bromochloromethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Bromodichloromethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Bromoform	U	1.0		µg/L	1	8/29/2006 6:53:00 AM

<b>Qualifiers:</b>	B Analyte detected in the associated Method Blank	E Value above quantitation range
	H Holding times for preparation or analysis exceeded	J Analyte detected below quantitation limits
	ND Not Detected at the Reporting Limit	S Spike Recovery outside accepted recovery limits
	U Indicates the compound was analyzed for but not detected	X Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Field Blank
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 11:10:00 AM
<b>Lab ID:</b>	0608273-05A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: LDS
Bromomethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Carbon disulfide	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Carbon tetrachloride	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Chlorobenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Chlorodifluoromethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Chloroethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Chloroform	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Chloromethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
cis-1,2-Dichloroethene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
cis-1,3-Dichloropropene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Dibromochloromethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Dibromomethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Dichlorodifluoromethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Diisopropyl ether	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Ethanol	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Ethyl acetate	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Ethylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Freon-114	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Hexachlorobutadiene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Isopropyl acetate	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Isopropylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
m,p-Xylene	U	2.0		µg/L	1	8/29/2006 6:53:00 AM
Methyl tert-butyl ether	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Methylene chloride	18	1.0	B	µg/L	1	8/29/2006 6:53:00 AM
n-Amyl acetate	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Naphthalene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
n-Butyl acetate	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
n-Butylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
n-Propyl acetate	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
n-Propylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
o-Xylene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
p-Diethylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
p-Ethyltoluene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
sec-Butylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Styrene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
t-Butyl alcohol	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
tert-Butylbenzene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Tetrachloroethene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Toluene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 01-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Field Blank
<b>Lab Order:</b>	0608273	<b>Tag Number:</b>	10005
<b>Project:</b>	PEN-Glenwood Landing	<b>Collection Date:</b>	8/24/2006 11:10:00 AM
<b>Lab ID:</b>	0608273-05A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: LDS
trans-1,2-Dichloroethene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
trans-1,3-Dichloropropene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Trichloroethene	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Trichlorofluoromethane	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Vinyl acetate	U	1.0		µg/L	1	8/29/2006 6:53:00 AM
Vinyl chloride	U	1.0		µg/L	1	8/29/2006 6:53:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing  
 Test No: SW8260B

**QC SUMMARY REPORT**  
**SURROGATE RECOVERIES**  
**VOLATILE SW-846 METHOD 8260**

Matrix: W

Sample ID	BR4FBZ	BZMED8	DBFM
0608273-01A	97.2	99.1	134
0608273-02A	96.3	94.5	137
0608273-03A	88.9	88.1	114
0608273-04A	88.5	99.3	124
0608273-05A	130	101	130
LCS082806BY	92.7	96.0	88.2
VBLK082806BY	96.6	96.2	95.7

Acronym	Surrogate	QC Limits
BR4FBZ	= Surr: 4-Bromofluorobenzene	69-134
BZMED8	= Surr: Toluene-d8	71-123
DBFM	= Surr: Dibromofluoromethane	70-135

\* Surrogate recovery outside acceptance limits

American Analytical Laboratories, LLC.

Date: 01-Sep-06

# ANALYTICAL QC SUMMARY REPORT

CLIENT: P.W. Grosser Consulting  
Work Order: 0608273  
Project: PEN-Glenwood Landing

BatchID: R20119

Sample ID: VBLK082806BY	SampType: MBLK	TestCode: DryFull8260_	Units: µg/Kg	Prep Date:	RunNo: 20119
Client ID: PBS	Batch ID: R20119	TestNo: SW8260B		Analysis Date: 8/29/2006	SeqNo: 244607

Analyte	Result	PQL	SPK value	SPK RefVal	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPD limit	Qual
1,1,1,2-Tetrachloroethane	U	5.0									
1,1,1-Trichloroethane	U	5.0									
1,1,2,2-Tetrachloroethane	U	5.0									
1,1,2-Trichloro-1,2,2-trifluoroethane	U	5.0									
1,1,2-Trichloroethane	U	5.0									
1,1-Dichloroethane	U	5.0									
1,1-Dichloroethene	U	5.0									
1,1-Dichloropropene	U	5.0									
1,2,3-Trichlorobenzene	U	5.0									
1,2,3-Trichloropropane	U	5.0									
1,2,4,5-Tetramethylbenzene	U	5.0									
1,2,4-Trichlorobenzene	U	5.0									
1,2,4-Trimethylbenzene	U	5.0									
1,2-Dibromo-3-chloropropane	U	5.0									
1,2-Dibromoethane	U	5.0									
1,2-Dichlorobenzene	U	5.0									
1,2-Dichloroethane	U	5.0									
1,2-Dichloropropane	U	5.0									
1,3,5-Trimethylbenzene	U	5.0									
1,3-Dichlorobenzene	U	5.0									
1,3-dichloropropane	U	5.0									
1,4-Dichlorobenzene	U	5.0									
2,2-Dichloropropane	U	5.0									
2-Butenone	U	5.0									
2-Chloroethyl vinyl ether	U	5.0									
2-Chlorotoluene	U	5.0									
2-Hexanone	U	5.0									
2-Propanol	U	50									
4-Chlorotoluene	U	5.0									
4-Isopropyltoluene	U	5.0									

Qualifiers: E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608273  
**Project:** PEN-Glenwood Landing

**ANALYTICAL QC SUMMARY REPORT**

**BatchID:** R20119

Sample ID: VBLK082806BY	SampType: MBLK	TestCode: DryFu118260_	Units: µg/Kg	Prep Date:	RunNo: 20119						
Client ID: PBS	Batch ID: R20119	TestNo: SW8260B		Analysis Date: 8/29/2006	SeqNo: 244607						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
4-Methyl-2-pentanone	U	5.0									
Acetone	U	5.0									
Acrolein	U	25									
Acrylonitrile	U	5.0									
Benzene	U	5.0									
Bromobenzene	U	5.0									
Bromochloromethane	U	5.0									
Bromodichloromethane	U	5.0									
Bromoform	U	5.0									
Bromomethane	U	5.0									
Carbon disulfide	U	5.0									
Carbon tetrachloride	U	5.0									
Chlorobenzene	U	5.0									
Chlorodifluoromethane	U	5.0									
Chloroethane	U	5.0									
Chloroform	U	5.0									
Chloromethane	U	5.0									
cis-1,2-Dichloroethene	U	5.0									
cis-1,3-Dichloropropene	U	5.0									
Dibromochloromethane	U	5.0									
Dibromomethane	U	5.0									
Dichlorodifluoromethane	U	5.0									
Diisopropyl ether	U	5.0									
Ethanol	U	25									
Ethyl acetate	U	5.0									
Ethylbenzene	U	5.0									
Freon-114	U	5.0									
Hexachlorobutadiene	U	5.0									
Isopropyl acetate	U	5.0									
Isopropylbenzene	U	5.0									
m,p-Xylene	U	10									

**Qualifiers:** E Value above quantitation range  
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H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

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CLIENT: P.W. Grosser Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing

**ANALYTICAL QC SUMMARY REPORT**

BatchID: R20119

Sample ID: VBLK082806BY    SampType: MBLK    TestCode: DryFull8260\_    Units: µg/Kg    Prep Date:    RunNo: 20119  
 Client ID: PBS    Batch ID: R20119    TestNo: SW8260B    Analysis Date: 8/29/2006    SeqNo: 244608

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Methyl tert-butyl ether	U	5.0									
Methylene chloride	15	5.0									
n-Amyl acetate	U	5.0									
Naphthalene	U	5.0									
n-Butyl acetate	U	5.0									
n-Butylbenzene	U	5.0									
n-Propyl acetate	U	5.0									
n-Propylbenzene	U	5.0									
o-Xylene	U	5.0									
p-Diethylbenzene	U	5.0									
p-Ethyltoluene	U	5.0									
sec-Butylbenzene	U	5.0									
Styrene	U	5.0									
t-Butyl alcohol	U	5.0									
tert-Butylbenzene	U	5.0									
Tetrachloroethene	U	5.0									
Toluene	U	5.0									
trans-1,2-Dichloroethene	U	5.0									
trans-1,3-Dichloropropene	U	5.0									
Trichloroethene	U	5.0									
Trichlorofluoromethane	U	5.0									
Vinyl acetate	U	5.0									
Vinyl chloride	U	5.0									

Sample ID: LCS082806BY    SampType: LCS    TestCode: DryFull8260\_    Units: µg/Kg    Prep Date:    RunNo: 20119  
 Client ID: LCSS    Batch ID: R20119    TestNo: SW8260B    Analysis Date: 8/28/2006    SeqNo: 244608

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	46	5.0	50.00	0	92.5	42	138				
Benzene	56	5.0	50.00	0	111	45	137				
Chlorobenzene	64	5.0	50.00	0	128	41	143				

**Qualifiers:** E Value above quantitation range    H Holding times for preparation or analysis exceeded    J Analyte detected below quantitation li  
 ND Not Detected at the Reporting Limit    R RPD outside accepted recovery limits    S Spike Recovery outside accepted reco  
 U Indicates the compound was analyzed for but not detect

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# ANALYTICAL QC SUMMARY REPORT

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing

BatchID: R20119

Sample ID: LCS082806BY	SampType: LCS	TestCode: DryFul8260	Units: µg/Kg	Prep Date:	RunNo: 20119						
Client ID: LCSS	Batch ID: R20119	TestNo: SW8260B		Analysis Date: 8/28/2006	SeqNo: 244608						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Toluene	59	5.0	50.00	0	118	38	141				
Trichloroethene	48	5.0	50.00	0	95.6	39	136				

Sample ID: VBLK082806BY	SampType: MBLK	TestCode: Full8260_W	Units: µg/L	Prep Date:	RunNo: 20119						
Client ID: PBW	Batch ID: R20119	TestNo: SW8260B		Analysis Date: 8/29/2006	SeqNo: 244603						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1,1,2-Tetrachloroethane	U	1.0									
1,1,1-Trichloroethane	U	1.0									
1,1,2,2-Tetrachloroethane	U	1.0									
1,1,2-Trichloro-1,2,2-trifluoroethane	U	1.0									
1,1,2-Trichloroethane	U	1.0									
1,1-Dichloroethane	U	1.0									
1,1-Dichloroethene	U	1.0									
1,1-Dichloropropene	U	1.0									
1,2,3-Trichlorobenzene	U	1.0									
1,2,3-Trichloropropane	U	1.0									
1,2,4,5-Tetramethylbenzene	U	1.0									
1,2,4-Trichlorobenzene	U	1.0									
1,2,4-Trimethylbenzene	U	1.0									
1,2-Dibromo-3-chloropropane	U	1.0									
1,2-Dibromoethane	U	1.0									
1,2-Dichlorobenzene	U	1.0									
1,2-Dichloroethane	U	1.0									
1,2-Dichloropropane	U	1.0									
1,3,5-Trimethylbenzene	U	1.0									
1,3-Dichlorobenzene	U	1.0									
1,3-dichloropropane	U	1.0									
1,4-Dichlorobenzene	U	1.0									
2,2-Dichloropropane	U	1.0									
2-Butanone	U	1.0									

**Qualifiers:** E Value above quantitation range    H Holding times for preparation or analysis exceeded    J Analyte detected below quantitation li  
 ND Not Detected at the Reporting Limit    R RPD outside accepted recovery limits    S Spike Recovery outside accepted reco  
 U Indicates the compound was analyzed for but not detected

CLIENT: P. W. Grosser Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing

**ANALYTICAL QC SUMMARY REPORT**

BatchID: R20119

Sample ID: VBLK082806BY	Sample Type: MBLK	Test Code: FUI8260_W	Units: µg/L	Prep Date:	Run No: 20119						
Client ID: PBW	Batch ID: R20119	Test No: SW8260B		Analysis Date: 8/29/2006	Seq No: 244603						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
2-Chloroethyl vinyl ether	U	1.0									
2-Chlorotoluene	U	1.0									
2-Hexanone	U	1.0									
2-Propanol	U	50									
4-Chlorotoluene	U	1.0									
4-Isopropyltoluene	U	1.0									
4-Methyl-2-pentanone	U	1.0									
Acetone	U	1.0									
Acrolein	U	1.0									
Acrylonitrile	U	1.0									
Benzene	U	1.0									
Bromobenzene	U	1.0									
Bromochloromethane	U	1.0									
Bromodichloromethane	U	1.0									
Bromoform	U	1.0									
Bromomethane	U	1.0									
Carbon disulfide	U	1.0									
Carbon tetrachloride	U	1.0									
Chlorobenzene	U	1.0									
Chlorodifluoromethane	U	1.0									
Chloroethane	U	1.0									
Chloroform	U	1.0									
Chloromethane	U	1.0									
cis-1,2-Dichloroethene	U	1.0									
cis-1,3-Dichloropropene	U	1.0									
Dibromochloromethane	U	1.0									
Dibromomethane	U	1.0									
Dichlorodifluoromethane	U	1.0									
Diisopropyl ether	U	1.0									
Ethanol	U	1.0									
Ethyl acetate	U	1.0									

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# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P. W. Grosser Consulting  
**Work Order:** 0608273  
**Project:** PEN-Glenwood Landing

**BatchID:** R20119

Sample ID: <b>VBLK082806BY</b>	SampType: <b>MBLK</b>	TestCode: <b>Full8260_W</b>	Units: <b>µg/L</b>	Prep Date:	RunNo: <b>20119</b>						
Client ID: <b>PBW</b>	Batch ID: <b>R20119</b>	TestNo: <b>SW8260B</b>		Analysis Date: <b>8/29/2006</b>	SeqNo: <b>244603</b>						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Ethylbenzene	U	1.0									
Freon-114	U	1.0									
Hexachlorobutadiene	U	1.0									
Isopropyl acetate	U	1.0									
Isopropylbenzene	U	1.0									
m,p-Xylene	U	2.0									
Methyl tert-butyl ether	U	1.0									
Methylene chloride	15	1.0									
n-Amyl acetate	U	1.0									
Naphthalene	U	1.0									
n-Butyl acetate	U	1.0									
n-Butylbenzene	U	1.0									
n-Propyl acetate	U	1.0									
n-Propylbenzene	U	1.0									
o-Xylene	U	1.0									
p-Diethylbenzene	U	1.0									
p-Ethyltoluene	U	1.0									
sec-Butylbenzene	U	1.0									
Styrene	U	1.0									
t-Butyl alcohol	U	1.0									
tert-Butylbenzene	U	1.0									
Tetrachloroethene	U	1.0									
Toluene	U	1.0									
trans-1,2-Dichloroethene	U	1.0									
trans-1,3-Dichloropropene	U	1.0									
Trichloroethene	U	1.0									
Trichlorofluoromethane	U	1.0									
Vinyl acetate	U	1.0									
Vinyl chloride	U	1.0									

**Qualifiers:** E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
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 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing

# ANALYTICAL QC SUMMARY REPORT

BatchID: R20119

Sample ID: LCS082806BY	SampType: LCS	TestCode: Full8260_W	Units: µg/L	Prep Date:	RunNo: 20119
Client ID: LCSW	Batch ID: R20119	TestNo: SW8260B		Analysis Date: 8/28/2006	SeqNo: 244604

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1,1,2-Tetrachloroethane	U	1.0	0	0	0						
1,1,1-Trichloroethane	U	1.0	0	0	0						
1,1,2,2-Tetrachloroethane	U	1.0	0	0	0						
1,1,2-Trichloro-1,2,2-trifluoroethane	U	1.0	0	0	0						
1,1,2-Trichloroethane	U	1.0	0	0	0						
1,1-Dichloroethane	U	1.0	0	0	0						
1,1-Dichloroethene	46	1.0	50.00	0	92.5	51	139				
1,1-Dichloropropene	U	1.0	0	0	0						
1,2,3-Trichlorobenzene	U	1.0	0	0	0						
1,2,3-Trichloropropane	U	1.0	0	0	0						
1,2,4,5-Tetramethylbenzene	U	1.0	0	0	0						
1,2,4-Trichlorobenzene	U	1.0	0	0	0						
1,2,4-Trimethylbenzene	U	1.0	0	0	0						
1,2-Dibromo-3-chloropropane	U	1.0	0	0	0						
1,2-Dibromoethane	U	1.0	0	0	0						
1,2-Dichlorobenzene	U	1.0	0	0	0						
1,2-Dichloroethane	U	1.0	0	0	0						
1,2-Dichloropropane	U	1.0	0	0	0						
1,3,5-Trimethylbenzene	U	1.0	0	0	0						
1,3-Dichlorobenzene	U	1.0	0	0	0						
1,3-dichloropropane	U	1.0	0	0	0						
1,4-Dichlorobenzene	U	1.0	0	0	0						
2,2-Dichloropropane	U	1.0	0	0	0						
2-Butanone	U	1.0	0	0	0						
2-Chloroethyl vinyl ether	U	1.0	0	0	0						
2-Chlorotoluene	U	1.0	0	0	0						
2-Hexanone	U	1.0	0	0	0						
2-Propanol	U	50	0	0	0						
4-Chlorotoluene	U	1.0	0	0	0						
4-Isopropyltoluene	U	1.0	0	0	0						
4-Methyl-2-pentanone	U	1.0	0	0	0						

Qualifiers: E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing

**ANALYTICAL QC SUMMARY REPORT**

BatchID: R20119

Sample ID: LCS082806BY	SampType: LCS	TestCode: Full8260_W	Units: µg/L	Prep Date:	RunNo: 20119						
Client ID: LCSW	Batch ID: R20119	TestNo: SW8260B		Analysis Date: 8/28/2006	SeqNo: 244604						
Analyte	Result	PQL	SPK value	SPK RefVal	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Acetone	U	1.0	0	0	0						
Acrolein	U	1.0	0	0	0						
Acrylonitrile	U	1.0	0	0	0						
Benzene	56	1.0	50.00	0	111	53	135				
Bromobenzene	U	1.0	0	0	0						
Bromochloromethane	U	1.0	0	0	0						
Bromodichloromethane	U	1.0	0	0	0						
Bromoform	U	1.0	0	0	0						
Bromomethane	U	1.0	0	0	0						
Carbon disulfide	U	1.0	0	0	0						
Carbon tetrachloride	U	1.0	0	0	0						
Chlorobenzene	64	1.0	50.00	0	128	58	142				
Chlorodifluoromethane	U	1.0	0	0	0						
Chloroethane	U	1.0	0	0	0						
Chloroform	U	1.0	0	0	0						
Chloromethane	U	1.0	0	0	0						
cis-1,2-Dichloroethene	U	1.0	0	0	0						
cis-1,3-Dichloropropene	U	1.0	0	0	0						
Dibromochloromethane	U	1.0	0	0	0						
Dibromomethane	U	1.0	0	0	0						
Dichlorodifluoromethane	U	1.0	0	0	0						
Diisopropyl ether	U	1.0	0	0	0						
Ethanol	U	1.0	0	0	0						
Ethyl acetate	U	1.0	0	0	0						
Ethylbenzene	U	1.0	0	0	0						
Freon-114	U	1.0	0	0	0						
Hexachlorobutadiene	U	1.0	0	0	0						
Isopropyl acetate	U	1.0	0	0	0						
Isopropylbenzene	U	1.0	0	0	0						
m,p-Xylene	U	2.0	0	0	0						
Methyl tert-butyl ether	U	1.0	0	0	0						

**Qualifiers:** E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

J Analyte detected below quantitation limit  
 S Spike Recovery outside accepted recovery

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CLIENT: P. W. Grosser Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing

# ANALYTICAL QC SUMMARY REPORT

BatchID: R20119

Sample ID: LCS082806BY    SampType: LCS    TestCode: Full8260\_W    Units: µg/L    Prep Date:    RunNo: 20119  
 Client ID: LCSW    Batch ID: R20119    TestNo: SW8260B    Analysis Date: 8/28/2006    SeqNo: 244604

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Methylene chloride	19	1.0	0	0	0						B
Naphthalene	U	1.0	0	0	0						
n-Butyl acetate	U	1.0	0	0	0						
n-Butylbenzene	U	1.0	0	0	0						
n-Propyl acetate	U	1.0	0	0	0						
n-Propylbenzene	U	1.0	0	0	0						
o-Xylene	U	1.0	0	0	0						
p-Diethylbenzene	U	1.0	0	0	0						
p-Ethyltoluene	U	1.0	0	0	0						
sec-Butylbenzene	U	1.0	0	0	0						
Styrene	U	1.0	0	0	0						
t-Butyl alcohol	U	1.0	0	0	0						
tert-Butylbenzene	U	1.0	0	0	0						
Tetrachloroethene	U	1.0	0	0	0						
Toluene	59	1.0	50.00	0	118	50	142				
trans-1,2-Dichloroethene	U	1.0	0	0	0						
trans-1,3-Dichloropropene	U	1.0	0	0	0						
Trichloroethene	48	1.0	50.00	0	95.6	53	131				
Trichlorofluoromethane	U	1.0	0	0	0						
Vinyl acetate	U	1.0	0	0	0						
Vinyl chloride	U	1.0	0	0	0						

**Qualifiers:** E Value above quantitation range    H Holding times for preparation or analysis exceeded    J Analytic detected below quantitation li  
 ND Not Detected at the Reporting Limit    R RPD outside accepted recovery limits    S Spike Recovery outside accepted reco  
 U Indicates the compound was analyzed for but not detected

American Analytical Laboratories, LLC.

Date: 31-Aug-06

ANALYTICAL QC SUMMARY REPORT

CLIENT: P.W. Grosser Consulting

Work Order: 0608273

Project: PEN-Glenwood Landing

BatchID: 11897

Sample ID:	MB-11897	Samp Type:	MBLK	Test Code:	Dry8270_Soil	Units:	µg/Kg	Prep Date:	8/27/2006	Run No:	20070		
Client ID:	PBS	Batch ID:	11897	Test No:	SW8270D	SW3550A		Analysis Date:	8/29/2006	Seq No:	243771		
Analyte		Result		FQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene		U		120									
1,2-Dichlorobenzene		U		120									
1,3-Dichlorobenzene		U		120									
1,4-Dichlorobenzene		U		120									
2,4,5-Trichlorophenol		U		120									
2,4,6-Trichlorophenol		U		120									
2,4-Dichlorophenol		U		120									
2,4-Dimethylphenol		U		120									
2,4-Dinitrophenol		U		120									
2,4-Dinitrotoluene		U		120									
2,6-Dinitrotoluene		U		120									
2-Chloronaphthalene		U		120									
2-Chlorophenol		U		120									
2-Methylnaphthalene		U		120									
2-Methylphenol		U		120									
2-Nitroaniline		U		120									
2-Nitrophenol		U		120									
3,3'-Dichlorobenzidine		U		120									
3+4-Methylphenol		U		120									
3-Nitroaniline		U		120									
4,6-Dinitro-2-methylphenol		U		120									
4-Bromophenyl phenyl ether		U		120									
4-Chloro-3-methylphenol		U		120									
4-Chloroaniline		U		120									
4-Chlorophenyl phenyl ether		U		120									
4-Nitroaniline		U		120									
4-Nitrophenol		U		120									
Acenaphthene		U		120									
Acenaphthylene		U		120									
Aniline		U		120									

Qualifiers: E Value above quantitation range H Holding times for preparation or analysis exceeded J Analyte detected below quantitation li  
 ND Not Detected at the Reporting Limit R RPD outside accepted recovery limits S Spike Recovery outside accepted reco  
 U Indicates the compound was analyzed for but not detect

CLIENT: P.W. Grosset Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing

# ANALYTICAL QC SUMMARY REPORT

BatchID: 11897

Sample ID: MB-11897	Sample Type: MBLK	Test Code: Dry8270_Soil	Units: µg/Kg	Prep Date: 8/27/2006	Run No: 20070						
Client ID: PBS	Batch ID: 11897	Test No: SW8270D	SW3550A	Analysis Date: 8/29/2006	Seq No: 243771						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Anthracene	U	120									
Azobenzene	U	120									
Benzidine	U	120									
Benzo(a)anthracene	U	120									
Benzo(a)pyrene	U	120									
Benzo(b)fluoranthene	U	120									
Benzo(g,h,i)perylene	U	120									
Benzo(k)fluoranthene	U	120									
Benzoic acid	U	120									
Benzyl alcohol	U	120									
Bis(2-chloroethoxy)methane	U	120									
Bis(2-chloroethyl)ether	U	120									
Bis(2-chloroisopropyl)ether	U	120									
Bis(2-ethylhexyl)phthalate	U	120									
Butyl benzyl phthalate	U	120									
Carbazole	U	120									
Chrysene	U	120									
Dibenzo(a,h)anthracene	U	120									
Dibenzofuran	U	120									
Diethyl phthalate	U	120									
Dimethyl phthalate	U	120									
Di-n-butyl phthalate	U	120									
Di-n-octyl phthalate	U	120									
Fluoranthene	U	120									
Fluorene	U	120									
Hexachlorobenzene	U	120									
Hexachlorobutadiene	U	120									
Hexachlorocyclopentadiene	U	120									
Hexachloroethane	U	120									
Indeno(1,2,3-c,d)pyrene	U	120									
Isophorone	U	120									

Qualifiers: E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits

J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608273  
**Project:** PEN-Glenwood Landing

**BatchID: 11897**

Sample ID: MB-11897	SampType: MBLK	TestCode: Dry8270_Soil	Units: µg/Kg	Prep Date: 8/27/2006	RunNo: 20070
Client ID: PBS	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/29/2006	SeqNo: 243771

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Naphthalene	U	120									
Nitrobenzene	U	120									
N-Nitrosodimethylamine	U	120									
N-Nitrosodi-n-propylamine	U	120									
N-Nitrosodiphenylamine	U	120									
Pentachlorophenol	U	120									
Phenanthrene	U	120									
Phenol	U	120									
Pyrene	U	120									
Pyridine	U	120									

Sample ID: LCS-11897	SampType: LCS	TestCode: Dry8270_Soil	Units: µg/Kg	Prep Date: 8/27/2006	RunNo: 20070
Client ID: LCSS	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/29/2006	SeqNo: 243772

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	3000	120	3999	0	75.0	23	122				
1,4-Dichlorobenzene	3200	120	3999	0	79.6	22	124				
2,4-Dinitrotoluene	2300	120	3999	0	58.4	21	119				
2-Chlorophenol	3000	120	3999	0	75.6	23	125				
4-Chloro-3-methylphenol	3100	120	3999	0	77.2	28	121				
4-Nitrophenol	1500	120	3999	0	37.8	15	115				
Acenaphthene	2900	120	3999	0	73.4	24	124				
N-Nitrosodi-n-propylamine	3800	120	3999	0	95.9	25	127				
Pentachlorophenol	850	120	3999	0	21.3	14	112				
Phenol	2300	120	3999	0	58.5	19	125				
Pyrene	3500	120	3999	0	88.3	31	125				

Sample ID: MB-11897	SampType: MBLK	TestCode: Dry8270BN_S	Units: µg/Kg	Prep Date: 8/27/2006	RunNo: 20070
Client ID: PBS	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/29/2006	SeqNo: 243841

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	3000	120	3999	0	75.0	23	122				
1,4-Dichlorobenzene	3200	120	3999	0	79.6	22	124				
2,4-Dinitrotoluene	2300	120	3999	0	58.4	21	119				
2-Chlorophenol	3000	120	3999	0	75.6	23	125				
4-Chloro-3-methylphenol	3100	120	3999	0	77.2	28	121				
4-Nitrophenol	1500	120	3999	0	37.8	15	115				
Acenaphthene	2900	120	3999	0	73.4	24	124				
N-Nitrosodi-n-propylamine	3800	120	3999	0	95.9	25	127				
Pentachlorophenol	850	120	3999	0	21.3	14	112				
Phenol	2300	120	3999	0	58.5	19	125				
Pyrene	3500	120	3999	0	88.3	31	125				

**Qualifiers:** E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detect

J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

Page 3 of 9

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608273  
**Project:** PEN-Glenwood Landing

# ANALYTICAL QC SUMMARY REPORT

**BatchID: 11897**

Sample ID: MB-11897	Samp Type: MBLK	TestCode: Dry8270BN_S	Units: µg/Kg	Prep Date: 8/27/2006	RunNo: 20070						
Client ID: PBS	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/29/2006	SeqNo: 243841						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,2,4-Trichlorobenzene	U	120									
1,2-Dichlorobenzene	U	120									
1,3-Dichlorobenzene	U	120									
1,4-Dichlorobenzene	U	120									
2,4-Dinitrotoluene	U	120									
2,6-Dinitrotoluene	U	120									
2-Chloronaphthalene	U	120									
2-Methylnaphthalene	U	120									
2-Nitroaniline	U	120									
3,3'-Dichlorobenzidine	U	120									
3-Nitroaniline	U	120									
4-Bromophenyl phenyl ether	U	120									
4-Chloroaniline	U	120									
4-Chlorophenyl phenyl ether	U	120									
4-Nitroaniline	U	150									
Acenaphthene	U	120									
Acenaphthylene	U	120									
Aniline	U	120									
Anthracene	U	120									
Azobenzene	U	120									
Benzidine	U	120									
Benzo(a)anthracene	U	120									
Benzo(a)pyrene	U	120									
Benzo(b)fluoranthene	U	120									
Benzo(g,h,i)perylene	U	120									
Benzo(k)fluoranthene	U	120									
Benzyl alcohol	U	120									
Bis(2-chloroethoxy)methane	U	120									
Bis(2-chloroethyl)ether	U	120									
Bis(2-chloroisopropyl)ether	U	120									
Bis(2-ethylhexyl)phthalate	U	120									

**Qualifiers:** E Value above quantitation range    H Holding times for preparation or analysis exceeded    J Analyte detected below quantitation limit  
 ND Not Detected at the Reporting Limit    R RPD outside accepted recovery limits    S Spike Recovery outside accepted recovery  
 U Indicates the compound was analyzed for but not detected

**ANALYTICAL QC SUMMARY REPORT**

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608273  
**Project:** PEN-Glenwood Landing

**BatchID:** 11897

Sample ID: MB-11897	SampType: MBLK	TestCode: Dry8270BN_S	Units: µg/Kg	Prep Date: 8/27/2006	RunNo: 20070						
Client ID: PBS	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/29/2006	SeqNo: 243841						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Butyl benzyl phthalate	U	120									
Carbazole	U	120									
Chrysene	U	120									
Dibenzo(a,h)anthracene	U	120									
Dibenzofuran	U	120									
Diethyl phthalate	U	120									
Dimethyl phthalate	U	120									
Di-n-butyl phthalate	U	120									
Di-n-octyl phthalate	U	120									
Fluoranthene	U	120									
Fluorene	U	120									
Hexachlorobenzene	U	120									
Hexachlorobutadiene	U	120									
Hexachlorocyclopentadiene	U	120									
Hexachloroethane	U	120									
Indeno(1,2,3-c,d)pyrene	U	120									
Isophorone	U	120									
Naphthalene	U	120									
Nitrobenzene	U	120									
N-Nitrosodimethylamine	U	120									
N-Nitrosodi-n-propylamine	U	120									
N-Nitrosodiphenylamine	U	120									
Phenanthrene	U	120									
Pyrene	U	120									
Pyridine	U	120									

Sample ID: LCS-11897	SampType: LCS	TestCode: Dry8270BN_S	Units: µg/Kg	Prep Date: 8/27/2006	RunNo: 20070						
Client ID: LCSS	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/29/2006	SeqNo: 243842						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	3000	120	3999	0	75.0	23	122				

**Qualifiers:** E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

# ANALYTICAL QC SUMMARY REPORT

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608273  
 Project: PEN-Glenwood Landing

BatchID: 11897

Sample ID: LCS-11897	SampType: LCS	TestCode: Dry8270BN_S	Units: µg/Kg	Prep Date: 8/27/2006	RunNo: 20070
Client ID: LCS	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/29/2006	SeqNo: 243842

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,4-Dichlorobenzene	3200	120	3999	0	79.6	22	124				
2,4-Dinitrotoluene	2300	120	3999	0	58.4	21	119				
Acenaphthene	2900	120	3999	0	73.4	24	124				
N-Nitrosodi-n-propylamine	3800	120	3999	0	95.9	25	127				
Pyrene	3500	120	3999	0	88.3	31	125				

Sample ID: 0608210-19A-MS	SampType: MS	TestCode: Dry8270BN_S	Units: µg/Kg-dry	Prep Date: 8/27/2006	RunNo: 20070
Client ID: ZZZZZZ	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/30/2006	SeqNo: 243844

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	3300	120	4125	0	79.7	23	122				
1,4-Dichlorobenzene	3400	120	4125	0	82.9	22	124				
2,4-Dinitrotoluene	2500	120	4125	0	61.7	21	119				
Acenaphthene	3400	120	4125	0	82.4	24	124				
N-Nitrosodi-n-propylamine	4300	120	4125	0	104	25	127				
Pyrene	3900	120	4125	0	94.3	31	125				

Sample ID: 0608210-19A-MSD	SampType: MSD	TestCode: Dry8270BN_S	Units: µg/Kg-dry	Prep Date: 8/27/2006	RunNo: 20070
Client ID: ZZZZZZ	Batch ID: 11897	TestNo: SW8270D	SW3550A	Analysis Date: 8/30/2006	SeqNo: 243845

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	3200	120	4125	0	78.7	23	122	3286	1.22	20	
1,4-Dichlorobenzene	3400	120	4125	0	81.5	22	124	3421	1.70	20	
2,4-Dinitrotoluene	2600	120	4125	0	62.2	21	119	2546	0.802	20	
Acenaphthene	3200	120	4125	0	78.6	24	124	3399	4.66	20	
N-Nitrosodi-n-propylamine	4200	120	4125	0	101	25	127	4299	3.19	20	
Pyrene	3800	120	4125	0	91.2	31	125	3891	3.30	20	

Qualifiers: E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation limit  
 S Spike Recovery outside accepted recovery limits

# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608273  
**Project:** PEN-Glenwood Landing

**BatchID: R20053**

Sample ID: PBS	SampType: MIBLK	TestCode: TAGM_MET_	Units: mg/Kg	Prep Date:	RunNo: 20053						
Client ID: PBS	Batch ID: R20053	TestNo: SW6010B		Analysis Date: 8/29/2006	SeqNo: 243065						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Arsenic	U	0.500									
Barium	U	0.400									
Cadmium	U	0.200									
Chromium	U	0.400									
Lead	U	0.300									
Selenium:	U	0.500									
Silver	U	0.400									

Sample ID: LCSS	SampType: LCS	TestCode: TAGM_MET_	Units: mg/Kg	Prep Date:	RunNo: 20053						
Client ID: LCSS	Batch ID: R20053	TestNo: SW6010B		Analysis Date: 8/29/2006	SeqNo: 243066						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Arsenic	40.8	0.500	40.00	0	102	68	125				
Barium	41.3	0.400	40.00	0	103	69	123				
Cadmium	40.7	0.200	40.00	0	102	67	123				
Chromium	41.8	0.400	40.00	0	105	68	124				
Lead	41.5	0.300	40.00	0	104	66	125				
Selenium	41.2	0.500	40.00	0	103	67	124				
Silver	39.6	0.400	40.00	0	99.0	69	121				

Sample ID: 0608214-11A MS	SampType: MS	TestCode: TAGM_MET_	Units: mg/Kg-dry	Prep Date:	RunNo: 20053						
Client ID: ZZZZZZ	Batch ID: R20053	TestNo: SW6010B		Analysis Date: 8/29/2006	SeqNo: 243095						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Arsenic	21.5	0.549	21.95	3.658	81.3	68	125				
Barium	85.9	0.439	21.95	76.02	44.9	69	123				S
Cadmium	18.5	0.220	21.95	0.6971	81.3	67	123				
Chromium	33.2	0.439	21.95	18.30	68.0	68	124				
Lead	74.1	0.329	21.95	65.24	40.6	66	125				S
Selenium	15.3	0.549	21.95	0	69.5	67	124				
Silver	17.6	0.439	21.95	0	80.3	69	121				

**Qualifiers:** E Value above quantitation range    H Holding times for preparation or analysis exceeded    J Analyte detected below quantitation li  
 ND Not Detected at the Reporting Limit    R RPD outside accepted recovery limits    S Spike Recovery outside accepted reco  
 U Indicates the compound was analyzed for but not detected

# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608273  
**Project:** PEN-Glenwood Landing

**BatchID:** R20053

Sample ID:	0608214-11A MSD	SampType	MSD	TestCode:	TAGM_MET_	Units:	mg/Kg-dry	Prep Date:	RunNo:	20053	
Client ID:	ZZZZZZ	Batch ID:	R20053	TestID:	SW6010B			Analyz Date:	8/29/2006	Sample:	243096
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Arsenic	21.9	0.549	21.95	3.658	83.2	68	125	21.50	1.97	20	
Barium	86.3	0.439	21.95	76.02	46.9	69	123	85.87	0.507	20	S
Cadmium	18.6	0.220	21.95	0.6971	81.5	67	123	18.53	0.264	20	
Chromium	33.5	0.439	21.95	18.30	69.4	68	124	33.23	0.893	20	
Lead	74.3	0.329	21.95	65.24	41.3	66	125	74.14	0.228	20	S
Selenium	15.0	0.549	21.95	0	68.3	67	124	15.26	1.74	20	
Silver	17.5	0.439	21.95	0	79.7	69	121	17.62	0.686	20	

Squalific - Sample matrix interference KBK

**Qualifiers:** E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected  
 H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608273  
**Project:** PEN-Glenwood Landing

**BatchID:** R20064

Sample ID: PBS	Sample Type: MBLK	Test Code: DRYHG_S	Units: mg/Kg	Prep Date:	Run No: 20064						
Client ID: PBS	Batch ID: R20064	Test No: SW7471B		Analysis Date: 8/29/2006	Seq No: 243646						
Analyte	Result:	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Mercury	U	0.0100									

Sample ID: LCSS	Sample Type: LCS	Test Code: DRYHG_S	Units: mg/Kg	Prep Date:	Run No: 20064						
Client ID: LCSS	Batch ID: R20064	Test No: SW7471B		Analysis Date: 8/29/2006	Seq No: 243647						
Analyte	Result:	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Mercury	0.194	0.0100	0.2000	0	96.8	71.1	126				

Sample ID: 0608206-06A MS	Sample Type: MS	Test Code: DRYHG_S	Units: mg/Kg-dry	Prep Date:	Run No: 20064						
Client ID: ZZZZZZ	Batch ID: R20064	Test No: SW7471B		Analysis Date: 8/29/2006	Seq No: 243665						
Analyte	Result:	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Mercury	0.762	0.0175	0.3496	0.4942	76.7	71.1	126				

Sample ID: 0608206-06A MSD	Sample Type: MSD	Test Code: DRYHG_S	Units: mg/Kg-dry	Prep Date:	Run No: 20064						
Client ID: ZZZZZZ	Batch ID: R20064	Test No: SW7471B		Analysis Date: 8/29/2006	Seq No: 243666						
Analyte	Result:	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Mercury	0.775	0.0175	0.3496	0.4942	80.5	71.1	126	0.7623	1.71	20	

**Qualifiers:** E Value above quantitation range      H Holding times for preparation or analysis exceeded      J Analyte detected below quantitation li  
 ND Not Detected at the Reporting Limit      R RPD outside accepted recovery limits      S Spike Recovery outside accepted reco  
 U Indicates the compound was analyzed for but not detect

Friday, September 08, 2006

Jim Rhodes  
P. W. Grosser Consulting  
630 Johnson Avenue  
Suite 7  
Bohemia, NY 11716  
TEL: (631) 589-6353  
FAX (631) 589-8705

RE: PEN-1 Shore Rd., Glenwood Landing

Order No.: 0608317

Dear Jim Rhodes:

American Analytical Laboratories, LLC. received 12 sample(s) on 8/30/2006 for the analyses presented in the following report.

Samples were analyzed in accordance with the test procedures documented on the chain of custody and detailed throughout the text of this report.

The limits provided in the data package are analytical reporting limits and not Federal or Local mandated values to which the sample results should be compared.

There were no problems with the analyses and all data for associated QC met laboratory specifications. If there are any exceptions a Case Narrative is provided in the report.

If you have any questions regarding these tests results, please do not hesitate to call (631) 454-6100 or email me directly at lbeyer@american-analytical.com.

Sincerely,

  
Lori Beyer  
Lab Director

**American Analytical Laboratories, LLC.**

Date: 08-Sep-06

---

**CLIENT:** P.W. Grosser Consulting  
**Project:** PEN-1 Shore Rd., Glenwood Landing  
**Lab Order:** 0608317**Work Order Sample Summary**

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<b>Lab Sample ID</b>	<b>Client Sample ID</b>	<b>Tag Number</b>	<b>Date Collected</b>	<b>Date Received</b>
0608317-01A	SB-4 [12.5'-15']	11018	8/30/2006 9:15:00 AM	8/30/2006
0608317-02A	SB-4 [27.5'-30']	11018	8/30/2006 9:30:00 AM	8/30/2006
0608317-03A	SB-5 [12.5'-15']	11018	8/30/2006 10:40:00 AM	8/30/2006
0608317-04A	SB-5 [27.5'-30']	11018	8/30/2006 10:50:00 AM	8/30/2006
0608317-05A	SB-6 [12.5'-15']	11018	8/30/2006 12:40:00 PM	8/30/2006
0608317-06A	SB-6 [27.5'-30']	11018	8/30/2006 12:50:00 PM	8/30/2006
0608317-07A	SB-7 [17.5'-20']	11018	8/30/2006 2:00:00 PM	8/30/2006
0608317-08A	SB-7 [27.5'-30']	11018	8/30/2006 2:15:00 PM	8/30/2006
0608317-09A	SB-8 [17.5'-20']	11018	8/30/2006 2:50:00 PM	8/30/2006
0608317-10A	SB-8 [13'-15']	11018	8/30/2006 2:45:00 PM	8/30/2006
0608317-11A	Field Blank	11018	8/30/2006 12:00:00 PM	8/30/2006
0608317-12A	Trip Blank	11018	8/30/2006 12:00:00 PM	8/30/2006



**AMERICAN ANALYTICAL LABORATORIES, LLC**  
**56 TOLEDO STREET**  
**FARMINGDALE, NEW YORK 11735**  
**TELEPHONE: (631) 454-6100      FAX: (631) 454-8027**

**DATA REPORTING QUALIFIERS**

For reporting results, the following "Results Qualifiers" are used:

<b>Value</b>	If the result is greater than or equal to the detection limit, report the value
<b>U</b>	Indicates the compound was analyzed for but was not detected. Report the minimum detection limit for the sample with the U, i.e. "10U". This is not necessarily the instrument detection limit attainable for this particular sample based on any concentration or dilution that may have been required.
<b>J</b>	Indicates an estimated value. The flag is used: <ol style="list-style-type: none"><li>(1) When estimating a concentration for a tentatively identified compound (library search hits, where a 1:1 response is assumed.)</li><li>(2) When the mass spectral data indicated the identification, however the result was less than the specified detection limit greater than zero. If the detection limit was 10ug/L and a concentration of 3ug/L was calculated report as 3J. This flag is used when similar situations arise on any organic parameter i.e. Pesticide, PCBs and others.</li></ol>
<b>B</b>	Indicates the analyte was found in the blank as well as the sample report "10B".
<b>E</b>	Indicates the analytes concentration exceeds the calibrated range of the instrument for that specific analysis.
<b>D</b>	This flag identifies all compounds identified in an analysis at a secondary dilution factor.
<b>P</b>	This flag is used for Pesticide / PCB target analyte when there is >25% difference for detected concentrations between the two GC Columns. The higher of the two values is reported on Form I and flagged with a "P".
<b>N</b>	This flag indicates presumptive evidence of a compound. This is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It applies to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the flag is not used.
<b>H</b>	Indicates sample was received and/or analyzed outside of The method allowable holding time

**American Analytical Laboratories, LLC.**

Date: 08-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	SB-8 [13'-15']
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 2:45:00 PM
<b>Lab ID:</b>	0608317-10A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>MERCURY</b>		<b>SW7471B</b>		<b>SW7471B</b>		Analyst: <b>WN</b>
Mercury	U	0.0116		mg/Kg-dry	1	8/31/2006 3:20:00 PM
<b>PERCENT MOISTURE</b>		<b>D2216</b>				Analyst: <b>PA</b>
Percent Moisture	18.4	0		wt%	1	8/31/2006
<b>NCDH METALS</b>		<b>SW6010B</b>		<b>SW3050A</b>		Analyst: <b>JP</b>
Arsenic	U	0.601		mg/Kg-dry	1	8/31/2006 1:57:50 PM
Barium	3.06	0.481		mg/Kg-dry	1	8/31/2006 1:57:50 PM
Cadmium	0.835	0.240		mg/Kg-dry	1	8/31/2006 1:57:50 PM
Chromium	1.13	0.481		mg/Kg-dry	1	8/31/2006 1:57:50 PM
Lead	38.6	0.361		mg/Kg-dry	1	8/31/2006 1:57:50 PM
Selenium	U	0.601		mg/Kg-dry	1	8/31/2006 1:57:50 PM
Silver	U	0.481		mg/Kg-dry	1	8/31/2006 1:57:50 PM
<b>SEMIVOLATILE SW-846 METHOD 8270</b>		<b>SW8270D</b>		<b>SW3550A</b>		Analyst: <b>SB</b>
1,2,4-Trichlorobenzene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
1,2-Dichlorobenzene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
1,3-Dichlorobenzene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
1,4-Dichlorobenzene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2,4,5-Trichlorophenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2,4,6-Trichlorophenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2,4-Dichlorophenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2,4-Dimethylphenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2,4-Dinitrophenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2,4-Dinitrotoluene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2,6-Dinitrotoluene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2-Chloronaphthalene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2-Chlorophenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2-Methylnaphthalene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2-Methylphenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2-Nitroaniline	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
2-Nitrophenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
3,3'-Dichlorobenzidine	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
3+4-Methylphenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
3-Nitroaniline	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
4,6-Dinitro-2-methylphenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
4-Bromophenyl phenyl ether	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
4-Chloro-3-methylphenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
4-Chloroaniline	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
4-Chlorophenyl phenyl ether	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
4-Nitroaniline	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 08-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	SB-8 [13'-15']
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 2:45:00 PM
<b>Lab ID:</b>	0608317-10A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>SEMIVOLATILE SW-846 METHOD 8270</b>		<b>SW8270D</b>	<b>SW3550A</b>			Analyst: <b>SB</b>
4-Nitrophenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Acenaphthene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Acenaphthylene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Aniline	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Anthracene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Azobenzene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Benzidine	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Benzo(a)anthracene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Benzo(a)pyrene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Benzo(b)fluoranthene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Benzo(g,h,i)perylene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Benzo(k)fluoranthene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Benzoic acid	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Benzyl alcohol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Bis(2-chloroethoxy)methane	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Bis(2-chloroethyl)ether	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Bis(2-chloroisopropyl)ether	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Bis(2-ethylhexyl)phthalate	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Butyl benzyl phthalate	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Carbazole	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Chrysene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Dibenzo(a,h)anthracene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Dibenzofuran	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Diethyl phthalate	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Dimethyl phthalate	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Di-n-butyl phthalate	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Di-n-octyl phthalate	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Fluoranthene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Fluorene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Hexachlorobenzene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Hexachlorobutadiene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Hexachlorocyclopentadiene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Hexachloroethane	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Indeno(1,2,3-c,d)pyrene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Isophorone	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Naphthalene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Nitrobenzene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
N-Nitrosodimethylamine	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
N-Nitrosodi-n-propylamine	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 08-Sep-06

<b>CLIENT:</b>	P. W. Grosser Consulting	<b>Client Sample ID:</b>	SB-8 [13'-15']
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 2:45:00 PM
<b>Lab ID:</b>	0608317-10A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>SEMIVOLATILE SW-846 METHOD 8270</b>		<b>SW8270D</b>		<b>SW3550A</b>		Analyst: <b>SB</b>
N-Nitrosodiphenylamine	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Pentachlorophenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Phenanthrene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Phenol	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Pyrene	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
Pyridine	U	150		µg/Kg-dry	1	9/5/2006 7:04:00 PM
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: <b>MMR</b>
1,1,1,2-Tetrachloroethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,1,1-Trichloroethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,1,2,2-Tetrachloroethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,1,2-Trichloro-1,2,2-trifluoroethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,1,2-Trichloroethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,1-Dichloroethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,1-Dichloroethene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,1-Dichloropropene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2,3-Trichlorobenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2,3-Trichloropropane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2,4,5-Tetramethylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2,4-Trichlorobenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2,4-Trimethylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2-Dibromo-3-chloropropane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2-Dibromoethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2-Dichlorobenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2-Dichloroethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,2-Dichloropropane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,3,5-Trimethylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,3-Dichlorobenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,3-dichloropropane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
1,4-Dichlorobenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
2,2-Dichloropropane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
2-Butanone	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
2-Chloroethyl vinyl ether	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
2-Chlorotoluene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
2-Hexanone	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
2-Propanol	U	61		µg/Kg-dry	1	9/1/2006 6:50:00 AM
4-Chlorotoluene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
4-Isopropyltoluene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
4-Methyl-2-pentanone	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Acetone	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 08-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	SB-8 [13'-15']
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 2:45:00 PM
<b>Lab ID:</b>	0608317-10A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>						
		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: MMR
Acrolein	U	31		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Acrylonitrile	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Benzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Bromobenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Bromochloromethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Bromodichloromethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Bromoform	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Bromomethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Carbon disulfide	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Carbon tetrachloride	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Chlorobenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Chlorodifluoromethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Chloroethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Chloroform	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Chloromethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
cis-1,2-Dichloroethene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
cis-1,3-Dichloropropene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Dibromochloromethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Dibromomethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Dichlorodifluoromethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Diisopropyl ether	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Ethanol	U	31		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Ethyl acetate	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Ethylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Freon-114	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Hexachlorobutadiene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Isopropyl acetate	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Isopropylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
m,p-Xylene	U	12		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Methyl tert-butyl ether	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Methylene chloride	9.9	6.1	B	µg/Kg-dry	1	9/1/2006 6:50:00 AM
n-Amyl acetate	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Naphthalene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
n-Butyl acetate	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
n-Butylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
n-Propyl acetate	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
n-Propylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
o-Xylene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
p-Diethylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
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**American Analytical Laboratories, LLC.**

Date: 08-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	SB-8 [13'-15']
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 2:45:00 PM
<b>Lab ID:</b>	0608317-10A	<b>Matrix:</b>	SOIL

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		Analyst: MMR
p-Ethyltoluene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
sec-Butylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Styrene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
t-Butyl alcohol	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
tert-Butylbenzene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Tetrachloroethene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Toluene	26	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
trans-1,2-Dichloroethene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
trans-1,3-Dichloropropene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Trichloroethene	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Trichlorofluoromethane	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Vinyl acetate	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM
Vinyl chloride	U	6.1		µg/Kg-dry	1	9/1/2006 6:50:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
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# American Analytical Laboratories, LLC.

Date: 08-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Field Blank
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 12:00:00 PM
<b>Lab ID:</b>	0608317-11A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
1,1,1,2-Tetrachloroethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,1,1-Trichloroethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,1,2,2-Tetrachloroethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,1,2-Trichloro-1,2,2-trifluoroethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,1,2-Trichloroethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,1-Dichloroethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,1-Dichloroethene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,1-Dichloropropene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2,3-Trichlorobenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2,3-Trichloropropane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2,4,5-Tetramethylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2,4-Trichlorobenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2,4-Trimethylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2-Dibromo-3-chloropropane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2-Dibromoethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2-Dichlorobenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2-Dichloroethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,2-Dichloropropane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,3,5-Trimethylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,3-Dichlorobenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,3-dichloropropane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
1,4-Dichlorobenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
2,2-Dichloropropane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
2-Butanone	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
2-Chloroethyl vinyl ether	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
2-Chlorotoluene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
2-Hexanone	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
2-Propanol	U	50		µg/L	1	9/1/2006 2:54:00 AM
4-Chlorotoluene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
4-Isopropyltoluene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
4-Methyl-2-pentanone	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Acetone	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Acrolein	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Acrylonitrile	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Benzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Bromobenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Bromochloromethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Bromodichloromethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Bromoform	U	1.0		µg/L	1	9/1/2006 2:54:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 08-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Field Blank
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 12:00:00 PM
<b>Lab ID:</b>	0608317-11A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>						
		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
Bromomethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Carbon disulfide	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Carbon tetrachloride	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Chlorobenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Chlorodifluoromethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Chloroethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Chloroform	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Chloromethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
cis-1,2-Dichloroethene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
cis-1,3-Dichloropropene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Dibromochloromethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Dibromomethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Dichlorodifluoromethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Diisopropyl ether	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Ethanol	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Ethyl acetate	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Ethylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Freon-114	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Hexachlorobutadiene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Isopropyl acetate	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Isopropylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
m,p-Xylene	U	2.0		µg/L	1	9/1/2006 2:54:00 AM
Methyl tert-butyl ether	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Methylene chloride	5.1	1.0	B	µg/L	1	9/1/2006 2:54:00 AM
n-Amyl acetate	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Naphthalene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
n-Butyl acetate	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
n-Butylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
n-Propyl acetate	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
n-Propylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
o-Xylene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
p-Diethylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
p-Ethyltoluene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
sec-Butylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Styrene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
t-Butyl alcohol	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
tert-Butylbenzene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Tetrachloroethene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Toluene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM

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	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

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<b>CLIENT:</b>	P. W. Grosser Consulting	<b>Client Sample ID:</b>	Field Blank
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 12:00:00 PM
<b>Lab ID:</b>	0608317-11A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>						Analyst: LDS
trans-1,2-Dichloroethene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
trans-1,3-Dichloropropene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Trichloroethene	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Trichlorofluoromethane	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Vinyl acetate	U	1.0		µg/L	1	9/1/2006 2:54:00 AM
Vinyl chloride	U	1.0		µg/L	1	9/1/2006 2:54:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

**American Analytical Laboratories, LLC.**

Date: 08-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Trip Blank
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 12:00:00 PM
<b>Lab ID:</b>	0608317-12A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>		<b>SW5030A</b>		<b>Analyst: LDS</b>
1,1,1,2-Tetrachloroethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,1,1-Trichloroethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,1,2,2-Tetrachloroethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,1,2-Trichloro-1,2,2-trifluoroethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,1,2-Trichloroethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,1-Dichloroethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,1-Dichloroethene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,1-Dichloropropene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2,3-Trichlorobenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2,3-Trichloropropane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2,4,5-Tetramethylbenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2,4-Trichlorobenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2,4-Trimethylbenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2-Dibromo-3-chloropropane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2-Dibromoethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2-Dichlorobenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2-Dichloroethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,2-Dichloropropane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,3,5-Trimethylbenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,3-Dichlorobenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,3-dichloropropane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
1,4-Dichlorobenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
2,2-Dichloropropane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
2-Butanone	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
2-Chloroethyl vinyl ether	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
2-Chlorotoluene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
2-Hexanone	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
2-Propanol	U	50		µg/L	1	9/1/2006 3:41:00 AM
4-Chlorotoluene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
4-Isopropyltoluene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
4-Methyl-2-pentanone	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Acetone	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Acrolein	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Acrylonitrile	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Benzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Bromobenzene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Bromochloromethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Bromodichloromethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Bromoform	U	1.0		µg/L	1	9/1/2006 3:41:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detecte	X	Value exceeds Maximum Contaminant Level

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Date: 08-Sep-06

<b>CLIENT:</b>	P.W. Grosser Consulting	<b>Client Sample ID:</b>	Trip Blank
<b>Lab Order:</b>	0608317	<b>Tag Number:</b>	11018
<b>Project:</b>	PEN-1 Shore Rd., Glenwood Landing	<b>Collection Date:</b>	8/30/2006 12:00:00 PM
<b>Lab ID:</b>	0608317-12A	<b>Matrix:</b>	LIQUID

Analyses	Result	Limit	Qual	Units	DF	Date Analyzed
<b>VOLATILE SW-846 METHOD 8260</b>		<b>SW8260B</b>	<b>SW5030A</b>			Analyst: LDS
trans-1,2-Dichloroethene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
trans-1,3-Dichloropropene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Trichloroethene	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Trichlorofluoromethane	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Vinyl acetate	U	1.0		µg/L	1	9/1/2006 3:41:00 AM
Vinyl chloride	U	1.0		µg/L	1	9/1/2006 3:41:00 AM

<b>Qualifiers:</b>	B	Analyte detected in the associated Method Blank	E	Value above quantitation range
	H	Holding times for preparation or analysis exceeded	J	Analyte detected below quantitation limits
	ND	Not Detected at the Reporting Limit	S	Spike Recovery outside accepted recovery limits
	U	Indicates the compound was analyzed for but not detected	X	Value exceeds Maximum Contaminant Level

# American Analytical Laboratories, LLC.

Date: 08-Sep-06

CLIENT: P.W. Grosser Consulting

Work Order: 0608317

Project: PEN-1 Shore Rd., Glenwood Landing

## ANALYTICAL QC SUMMARY REPORT

TestNo: SW8260B

Sample ID: VBLKS083106H	SampType: MBLK	TestCode: DryFull8260_	Units: µg/Kg	Prep Date:	RunNo: 20139						
Client ID: PBS	Batch ID: 11999	TestNo: SW8260B	SW5030A	Analysis Date: 8/31/2006	SeqNo: 244853						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1,1,2-Tetrachloroethane	U	5.0									
1,1,1-Trichloroethane	U	5.0									
1,1,2,2-Tetrachloroethane	U	5.0									
1,1,2-Trichloro-1,2,2-trifluoroethane	U	5.0									
1,1,2-Trichloroethane	U	5.0									
1,1-Dichloroethane	U	5.0									
1,1-Dichloroethene	U	5.0									
1,1-Dichloropropene	U	5.0									
1,2,3-Trichlorobenzene	U	5.0									
1,2,3-Trichloropropane	U	5.0									
1,2,4,5-Tetramethylbenzene	U	5.0									
1,2,4-Trichlorobenzene	U	5.0									
1,2,4-Trimethylbenzene	U	5.0									
1,2-Dibromo-3-chloropropane	U	5.0									
1,2-Dibromoethane	U	5.0									
1,2-Dichlorobenzene	U	5.0									
1,2-Dichloroethane	U	5.0									
1,2-Dichloropropane	U	5.0									
1,3,5-Trimethylbenzene	U	5.0									
1,3-Dichlorobenzene	U	5.0									
1,3-dichloropropane	U	5.0									
1,4-Dichlorobenzene	U	5.0									
2,2-Dichloropropane	U	5.0									
2-Butanone	U	5.0									
2-Chloroethyl vinyl ether	U	5.0									
2-Chlorotoluene	U	5.0									
2-Hexanone	U	5.0									
2-Propanol	U	5.0									
4-Chlorotoluene	U	5.0									
4-Isopropyltoluene	U	5.0									

Qualifiers: E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detect

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation limit  
 S Spike Recovery outside accepted recovery

# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608317  
**Project:** PEN-1 Shore Rd., Glenwood Landing

**TestNo:** SW8260B

Sample ID: VBLKS083106H	SampType: MBLK	TestCode: DryFull8260_	Units: µg/Kg
Client ID: PBS	Batch ID: 11999	TestNo: SW8260B	SW5030A
Analyte	Result	PQL	SPK value
		SPK Ref Val	%REC
		LowLimit	HighLimit
		RPD Ref Val	%RPD
		RPDLimit	Qual
			RunNo: 20139
			SeqNo: 244853
			Prep Date:
			Analysis Date: 8/31/2006

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
4-Methyl-2-pentanone	U	5.0									
Acetone	U	5.0									
Acrolein	U	25									
Acrylonitrile	U	5.0									
Benzene	U	5.0									
Bromobenzene	U	5.0									
Bromochloromethane	U	5.0									
Bromodichloromethane	U	5.0									
Bromoform	U	5.0									
Bromomethane	U	5.0									
Carbon disulfide	U	5.0									
Carbon tetrachloride	U	5.0									
Chlorobenzene	U	5.0									
Chlorodifluoromethane	U	5.0									
Chloroethane	U	5.0									
Chloroform	U	5.0									
Chloromethane	U	5.0									
cis-1,2-Dichloroethene	U	5.0									
cis-1,3-Dichloropropene	U	5.0									
Dibromochloromethane	U	5.0									
Dibromomethane	U	5.0									
Dichlorodifluoromethane	U	5.0									
Diisopropyl ether	U	5.0									
Ethanol	U	25									
Ethyl acetate	U	5.0									
Ethylbenzene	U	5.0									
Freon-114	U	5.0									
Hexachlorobutadiene	U	5.0									
Isopropyl acetate	U	5.0									
Isopropylbenzene	U	5.0									
m,p-Xylene	U	10									

**Qualifiers:** E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608317  
**Project:** PEN-1 Shore Rd., Glenwood Landing

**ANALYTICAL QC SUMMARY REPORT**

**TestNo:** SW8260B

Sample ID: VBLKS083106H    SampType: MBLK    TestCode: DryFull8260\_    Units: µg/Kg    Prep Date:    RunNo: 20139  
 Client ID: PBS    Batch ID: 11999    TestNo: SW8260B    SW5030A    Analysis Date: 8/31/2006    SeqNo: 244853

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Methyl tert-butyl ether	U	5.0									
Methylene chloride	3.7	5.0									J
n-Amyl acetate	U	5.0									
Naphthalene	U	5.0									
n-Butyl acetate	U	5.0									
n-Butylbenzene	U	5.0									
n-Propyl acetate	U	5.0									
n-Propylbenzene	U	5.0									
o-Xylene	U	5.0									
p-Diethylbenzene	U	5.0									
p-Ethyltoluene	U	5.0									
sec-Butylbenzene	U	5.0									
Styrene	U	5.0									
t-Butyl alcohol	U	5.0									
tert-Butylbenzene	U	5.0									
Tetrachloroethene	U	5.0									
Toluene	U	5.0									
trans-1,2-Dichloroethene	U	5.0									
trans-1,3-Dichloropropene	U	5.0									
Trichloroethene	U	5.0									
Trichlorofluoromethane	U	5.0									
Vinyl acetate	U	5.0									
Vinyl chloride	U	5.0									

Sample ID: LCSS083106H    SampType: LCS    TestCode: DryFull8260\_    Units: µg/Kg    Prep Date:    RunNo: 20139  
 Client ID: LCSS    Batch ID: 11999    TestNo: SW8260B    SW5030A    Analysis Date: 8/31/2006    SeqNo: 244854

Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1-Dichloroethene	61	5.0	50.00	0	123	42	138				
Benzene	55	5.0	50.00	0	109	45	137				
Chlorobenzene	51	5.0	50.00	0	101	41	143				

**Qualifiers:** E Value above quantitation range    H Holding times for preparation or analysis exceeded    J Analyte detected below quantitation li  
 ND Not Detected at the Reporting Limit    R RPD outside accepted recovery limits    S Spike Recovery outside accepted reco  
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CLIENT: P.W. Grosser Consulting  
 Work Order: 0608317

Project: PEN-1 Shore Rd., Greenwood Landing

# ANALYTICAL QC SUMMARY REPORT

TestNo: SW8260B

Sample ID: LCSS083106H	SampType: LCS	TestCode: DryFull8260_	Units: µg/Kg	Prep Date:	RunNo: 20139						
Client ID: LCSS	Batch ID: 11999	TestNo: SW8260B	SW5030A	Analysis Date: 8/31/2006	SeqNo: 244854						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Toluene	52	5.0	50.00	0	103	38	141				
Trichloroethene	49	5.0	50.00	0	97.4	39	136				

Sample ID: 0608317-10AMS	SampType: MS	TestCode: DryFull8260_	Units: µg/Kg-dry	Prep Date:	RunNo: 20139						
Client ID: SB-8 [13'-15']	Batch ID: 11999	TestNo: SW8260B	SW5030A	Analysis Date: 9/1/2006	SeqNo: 244865						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene	33	6.1	61.29	0	54.1	42	138				
Benzene	40	6.1	61.29	0	65.8	45	137				
Chlorobenzene	38	6.1	61.29	0	61.9	41	143				
Toluene	64	6.1	61.29	25.62	63.1	38	141				
Trichloroethene	25	6.1	61.29	0	41.5	39	136				

Sample ID: 0608317-10AMSD	SampType: MSD	TestCode: DryFull8260_	Units: µg/Kg-dry	Prep Date:	RunNo: 20139						
Client ID: SB-8 [13'-15']	Batch ID: 11999	TestNo: SW8260B	SW5030A	Analysis Date: 9/1/2006	SeqNo: 244866						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,1-Dichloroethene	2.2	6.1	61.29	0	3.60	42	138	33.15	0	20	JS
Benzene	25	6.1	61.29	0	40.7	45	137	40.30	47.2	20	SR
Chlorobenzene	23	6.1	61.29	0	37.6	41	143	37.91	48.7	20	SR
Toluene	35	6.1	61.29	25.62	15.1	38	141	64.29	59.4	20	SR
Trichloroethene	10	6.1	61.29	0	16.4	39	136	25.45	86.6	20	SR

**Qualifiers:** E Value above quantitation range    H Holding times for preparation or analysis exceeded    J Analyte detected below quantitation li  
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**CLIENT:** P. W. Grosser Consulting  
**Work Order:** 0608317

**Project:** PEN-1 Shore Rd., Glenwood Landing

# ANALYTICAL QC SUMMARY REPORT

**TestNo:** SW8270D

Sample ID: <b>MB-11979</b>	SampType: <b>MBLK</b>	TestCode: <b>Dry8270_Soil</b>	Units: <b>µg/Kg</b>	Prep Date: <b>9/2/2006</b>	RunNo: <b>20167</b>						
Client ID: <b>PBS</b>	Batch ID: <b>11979</b>	TestNo: <b>SW8270D</b>	SW <b>3550A</b>	Analysis Date: <b>9/5/2006</b>	SeqNo: <b>245202</b>						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,2,4-Trichlorobenzene	U	120									
1,2-Dichlorobenzene	U	120									
1,3-Dichlorobenzene	U	120									
1,4-Dichlorobenzene	U	120									
2,4,5-Trichlorophenol	U	120									
2,4,6-Trichlorophenol	U	120									
2,4-Dichlorophenol	U	120									
2,4-Dimethylphenol	U	120									
2,4-Dinitrophenol	U	120									
2,4-Dinitrotoluene	U	120									
2,6-Dinitrotoluene	U	120									
2-Chloronaphthalene	U	120									
2-Chlorophenol	U	120									
2-Methylnaphthalene	U	120									
2-Methylphenol	U	120									
2-Nitroaniline	U	120									
2-Nitrophenol	U	120									
3,3'-Dichlorobenzidine	U	120									
3+4-Methylphenol	U	120									
3-Nitroaniline	U	120									
4,6-Dinitro-2-methylphenol	U	120									
4-Bromophenyl phenyl ether	U	120									
4-Chloro-3-methylphenol	U	120									
4-Chloroaniline	U	120									
4-Chlorophenyl phenyl ether	U	120									
4-Nitroaniline	U	120									
4-Nitrophenol	U	120									
Acenaphthene	U	120									
Acenaphthylene	U	120									
Aniline	U	120									
Anthracene	U	120									

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 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608317  
 Project: PEN-1 Shore Rd., Glenwood Landing

# ANALYTICAL QC SUMMARY REPORT

TestNo: SW8270D

Sample ID: MB-11979	SampType: MBLK	TestCode: Dry8270_Soil	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20167						
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245202						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Azobenzene	U	120									
Benzidine	U	120									
Benzo(a)anthracene	U	120									
Benzo(a)pyrene	U	120									
Benzo(b)fluoranthene	U	120									
Benzo(g,h,i)perylene	U	120									
Benzo(k)fluoranthene	U	120									
Benzoic acid	U	120									
Benzyl alcohol	U	120									
Bis(2-chloroethoxy)methane	U	120									
Bis(2-chloroethyl)ether	U	120									
Bis(2-chloroisopropyl)ether	U	120									
Bis(2-ethylhexyl)phthalate	U	120									
Butyl benzyl phthalate	U	120									
Carbazole	U	120									
Chrysene	U	120									
Dibenzo(a,h)anthracene	U	120									
Dibenzofuran	U	120									
Diethyl phthalate	U	120									
Dimethyl phthalate	U	120									
Di-n-butyl phthalate	U	120									
Di-n-octyl phthalate	U	120									
Fluoranthene	U	120									
Fluorene	U	120									
Hexachlorobenzene	U	120									
Hexachlorobutadiene	U	120									
Hexachlorocyclopentadiene	U	120									
Hexachloroethane	U	120									
Indeno(1,2,3-c,d)pyrene	U	120									
Isophorone	U	120									
Naphthalene	U	120									

Qualifiers: E Value above quantitation range  
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 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608317  
**Project:** PEN-1 Shore Rd., Glenwood Landing

**ANALYTICAL QC SUMMARY REPORT**

**TestNo:** SW8270D

Sample ID: MB-11979	SampType: MBLK	TestCode: Dry8270_Soil	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20167						
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245202						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Nitrobenzene	U	120									
N-Nitrosodimethylamine	U	120									
N-Nitrosodi-n-propylamine	U	120									
N-Nitrosodiphenylamine	U	120									
Pentachlorophenol	U	120									
Phenanthrene	U	120									
Phenol	U	120									
Pyrene	U	120									
Pyridine	U	120									

Sample ID: LCS-11979	SampType: LCS	TestCode: Dry8270_Soil	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20167						
Client ID: LCSS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245203						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,2,4-Trichlorobenzene	2200	120	3999	0	54.9	23	122				
1,4-Dichlorobenzene	2400	120	3999	0	60.0	22	124				
2,4-Dinitrotoluene	1500	120	3999	0	38.3	21	119				
2-Chlorophenol	2100	120	3999	0	53.6	23	125				
4-Chloro-3-methylphenol	1900	120	3999	0	46.6	28	121				
4-Nitrophenol	630	120	3999	0	15.7	15	115				
Acenaphthene	2300	120	3999	0	58.0	24	124				
N-Nitrosodi-n-propylamine	3200	120	3999	0	79.5	25	127				
Pentachlorophenol	580	120	3999	0	14.5	14	112				
Phenol	1700	120	3999	0	42.2	19	125				
Pyrene	2400	120	3999	0	60.5	31	125				

Sample ID: MB-11979	SampType: MBLK	TestCode: Dry8270_Soil	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245629						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,2,4-Trichlorobenzene	U	120									
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 ND Not Detected at the Reporting Limit R RPD outside accepted recovery limits S Spike Recovery outside accepted recovery  
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CLIENT: P.W. Grosser Consulting  
 Work Order: 0608317

# ANALYTICAL QC SUMMARY REPORT

Project: PEN-1 Shore Rd., Glenwood Landing

TestNo: SW8270D

Sample ID: MB-11979	SampType: MBLK	TestCode: Dry8270_Soil	Units: µg/Kg	RunNo: 20194
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	SeqNo: 245629
Analyte	Result	PQL	SPK value	SPK RefVal
			%REC	LowLimit
			HighLimit	RPD RefVal
			%RPD	RPDLimit
				Qual

Analyte	Result	PQL	SPK value	SPK RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2-Dichlorobenzene	U	120									
1,3-Dichlorobenzene	U	120									
1,4-Dichlorobenzene	U	120									
2,4,5-Trichlorophenol	U	120									
2,4,6-Trichlorophenol	U	120									
2,4-Dichlorophenol	U	120									
2,4-Dimethylphenol	U	120									
2,4-Dinitrophenol	U	120									
2,4-Dinitrotoluene	U	120									
2,6-Dinitrotoluene	U	120									
2-Chloronaphthalene	U	120									
2-Chlorophenol	U	120									
2-Methylnaphthalene	U	120									
2-Methylphenol	U	120									
2-Nitroaniline	U	120									
2-Nitrophenol	U	120									
3,3'-Dichlorobenzidine	U	120									
3+4-Methylphenol	U	120									
3-Nitroaniline	U	120									
4,6-Dinitro-2-methylphenol	U	120									
4-Bromophenyl phenyl ether	U	120									
4-Chloro-3-methylphenol	U	120									
4-Chloroaniline	U	120									
4-Chlorophenyl phenyl ether	U	120									
4-Nitroaniline	U	120									
4-Nitrophenol	U	120									
Acenaphthene	U	120									
Acenaphthylene	U	120									
Aniline	U	120									
Anthracene	U	120									
Azobenzene	U	120									

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 S Spike Recovery outside accepted recovery

# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608317  
**Project:** PEN-1 Shore Rd., Glenwood Landing

**TestNo:** SW8270D

Sample ID:	MB-11979	Samp Type:	MBLK	TestCode:	Dry8270_Soil	Units:	µg/Kg	Prep Date:	9/2/2006	RunNo:	20194
Client ID:	PBS	Batch ID:	11979	TestNo:	SW8270D	SW3550A		Analysis Date:	9/5/2006	SeqNo:	245629
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Benzidine	U	120									
Benzo(a)anthracene	U	120									
Benzo(a)pyrene	U	120									
Benzo(b)fluoranthene	U	120									
Benzo(g,h,i)perylene	U	120									
Benzo(k)fluoranthene	U	120									
Benzoic acid	U	120									
Benzyl alcohol	U	120									
Bis(2-chloroethoxy)methane	U	120									
Bis(2-chloroethyl)ether	U	120									
Bis(2-chloroisopropyl)ether	U	120									
Bis(2-ethylhexyl)phthalate	U	120									
Butyl benzyl phthalate	U	120									
Carbazole	U	120									
Chrysene	U	120									
Dibenzo(a,h)anthracene	U	120									
Dibenzofuran	U	120									
Diethyl phthalate	U	120									
Dimethyl phthalate	U	120									
D,n-butyl phthalate	U	120									
D,n-octyl phthalate	U	120									
Fluoranthene	U	120									
Fluorene	U	120									
Hexachlorobenzene	U	120									
Hexachlorobutadiene	U	120									
Hexachlorocyclopentadiene	U	120									
Hexachloroethane	U	120									
Indeno(1,2,3-c,d)pyrene	U	120									
Isophorone	U	120									
Naphthalene	U	120									
Nitrobenzene	U	120									

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# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608317  
**Project:** PEN-1 Shore Rd., Glenwood Landing

**TestNo:** SW8270D

Sample ID: MB-11979	SampType: MBLK	TestCode: Dry8270_Soil	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245629						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

N-Nitrosodimethylamine	U	120									
N-Nitrosodi-n-propylamine	U	120									
N-Nitrosodiphenylamine	U	120									
Pentachlorophenol	U	120									
Phenanthrene	U	120									
Phenol	U	120									
Pyrene	U	120									
Pyridine	U	120									

Sample ID: LCS-11979	SampType: LCS	TestCode: Dry8270_Soil	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: LCSS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245631						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,2,4-Trichlorobenzene	2200	120	3999	0	54.9	23	122				
1,4-Dichlorobenzene	2400	120	3999	0	60.0	22	124				
2,4-Dinitrotoluene	1500	120	3999	0	38.3	21	119				
2-Chlorophenol	2100	120	3999	0	53.6	23	125				
4-Chloro-3-methylphenol	1900	120	3999	0	46.6	28	121				
4-Nitrophenol	630	120	3999	0	15.7	15	115				
Acenaphthene	2300	120	3999	0	58.0	24	124				
N-Nitrosodi-n-propylamine	3200	120	3999	0	79.5	25	127				
Pentachlorophenol	580	120	3999	0	14.5	14	112				
Phenol	1700	120	3999	0	42.2	19	125				
Pyrene	2400	120	3999	0	60.5	31	125				

Sample ID: 0608317-10A-MS	SampType: MS	TestCode: Dry8270_Soil	Units: µg/Kg-dry	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: SB-8 [13'-15']	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245641						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

1,2,4-Trichlorobenzene	3200	150	4902	0	65.6	23	122				
1,4-Dichlorobenzene	3500	150	4902	0	71.8	22	124				

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 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608317

Project: PEN-1 Shore Rd., Glenwood Landing

# ANALYTICAL QC SUMMARY REPORT

TestNo: SW8270D

Sample ID: 0608317-10A-MS	SampType: MS	TestCode: Dry8270_Soil	Units: µg/Kg-dry	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: SB-8 [13'-15']	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245641						
Analyte	Result	PQL	SPK value	SPK RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual

2,4-Dinitrotoluene	2200	150	4902	0	45.5	21	119				
2-Chlorophenol	3300	150	4902	0	67.9	23	125				
4-Chloro-3-methylphenol	2800	150	4902	0	56.6	28	121				
4-Nitrophenol	2000	150	4902	0	41.2	15	115				
Acenaphthene	3200	150	4902	0	64.9	24	124				
N-Nitrosodi-n-propylamine	4400	150	4902	0	90.0	25	127				
Pentachlorophenol	1600	150	4902	0	33.3	14	112				
Phenol	3100	150	4902	0	63.7	19	125				
Pyrene	3300	150	4902	0	68.0	31	125				

Sample ID: 0608317-10A-MSD	SampType: MSD	TestCode: Dry8270_Soil	Units: µg/Kg-dry	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: SB-8 [13'-15']	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245642						
Analyte	Result	PQL	SPK value	SPK RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual

1,2,4-Trichlorobenzene	2700	150	4902	0	56.0	23	122				
1,4-Dichlorobenzene	2800	150	4902	0	57.6	22	124				
2,4-Dinitrotoluene	1900	150	4902	0	39.7	21	119				
2-Chlorophenol	2700	150	4902	0	54.5	23	125				
4-Chloro-3-methylphenol	2400	150	4902	0	49.1	28	121				
4-Nitrophenol	1300	150	4902	0	27.4	15	115				
Acenaphthene	2700	150	4902	0	54.2	24	124				
N-Nitrosodi-n-propylamine	3600	150	4902	0	73.2	25	127				
Pentachlorophenol	1400	150	4902	0	28.9	14	112				
Phenol	2100	150	4902	0	42.8	19	125				
Pyrene	2900	150	4902	0	59.8	31	125				

Sample ID: MB-11979	SampType: MBLK	TestCode: Dry8270stars	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245625						
Analyte	Result	PQL	SPK value	SPK RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual

Acenaphthene	U	120									
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CLIENT: P.W. Grosser Consulting  
 Work Order: 0608317

# ANALYTICAL QC SUMMARY REPORT

Project: PEN-1 Shore Rd., Glenwood Landing

TestNo: SW8270D

Sample ID: MB-11979	SampType: MBLK	TestCode: Dry8270stars	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245625						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Anthracene	U	120									
Benzo(a)anthracene	U	120									
Benzo(a)pyrene	U	120									
Benzo(b)fluoranthene	U	120									
Benzo(g,h,i)perylene	U	120									
Benzo(k)fluoranthene	U	120									
Chrysene	U	120									
Dibenzo(a,h)anthracene	U	120									
Fluoranthene	U	120									
Fluorene	U	120									
Indeno(1,2,3-c,d)pyrene	U	120									
Phenanthrene	U	120									
Pyrene	U	120									

Sample ID: LCS-11979	SampType: LCS	TestCode: Dry8270stars	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: LCSS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245626						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Acenaphthene	2300	120	3999	0	58.0	24	124				
Pyrene	2400	120	3999	0	60.5	31	125				

Sample ID: MB-11979	SampType: MBLK	TestCode: TAGM8270Dr	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245643						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

2,4,5-Trichlorophenol	U	120									
2,4-Dichlorophenol	U	120									
2,4-Dinitrophenol	U	120									
2,6-Dinitrotoluene	U	120									
2-Chlorophenol	U	120									
2-Methylnaphthalene	U	120									

Qualifiers: E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

# ANALYTICAL QC SUMMARY REPORT

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608317  
**Project:** PEN-1 Shore Rd., Glenwood Landing

**TestNo:** SW8270D

Sample ID:	MB-11979	SampType:	MBLK	TestCode:	TAGM8270Dr	Units:	µg/Kg	Prep Date:	9/2/2006	RunNo:	20194
Client ID:	PBS	Batch ID:	11979	TestNo:	SW8270D	SW3550A		Analysis Date:	9/5/2006	SeqNo:	245643
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
2-Methylphenol	U	120									
2-Nitroaniline	U	120									
2-Nitrophenol	U	120									
3,3'-Dichlorobenzidine	U	120									
3+4-Methylphenol	U	120									
3-Nitroaniline	U	120									
4-Chloro-3-methylphenol	U	120									
4-Chloroaniline	U	120									
4-Nitrophenol	U	120									
Acenaphthene	U	120									
Acenaphthylene	U	120									
Aniline	U	120									
Anthracene	U	120									
Benzo(a)anthracene	U	120									
Benzo(a)pyrene	U	120									
Benzo(b)fluoranthene	U	120									
Benzo(g,h,i)perylene	U	120									
Benzo(k)fluoranthene	U	120									
Benzoic acid	U	120									
Bis(2-ethylhexyl)phthalate	U	120									
Butyl benzyl phthalate	U	120									
Chrysene	U	120									
Dibenzo(a,h)anthracene	U	120									
Dibenzofuran	U	120									
Diethyl phthalate	U	120									
Dimethyl phthalate	U	120									
Di-n-butyl phthalate	U	120									
Di-n-octyl phthalate	U	120									
Fluoranthene	U	120									
Fluorene	U	120									
Hexachlorobenzene	U	120									

**Qualifiers:** E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
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 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation li  
 S Spike Recovery outside accepted reco

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608317

# ANALYTICAL QC SUMMARY REPORT

Project: PEN-1 Shore Rd., Glenwood Landing

TestNo: SW8270D

Sample ID: MB-11979	SampType: MBLK	TestCode: TAGM8270Dr	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: PBS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245643						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Indeno(1,2,3-c,d)pyrene	U	120									
Isophorone	U	120									
Naphthalene	U	120									
Nitrobenzene	U	120									
Pentachlorophenol	U	120									
Phenanthrene	U	120									
Phenol	U	120									
Pyrene	U	120									

Sample ID: LCS-11979	SampType: LCS	TestCode: TAGM8270Dr	Units: µg/Kg	Prep Date: 9/2/2006	RunNo: 20194						
Client ID: LCSS	Batch ID: 11979	TestNo: SW8270D	SW3550A	Analysis Date: 9/5/2006	SeqNo: 245644						
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
2-Chlorophenol	2100	120	3999	0	53.6	23	125				
4-Chloro-3-methylphenol	1900	120	3999	0	46.6	28	121				
4-Nitrophenol	630	120	3999	0	15.7	15	115				
Acenaphthene	2300	120	3999	0	58.0	24	124				
Pentachlorophenol	560	120	3999	0	14.5	14	112				
Phenol	1700	120	3999	0	42.2	19	125				
Pyrene	2400	120	3999	0	60.5	31	125				

Qualifiers: E Value above quantitation range  
 ND Not Detected at the Reporting Limit  
 U Indicates the compound was analyzed for but not detected

H Holding times for preparation or analysis exceeded  
 R RPD outside accepted recovery limits  
 J Analyte detected below quantitation limit  
 S Spike Recovery outside accepted recovery

**CLIENT:** P.W. Grosser Consulting  
**Work Order:** 0608317  
**Project:** PEN-1 Shore Rd., Glenwood Landi  
**Test No:** SW8260B **Matrix:** S

**QC SUMMARY REPORT  
 SURROGATE RECOVERIES  
 VOLATILE SW-846 METHOD 8260**

Sample ID	BR4FBZ	BZMED8	DBFM					
0608317-01A	95.2	94.7	107					
0608317-02A	90.0	98.0	105					
0608317-03A	83.9	92.5	113					
0608317-04A	95.4	95.2	105					
0608317-05A	81.1	94.7	113					
0608317-06A	94.8	93.7	104					
0608317-07A	112	92.2	107					
0608317-08A	97.0	95.5	104					
0608317-09A	95.8	96.7	103					
0608317-10A	92.4	96.9	111					
0608317-10AMS	86.3	98.9	97.8					
0608317-10AMSD	76.0	98.2	81.3					
0608317-11A	104	86.9	114					
0608317-12A	106	90.3	111					
LCSS083106H	92.2	95.2	106					
VBLKS083106H	92.1	92.9	109					

Acronym	Surrogate	QC Limits
BR4FBZ	= Surr: 4-Bromofluorobenzene	61-123
BZMED8	= Surr: Toluene-d8	63-121
DBFM	= Surr: Dibromofluoromethane	68-142

\* Surrogate recovery outside acceptance limits

CLIENT: P.W. Grosser Consulting  
 Work Order: 0608317  
 Project: PEN-1 Shore Rd., Glenwood Landi  
 Test No: SW8270D Matrix: S

**QC SUMMARY REPORT  
 SURROGATE RECOVERIES**

**SEMIVOLATILE SW-846 METHOD 8270**

Sample ID	NO2BZD5	PH246BR	PH2F	PHEN2F	PHEND14	PHENOLD6		
0608317-10A	66.5	53.2	41.9	67.1	65.6	54.0		
0608317-10A-MS	72.3	63.0	57.6	75.5	71.5	66.1		
0608317-10A-MSD	62.0	53.1	34.3	61.8	63.6	44.9		
LCS-11979	62.7	36.5	31.2	65.8	61.4	43.6		
MB-11979	63.9	42.0	33.3	62.6	59.9	44.1		

Acronym	Surrogate	QC Limits
NO2BZD5	= Surr: Nitrobenzene-d5	21-118
PH246BR	= Surr: 2,4,6-Tribromophenol	22-124
PH2F	= Surr: 2-Fluorophenol	21-123
PHEN2F	= Surr: 2-Fluorobiphenyl	27-119
PHEND14	= Surr: 4-Terphenyl-d14	28-126
PHENOLD6	= Surr: Phenol-d6	18-129

\* Surrogate recovery outside acceptance limits