ENVIRO-PRO
PROFESSIONAL CORPORATION

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92	10/31/01	DIN

November 1, 2001

Mr. Larry Griffin
North Mecklenburg Landfill
15300 Holbrooks Road
Huntersville, NC 28078

RE: Semi-annual Monitor Well Sampling Results
North Mecklenburg Landfill
Huntersville, NC
Project No. EP-1217

Dear Mr. Griffin:

Enviro-Pro, P.C. is pleased to submit this report which describes the field sampling activities and summarizes the semi-annual monitor well analytical results for the subject site.

Field Sampling Activities

On October 6, 2001, Enviro-Pro personnel collected groundwater samples from on-site perimeter monitor wells MW-l, MW-4, MW-5, MW-6, MW-10 and MW-ll. The approximate locations of these wells are indicated on Figure l.

The procedures for groundwater measurement and sampling were as follows:

1) Initially, the monitor well caps were removed to allow the groundwater levels to equilibrate to the ambient atmospheric pressure. Next, the depth to groundwater from a measuring point on top of the well casing was recorded. Water level measurements were obtained using an electronic water level meter. The water level probe was decontaminated between monitor wells with deionized water and isopropyl alcohol.

15101 Regena Lane • Charlotte, NC 28278 • (704) 583-0075 • FAX (704) 583-2673

2) At least three well volumes were removed from each monitor well to purge stagnant water and to ensure that fresh formation water would be sampled. Purging was conducted using dedicated disposable Teflon bailers.

3) Each well was then sampled utilizing laboratory prepared containers, labeled, and packed on ice in a portable cooler for shipment to Shealy Environmental Services, Inc., a North Carolina-certified laboratory located in Cayce, South Carolina. Chain-of-custody documentation is included with the analytical reports in Appendix A.

4) Quality assurance/quality control (QA/QC) measures in the field included wearing disposable sample gloves during sampling activities and changing them between sample locations to protect the groundwater samples from cross-contamination. Analytical QA/QC included a trip blank and a field (rinse) blank analyzed for volatile organic compounds (VOCs) by Method 8260B. Only clean, laboratory-supplied sample containers were utilized.

The field information obtained during well purging is summarized on the Well Development, Purge, and Sample Record included as Appendix B. Groundwater levels decreased in four of the wells, from 0.61 foot in well MW-6 to 1.58 foot in well MW-5 since the previous sampling event on March 23, 2001. The average decrease in water levels for the four wells was slightly over 1 foot. The water levels in MW-1 and MW-11 increased 1.2 and 0.1 foot, respectively, over this same period of time.

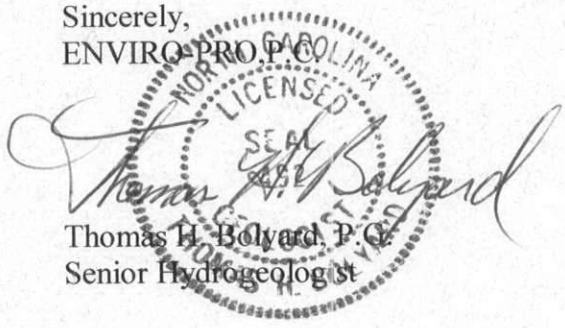
Laboratory Test Results

In accordance with regulatory requirements, the six monitor well samples were analyzed for the eight RCRA metals and volatile organic compounds (VOCs) via EPA Method 8260B by Shealy Environmental Services. Shealy's Report of Analysis is attached as Appendix A, with a summary of groundwater analytical results for this sampling event presented in Table 1.

Laboratory test results indicated that no metals or VOC compounds were present above their respective regulatory limits in any of the monitor well or QA/QC samples tested.

Enviro-Pro appreciates the opportunity to continue to provide our environmental services on your project. Please contact me at (704) 583-0075 if you have any questions concerning this report or when we can be of further service.

Sincerely,
ENVIRO-PRO, P.C.



Thomas H. Bolyard, P.G.
Senior Hydrogeologist

TABLE 1

SUMMARY OF GROUNDWATER ANALYTICAL RESULTS
North Mecklenburg C&D Landfill
Huntersville, North Carolina
October 6, 2001 Sampling Event

	Sample I.D.								
Parameter	MW-1	MW-4	MW-5	MW-6	MW-10	MW-11	Field Blank	Trip Blank	NCAC 2L Standard
Arsenic	BDL	BDL	BDL	BDL	BDL	BDL	NT	NT	0.05
Barium	0.075	0.12	0.15	0.14	0.12	0.22	NT	NT	2.0
Cadmium	BDL	BDL	0.0051	BDL	BDL	BDL	NT	NT	0.005
Chromium	BDL	BDL	BDL	BDL	BDL	BDL	NT	NT	0.05
Lead	BDL	BDL	BDL	0.0047	BDL	BDL	NT	NT	0.015
Mercury	BDL	BDL	BDL	BDL	BDL	0.00093	NT	NT	0.0011
Selenium	BDL	BDL	BDL	BDL	BDL	BDL	NT	NT	0.05
Silver	BDL	BDL	BDL	BDL	BDL	BDL	NT	NT	0.018
All 8260B Compounds	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	

Notes: BDL = Below Detection Limits
NT = Not Tested
All Metal Results Are Presented in Parts Per Million

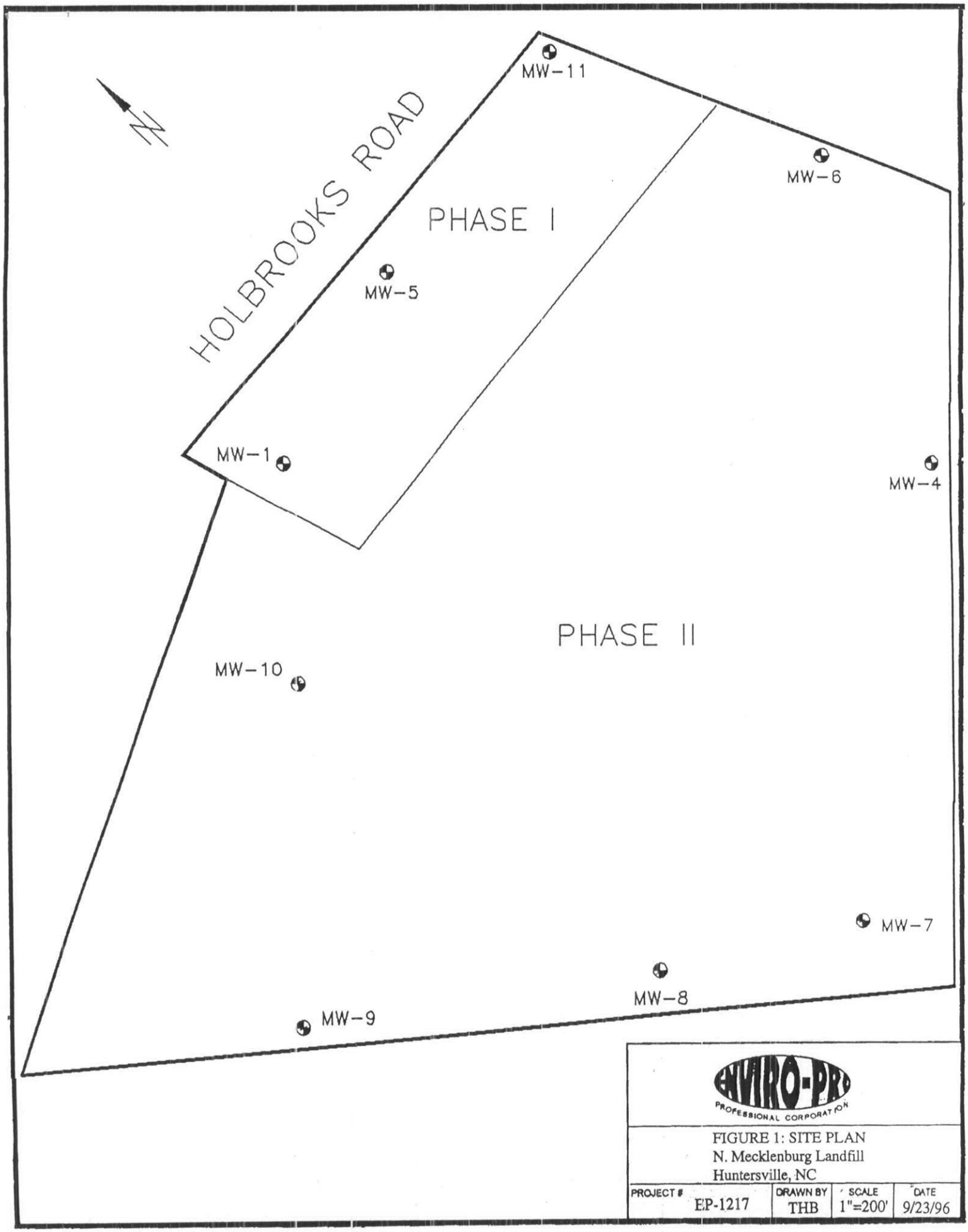


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FIGURE 1: SITE PLAN
N. Mecklenburg Landfill
Huntersville, NC

PROJECT #	DRAWN BY	SCALE	DATE
EP-1217	THB	1"=200'	9/23/96

SHEALY ENVIRONMENTAL SERVICES, INC.
Scientists and Consultants

Report of Analysis

Enviro-Pro, P.C.
15101 Regena Lane
Charlotte, NC 28278
Attention : Tom Bolyard

Project Name : **North Mecklenburg Landfill**

Project Number : **EP -1217**

Lot Number : **CJ09013**

Date Completed : **10/15/2001**

Daniel J. Wright
Project Manager

Total number of pages in report : **15**

This report has been prepared and reviewed in accordance with Shealy's Quality Assurance Management Plan. Any data qualifiers associated with sample analysis are footnoted on the analytical results page(s). This report shall not be reproduced, except in its entirety, without the written approval of Shealy Environmental Services, Inc.

The following non-paginated documents are considered part of this report: Chain Of Custody Record and Sample Receipt Checklist.

SHEALY ENVIRONMENTAL SERVICES, INC.

Report of Analysis

SC DHEC No. 32010 NC DEHNR No. 329

Client : **Enviro-Pro, P.C.**
15101 Regena Lane
Charlotte, NC 28278

Attention : **Tom Bolyard**
Project Name : **North Mecklenburg Landfill** Project Number : **EP -1217**

Description : **MW-1** Date Sampled : 10/06/2001 1345
Lot Number : CJ09013-001 Matrix : Aqueous

Date Received : 10/08/2001

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Acetone	5030B	8260B	ND		20	ug/L		10/12/2001 0331	RED
Benzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Bromodichloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Bromoform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Bromomethane (Methyl bromide)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
2-Butanone (MEK)	5030B	8260B	ND		10	ug/L		10/12/2001 0331	RED
Carbon disulfide	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Carbon tetrachloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Chlorobenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Chloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Chloroform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Chloromethane (Methyl chloride)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Dibromochloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
1,1-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
1,2-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
1,1-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
cis-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
trans-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
1,2-Dichloropropane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
cis-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
trans-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Ethylbenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
2-Hexanone	5030B	8260B	ND		10	ug/L		10/12/2001 0331	RED
4-Methyl-2-pentanone	5030B	8260B	ND		10	ug/L		10/12/2001 0331	RED
Methylene chloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Styrene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
1,1,2,2-Tetrachloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Tetrachloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	: MW-1		Date Sampled	:	10/06/2001 1345
Lot Number	: CJ09013-001		Matrix	:	Aqueous
Client	: **Enviro-Pro, P.C.**				
Date Received	: 10/08/2001				

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Toluene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
1,1,1-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
1,1,2-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Trichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Vinyl chloride	5030B	8260B	ND		2.0	ug/L		10/12/2001 0331	RED
Xylenes (total)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0331	RED
Surrogates									
1,2-Dichloroethane-d4	5030B	8260B	96		70-130	%		10/12/2001 0331	RED
Bromofluorobenzene	5030B	8260B	85		70-130	%		10/12/2001 0331	RED
Toluene-d8	5030B	8260B	80		70-130	%		10/12/2001 0331	RED
CVAA									
Mercury		7470A	ND		0.00010	mg/L	10/10/2001 0900	10/10/2001 1151	NGG
ICP-AES									
Arsenic	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2012	FTS
Barium	3030 C	6010B	**0.075**		0.025	mg/L	10/09/2001 1155	10/11/2001 2012	FTS
Cadmium	3030 C	6010B	ND		0.0020	mg/L	10/09/2001 1155	10/11/2001 2012	FTS
Chromium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2012	FTS
Lead	3030 C	6010B	ND		0.0030	mg/L	10/09/2001 1155	10/12/2001 1615	FTS
Selenium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2012	FTS
Silver	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2012	FTS

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	: MW-4		Date Sampled	: 10/06/2001 1245	
Lot Number	: CJ09013-002		Matrix	: Aqueous	
Client	: **Enviro-Pro, P.C.**				
Date Received	: 10/08/2001				

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Acetone	5030B	8260B	ND		20	ug/L		10/12/2001 0405	RED
Benzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Bromodichloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Bromoform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Bromomethane (Methyl bromide)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
2-Butanone (MEK)	5030B	8260B	ND		10	ug/L		10/12/2001 0405	RED
Carbon disulfide	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Carbon tetrachloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Chlorobenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Chloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Chloroform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Chloromethane (Methyl chloride)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Dibromochloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
1,1-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
1,2-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
1,1-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
cis-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
trans-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
1,2-Dichloropropane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
cis-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
trans-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Ethylbenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
2-Hexanone	5030B	8260B	ND		10	ug/L		10/12/2001 0405	RED
4-Methyl-2-pentanone	5030B	8260B	ND		10	ug/L		10/12/2001 0405	RED
Methylene chloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Styrene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
1,1,2,2-Tetrachloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Tetrachloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Toluene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
1,1,1-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
1,1,2-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Trichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Vinyl chloride	5030B	8260B	ND		2.0	ug/L		10/12/2001 0405	RED
Xylenes (total)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0405	RED
Surrogates									
1,2-Dichloroethane-d4	5030B	8260B	91		70-130	%		10/12/2001 0405	RED
Bromofluorobenzene	5030B	8260B	84		70-130	%		10/12/2001 0405	RED
Toluene-d8	5030B	8260B	77		70-130	%		10/12/2001 0405	RED
CVAA									
Mercury		7470A	ND		0.00010	mg/L	10/10/2001 0900	10/10/2001 1152	NGG

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	:	**MW-4**	Date Sampled	:	10/06/2001 1245
Lot Number	:	CJ09013-002	Matrix	:	Aqueous
Client	:	**Enviro-Pro, P.C.**			
Date Received	:	10/08/2001			

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
ICP-AES									
Arsenic	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2015	FTS
Barium	3030 C	6010B	**0.12**		0.025	mg/L	10/09/2001 1155	10/11/2001 2015	FTS
Cadmium	3030 C	6010B	ND		0.0020	mg/L	10/09/2001 1155	10/11/2001 2015	FTS
Chromium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2015	FTS
Lead	3030 C	6010B	ND		0.0030	mg/L	10/09/2001 1155	10/12/2001 1618	FTS
Selenium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2015	FTS
Silver	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2015	FTS

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	: MW-5		Date Sampled	: 10/06/2001 1330	
Lot Number	: CJ09013-003		Matrix	: Aqueous	
Client	: **Enviro-Pro, P.C.**				
Date Received	: 10/08/2001				

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Acetone	5030B	8260B	ND		20	ug/L		10/12/2001 0438	RED
Benzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Bromodichloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Bromoform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Bromomethane (Methyl bromide)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
2-Butanone (MEK)	5030B	8260B	ND		10	ug/L		10/12/2001 0438	RED
Carbon disulfide	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Carbon tetrachloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Chlorobenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Chloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Chloroform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Chloromethane (Methyl chloride)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Dibromochloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
1,1-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
1,2-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
1,1-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
cis-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
trans-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
1,2-Dichloropropane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
cis-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
trans-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Ethylbenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
2-Hexanone	5030B	8260B	ND		10	ug/L		10/12/2001 0438	RED
4-Methyl-2-pentanone	5030B	8260B	ND		10	ug/L		10/12/2001 0438	RED
Methylene chloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Styrene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
1,1,2,2-Tetrachloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Tetrachloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Toluene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
1,1,1-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
1,1,2-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Trichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Vinyl chloride	5030B	8260B	ND		2.0	ug/L		10/12/2001 0438	RED
Xylenes (total)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0438	RED
Surrogates									
1,2-Dichloroethane-d4	5030B	8260B	94		70-130	%		10/12/2001 0438	RED
Bromofluorobenzene	5030B	8260B	85		70-130	%		10/12/2001 0438	RED
Toluene-d8	5030B	8260B	76		70-130	%		10/12/2001 0438	RED
CVAA									
Mercury		7470A	ND		0.00010	mg/L	10/10/2001 0900	10/10/2001 1153	NGG

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	: **MW-5**		Date Sampled	:	10/06/2001 1330
Lot Number	: CJ09013-003		Matrix	:	Aqueous
Client	: **Enviro-Pro, P.C.**				
Date Received	: 10/08/2001				

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
ICP-AES									
Arsenic	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2018	FTS
Barium	3030 C	6010B	**0.15**		0.025	mg/L	10/09/2001 1155	10/11/2001 2018	FTS
Cadmium	3030 C	6010B	ND		0.0020	mg/L	10/09/2001 1155	10/11/2001 2018	FTS
Chromium	3030 C	6010B	**0.0051**		0.0050	mg/L	10/09/2001 1155	10/11/2001 2018	FTS
Lead	3030 C	6010B	ND		0.0030	mg/L	10/09/2001 1155	10/12/2001 1621	FTS
Selenium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2018	FTS
Silver	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2018	FTS

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	:	**MW-6**	Date Sampled	:	10/06/2001 1300
Lot Number	:	CJ09013-004	Matrix	:	Aqueous
Client	:	**Enviro-Pro, P.C.**			
Date Received	:	10/08/2001			

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Acetone	5030B	8260B	ND		20	ug/L		10/12/2001 0512	RED
Benzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Bromodichloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Bromoform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Bromomethane (Methyl bromide)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
2-Butanone (MEK)	5030B	8260B	ND		10	ug/L		10/12/2001 0512	RED
Carbon disulfide	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Carbon tetrachloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Chlorobenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Chloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Chloroform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Chloromethane (Methyl chloride)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Dibromochloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
1,1-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
1,2-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
1,1-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
cis-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
trans-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
1,2-Dichloropropane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
cis-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
trans-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Ethylbenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
2-Hexanone	5030B	8260B	ND		10	ug/L		10/12/2001 0512	RED
4-Methyl-2-pentanone	5030B	8260B	ND		10	ug/L		10/12/2001 0512	RED
Methylene chloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Styrene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
1,1,2,2-Tetrachloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Tetrachloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Toluene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
1,1,1-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
1,1,2-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Trichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Vinyl chloride	5030B	8260B	ND		2.0	ug/L		10/12/2001 0512	RED
Xylenes (total)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0512	RED
Surrogates									
1,2-Dichloroethane-d4	5030B	8260B	87		70-130	%		10/12/2001 0512	RED
Bromofluorobenzene	5030B	8260B	83		70-130	%		10/12/2001 0512	RED
Toluene-d8	5030B	8260B	78		70-130	%		10/12/2001 0512	RED
CVAA									
Mercury		7470A	ND		0.00010	mg/L	10/10/2001 0900	10/10/2001 1154	NGG

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	:	**MW-6**	Date Sampled	:	10/06/2001 1300
Lot Number	:	CJ09013-004	Matrix	:	Aqueous
Client	:	**Enviro-Pro, P.C.**			
Date Received	:	10/08/2001			

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
ICP-AES									
Arsenic	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2021	FTS
Barium	3030 C	6010B	**0.14**		0.025	mg/L	10/09/2001 1155	10/11/2001 2021	FTS
Cadmium	3030 C	6010B	ND		0.0020	mg/L	10/09/2001 1155	10/11/2001 2021	FTS
Chromium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2021	FTS
Lead	3030 C	6010B	**0.0047**		0.0030	mg/L	10/09/2001 1155	10/12/2001 1624	FTS
Selenium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2021	FTS
Silver	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2021	FTS

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	: **MW-10**		Date Sampled	: 10/06/2001 1400	
Lot Number	: CJ09013-005		Matrix	: Aqueous	
Client	: **Enviro-Pro, P.C.**				
Date Received	: 10/08/2001				

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Acetone	5030B	8260B	ND		20	ug/L		10/12/2001 0545	RED
Benzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Bromodichloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Bromoform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Bromomethane (Methyl bromide)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
2-Butanone (MEK)	5030B	8260B	ND		10	ug/L		10/12/2001 0545	RED
Carbon disulfide	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Carbon tetrachloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Chlorobenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Chloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Chloroform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Chloromethane (Methyl chloride)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Dibromochloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
1,1-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
1,2-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
1,1-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
cis-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
trans-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
1,2-Dichloropropane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
cis-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
trans-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Ethylbenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
2-Hexanone	5030B	8260B	ND		10	ug/L		10/12/2001 0545	RED
4-Methyl-2-pentanone	5030B	8260B	ND		10	ug/L		10/12/2001 0545	RED
Methylene chloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Styrene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
1,1,2,2-Tetrachloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Tetrachloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Toluene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
1,1,1-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
1,1,2-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Trichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Vinyl chloride	5030B	8260B	ND		2.0	ug/L		10/12/2001 0545	RED
Xylenes (total)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0545	RED
Surrogates									
1,2-Dichloroethane-d4	5030B	8260B	94		70-130	%		10/12/2001 0545	RED
Bromofluorobenzene	5030B	8260B	87		70-130	%		10/12/2001 0545	RED
Toluene-d8	5030B	8260B	77		70-130	%		10/12/2001 0545	RED
CVAA									
Mercury		7470A	ND		0.00010	mg/L	10/10/2001 0900	10/10/2001 1155	NGG

PQL = Practical quantitation limit

ND = Not detected at or above the PQL

B = Detected in the method blank

J = Estimated result less than the PQL

E = Quantitation of compound exceeded the calibration range

Soils reported on a dry weight basis unless flagged with a "W"

Description	:	**MW-10**	Date Sampled	:	10/06/2001 1400
Lot Number	:	CJ09013-005	Matrix	:	Aqueous
Client	:	**Enviro-Pro, P.C.**			
Date Received	:	10/08/2001			

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
ICP-AES									
Arsenic	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2024	FTS
Barium	3030 C	6010B	**0.12**		0.025	mg/L	10/09/2001 1155	10/11/2001 2024	FTS
Cadmium	3030 C	6010B	ND		0.0020	mg/L	10/09/2001 1155	10/11/2001 2024	FTS
Chromium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2024	FTS
Lead	3030 C	6010B	ND		0.0030	mg/L	10/09/2001 1155	10/12/2001 1627	FTS
Selenium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2024	FTS
Silver	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2024	FTS

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	:	**MW-11**	Date Sampled	:	10/06/2001 1315
Lot Number	:	CJ09013-006	Matrix	:	Aqueous
Client	:	**Enviro-Pro, P.C.**			
Date Received	:	10/08/2001			

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Acetone	5030B	8260B	ND		20	ug/L		10/12/2001 0619	RED
Benzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Bromodichloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Bromoform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Bromomethane (Methyl bromide)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
2-Butanone (MEK)	5030B	8260B	ND		10	ug/L		10/12/2001 0619	RED
Carbon disulfide	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Carbon tetrachloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Chlorobenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Chloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Chloroform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Chloromethane (Methyl chloride)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Dibromochloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
1,1-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
1,2-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
1,1-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
cis-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
trans-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
1,2-Dichloropropane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
cis-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
trans-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Ethylbenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
2-Hexanone	5030B	8260B	ND		10	ug/L		10/12/2001 0619	RED
4-Methyl-2-pentanone	5030B	8260B	ND		10	ug/L		10/12/2001 0619	RED
Methylene chloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Styrene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
1,1,2,2-Tetrachloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Tetrachloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Toluene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
1,1,1-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
1,1,2-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Trichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Vinyl chloride	5030B	8260B	ND		2.0	ug/L		10/12/2001 0619	RED
Xylenes (total)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0619	RED
Surrogates									
1,2-Dichloroethane-d4	5030B	8260B	88		70-130	%		10/12/2001 0619	RED
Bromofluorobenzene	5030B	8260B	84		70-130	%		10/12/2001 0619	RED
Toluene-d8	5030B	8260B	77		70-130	%		10/12/2001 0619	RED
CVAA									
Mercury		7470A	**0.00093**		0.00010	mg/L	10/10/2001 0900	10/10/2001 1156	NGG

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	:	**MW-11**	Date Sampled	:	10/06/2001 1315
Lot Number	:	CJ09013-006	Matrix	:	Aqueous
Client	:	**Enviro-Pro, P.C.**			
Date Received	:	10/08/2001			

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
ICP-AES									
Arsenic	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2027	FTS
Barium	3030 C	6010B	**0.22**		0.025	mg/L	10/09/2001 1155	10/11/2001 2027	FTS
Cadmium	3030 C	6010B	ND		0.0020	mg/L	10/09/2001 1155	10/11/2001 2027	FTS
Chromium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2027	FTS
Lead	3030 C	6010B	ND		0.0030	mg/L	10/09/2001 1155	10/12/2001 1629	FTS
Selenium	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2027	FTS
Silver	3030 C	6010B	ND		0.0050	mg/L	10/09/2001 1155	10/11/2001 2027	FTS

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	:	**Field Blank**	Date Sampled	:	10/06/2001 1405
Lot Number	:	CJ09013-007	Matrix	:	Aqueous
Client	:	**Enviro-Pro, P.C.**			
Date Received	:	10/08/2001			

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Acetone	5030B	8260B	ND		20	ug/L		10/12/2001 0652	RED
Benzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Bromodichloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Bromoform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Bromomethane (Methyl bromide)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
2-Butanone (MEK)	5030B	8260B	ND		10	ug/L		10/12/2001 0652	RED
Carbon disulfide	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Carbon tetrachloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Chlorobenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Chloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Chloroform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Chloromethane (Methyl chloride)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Dibromochloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
1,1-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
1,2-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
1,1-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
cis-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
trans-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
1,2-Dichloropropane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
cis-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
trans-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Ethylbenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
2-Hexanone	5030B	8260B	ND		10	ug/L		10/12/2001 0652	RED
4-Methyl-2-pentanone	5030B	8260B	ND		10	ug/L		10/12/2001 0652	RED
Methylene chloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Styrene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
1,1,2,2-Tetrachloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Tetrachloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Toluene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
1,1,1-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
1,1,2-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Trichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Vinyl chloride	5030B	8260B	ND		2.0	ug/L		10/12/2001 0652	RED
Xylenes (total)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0652	RED
Surrogates									
1,2-Dichloroethane-d4	5030B	8260B	94		70-130	%		10/12/2001 0652	RED
Bromofluorobenzene	5030B	8260B	89		70-130	%		10/12/2001 0652	RED
Toluene-d8	5030B	8260B	78		70-130	%		10/12/2001 0652	RED

PQL = Practical quantitation limit
ND = Not detected at or above the PQL
B = Detected in the method blank
J = Estimated result less than the PQL
E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

Description	:	**Trip Blank**	Date Sampled	:	10/06/2001
Lot Number	:	CJ09013-008	Matrix	:	Aqueous
Client	:	**Enviro-Pro, P.C.**			
Date Received	:	10/08/2001			

Parameter	Prep Method	Analytical Method	Result	Q	PQL	Units	Prep Date	Analysis Date	Analyst
Volatile Organic Compounds by GC/MS									
Acetone	5030B	8260B	ND		20	ug/L		10/12/2001 0726	RED
Benzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Bromodichloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Bromoform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Bromomethane (Methyl bromide)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
2-Butanone (MEK)	5030B	8260B	ND		10	ug/L		10/12/2001 0726	RED
Carbon disulfide	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Carbon tetrachloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Chlorobenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Chloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Chloroform	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Chloromethane (Methyl chloride)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Dibromochloromethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
1,1-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
1,2-Dichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
1,1-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
cis-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
trans-1,2-Dichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
1,2-Dichloropropane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
cis-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
trans-1,3-Dichloropropene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Ethylbenzene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
2-Hexanone	5030B	8260B	ND		10	ug/L		10/12/2001 0726	RED
4-Methyl-2-pentanone	5030B	8260B	ND		10	ug/L		10/12/2001 0726	RED
Methylene chloride	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Styrene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
1,1,2,2-Tetrachloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Tetrachloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Toluene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
1,1,1-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
1,1,2-Trichloroethane	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Trichloroethene	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Vinyl chloride	5030B	8260B	ND		2.0	ug/L		10/12/2001 0726	RED
Xylenes (total)	5030B	8260B	ND		5.0	ug/L		10/12/2001 0726	RED
Surrogates									
1,2-Dichloroethane-d4	5030B	8260B	92		70-130	%		10/12/2001 0726	RED
Bromofluorobenzene	5030B	8260B	85		70-130	%		10/12/2001 0726	RED
Toluene-d8	5030B	8260B	79		70-130	%		10/12/2001 0726	RED

PQL = Practical quantitation limit
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E = Quantitation of compound exceeded the calibration range
Soils reported on a dry weight basis unless flagged with a "W"

SHEALY **Chain of Custody Record**

SHEALY ENVIRONMENTAL SERVICES, INC.
106 Vantage Point Drive
Cayce, South Carolina 29033
Telephone No. (803) 791-9700 Fax No. (803) 791-9111

Number 07521

Client: Enviro-Pro, P.C.	Project Manager: T. Bolyard	Telephone No. / Fax No. / E-mail: 704-583-0075	Quote No.: 7456
Address: 15101 Regena Ln.	Sampler's Signature: X T. Bolyard	Waybill No.	Page 1 of 1
City: Charlotte / State: NC / Zip Code: 28278	Printed Name: T. Bolyard	Analysis (Attach list if more space is needed.)	
Project Name: NMLF			
Project No.: EP-1217	P.O. No.		Lot No.: CJ09013

Sample ID / Description (Containers for each sample may be combined on one line.)	Date	Time	G=Grab C=Composite	Matrix: Aqueous	Solid	Non-Aqueous	No. of Containers by Preservative Type: Unpres.	H_2SO_4	HNO_3	HCl	NaOH	5035 Kit	8260B	RCRA Metals (3030C Prep.)	Remarks / Cooler I.D.
MW-1	10/6/01	13:45	G	X					1	3			✓	✓	7-Day T-AT
MW-4	"	12:45	G	X					1	3			✓	✓	
MW-5	"	13:30	G	X					1	3			✓	✓	
MW-6	"	13:00	G	X					1	3			✓	✓	
MW-10	"	14:00	G	X					1	3			✓	✓	
MW-11	"	13:15	G	X					1	3			✓	✓	
Field Blank	"	14:05		X						2			✓		
Trip Blank	"			X						2			✓		

Possible Hazard Identification: ☑ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison ☐ Unknown

Sample Disposal: ☐ Return to Client ☑ Disposal by Lab

Note: All samples are retained for six weeks from receipt unless other arrangements are made.

Turn Around Time Required (Prior lab approval required for expedited TAT.): ☑ Standard ☐ Rush (Specify)

QC Requirements (Specify)

	Date	Time		Date	Time
1. Relinquished by: T. Bolyard	10/8/01	1600	1. Received by:	10/8/01	1600
2. Relinquished by:	10/8/01	1635	2. Received by: Apex Express	10/8/01	1635
3. Relinquished by: APEX	10/8/01	1859	3. Laboratory received by: Angelique Torres	10-8-01	1859

Comments

LAB USE ONLY
Received on ice (Circle) (Yes) No Ice Pack

Receipt Temp. 4.1 °C

DISTRIBUTION: WHITE & YELLOW-Return to laboratory with Sample(s); PINK-Field/Client Copy

Shealy Form No. 1001

WELL DEVELOPMENT, PURGE, AND SAMPLE RECORD

CLIENT: North Mecklenburg Landfill
LOCATION: 15300 Holbrooks Rd.
PROJECT NAME/DESCRIPTION: North Meck Landfill
DATE OF WATER LEVEL MEASUREMENTS: 10-06-01

DEVELOP:
PURGING: ✓
SAMPLING: ✓

PROJECT NO.: EP-1217
PROJ. MGR.: FTB
CHECKED BY:
PREPARED BY: FTB

Sample ID	A Well No.	B Sample Date	C Time (hrs)	D Well Depth (ft)	E MP Elevation (ft)	F Depth to Water (ft-bmp)	(E - F) = G Water Level Elevation (ft)	(D - F) = H Hgt. of Water Col. (ft)	I Volume of Water (gal)*	Vol. of Water Pgd. (gal)**	Temp. (°C)	Specific Cond. (umhos)	pH	Turbidity or Appearance
	MW-1	10-06-01	13:45	25		14.98		10.02	1.6	5.0				
	MW-4	"	12:45	42		36.30		5.70	0.9	3.0				
	MW-5	"	13:30	55		30.28		24.72	4.0	12.0				
	MW-6	"	13:00	67		61.13		4.87	0.8	2.5				
	MW-10	"	14:00	19		12.67		6.33	1.0	3.0				
	MW-11	"	13:15	80		69.80		10.20	1.6	5.0				
	Field Blank	"												
	Trip Blank													

Comments:

*To calculate volume of water in the well multiply "H" by 0.163 for a 2" well, 0.652 for a 4" well, or 1.469 for a 6" well.

**To calculate purge volume, multiply "I" by the desired number of well volumes (e.g., 3 to 5).

(e.g., after each well volume).