FIELD INVESTIGATION LETTER REPORT SOIL-GAS CONDUIT INSTALLATION AND SAMPLING AT THE NORTH FRANKLIN STREET SITE

NORTH FRANKLIN STREET SITE SITE #8-49-002 VILLAGE OF WATKINS GLEN, NEW YORK

Prepared For:

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION DIVISION OF ENVIRONMENTAL REMEDIATION WORK ASSIGNMENT D003825-09.5

FINAL

Prepared By:

URS CORPORATION 77 GOODELL STREET BUFFALO, NEW YORK 14203

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October 20, 2005

Mr. David J. Chiusano, Project Manager New York State Department of Environmental Conservation Division of Environmental Remediation 625 Broadway 12th Floor Albany, New York 12233-7013

RE: NYSDEC Standby Contract Active Venting System Operation and Maintenance # D003825-09.5 Soil-Gas Conduit Installation and Sampling North Franklin Street Site, Site No. 8-49-002 Summary of Soil-Gas Conduit Installation and Sampling at the North Franklin Street Site: Letter Report

Dear Mr. Chiusano:

URS Corporation (URS) has completed a soil-gas investigation in the vicinity of the North Franklin Street Site (Figure 1). This work was performed in accordance with the Field Sampling Plan (URS, July 2005).

The fieldwork associated with this investigation consisted of the installation and sampling of thirteen new soil-gas conduits. URS personnel supervised the installation of the soil-gas conduits between July 12, 2005 and July 13, 2005, and conducted the soil-gas conduit sampling on July 18, 2005.

Soil-Gas Conduit Installation and Construction

A total of thirteen soil-gas conduits (SG-01 through SG-13) were installed at the locations shown on Figure 2. The soil-gas conduit locations were approved in the field by a representative of the New York State Department of Environmental Conservation (NYSDEC). The soil-gas conduits were installed by GeoLogic NY, Inc., using a truck-mounted Geoprobe® unit. URS personnel supervised the installation of the soil-gas conduits, which were constructed in accordance with the procedures outlined in the Field Sampling Plan (URS, July 2005). The soil gas conduit construction details are included in Attachment 1. A photographic log of the installation activities and the completed soil-gas conduit locations is included in Attachment 2.

Soil-Gas Sampling, Analysis and Data Usability

All samples were collected using six-liter SUMMA canisters, in accordance with the procedures outlined in the Field Sampling Plan (URS, July 2005). URS collected thirteen one-hour soil gas samples plus two field duplicate samples. A helium tracer gas was used during the collection of the soil-gas samples and no elevated concentrations of helium (>20%) were detected prior to or following the sample collection at any soil-gas conduit location. One outdoor air was collected from an upwind location, located near the SG-1 location. The outdoor air sample was collected from approximately 4 feet above the ground surface

by placing the Summa canister on an elevated platform. Completed sampling logs are provided in Attachment 3. A photographic log of the sampling activities is included in Attachment 2.

After the sampling was completed, the samples were shipped under chain-of-custody (COC) control for VOC analysis via EPA Method TO-15 to Severn Trent Laboratories (STL) located in Knoxville, Tennessee. STL is a New York State Department of Health (NYSDOH) approved laboratory. A copy of the COCs is included in Attachment 3.

The data packages were prepared by the laboratory in accordance with the NYSDEC's Category B Deliverables requirements. These deliverables were reviewed by a URS chemist for compliance with the referenced method, following the guidelines in United States Environmental Protection Agency (USEPA) Region II's *Validating Canisters of Volatile Organics in Ambient Air, Rev. 0*, April 1994. A Data Usability Summary Report (DUSR) was prepared by a URS chemist following the guidelines provided in NYSDEC Division of Environmental Remediation Guidance for the Development of Data Usability Summary Reports, dated June 1999. The DUSR may be found in Attachment 4.

Analytical Summary

The validated analytical results from the soil-gas samples are summarized in Table 1. Concentrations of detected compounds at each sample location are shown on Figure 3. The following is a summary of the analytical results from the soil-gas conduit sampling.

- Volatile organic compounds (VOCs) were detected at every soil-gas conduit location, with the highest concentration detected in the sample collected from SG-03 (benzene at 22,700 ug/m³ and m,p-xylene at 6,430ug/m³). However, chlorinated VOCs were only detected in the samples collected from soil-gas conduits SG-01, SG-02, SG-04, SG-06, SG-07 and SG09 through SG-12, with the highest concentration detected in the sample collected from SG-02 (chloroform at 214.83 ug/m³).
- Tetrachloroethene was detected in samples collected from conduits SG-06, SG-07, SG-09, SG-10 and SG-12, at concentrations ranging from 3.26 ug/m³ to 35.3 ug/m³.
- Trichloroethene was detected in the sample collected from conduit SG-09, at a concentration of 12.4 ug/m³.

A copy of the laboratory report is included in Attachment 4.

It should also be noted that the concentrations listed in Table 1 and Figure 3 were obtained by converting parts per billion, by volume (ppbv) to ug/m³. The ug/m³ results are slightly different from the ppbv values listed on the laboratory forms in Attachment 4.



The following tables, figures and attachments are included as part of this field investigation letter report:

<u>Tables</u> Table 1	Soil-Gas Conduit Analytical Results
<u>Figures</u> Figure 1	Project Site
Figure 2	Soil-Gas Conduit Locations
Figure 3	Soil-Gas Conduit Sampling Results
Attachments	
Attachment 1	Soil-Gas Conduit Construction Details
Attachment 2	Photographic Log
Attachment 3	Sampling Logs and Chain-of-Custody
Attachment 4	Data Usability Summary Report

Should you have any questions or comments, please do not hesitate to contact me at 716-856-5636.

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Sincerely,

URS Corporation

Charles E. Dusel, Jr. Sr. Project Manager

cc: File: 05.35388 (C-1) (11173258)

TABLES

Location ID		SG-01	SG-02	SG-03	SG-04	SG-05
Sample ID		SG-1	SG-2	SG-3	SG-4	SG-8
Matrix		Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas
Depth Interval (ft)		•	·•	•	-	-
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units				_	
Volatile Organic Compounds		···-			<u></u>	
Chloromethane	UG/M3	0.74 J	16.1 UJ	9,090 U	19.0 U	16.5 U
Bromomethane	UG/M3	1.28 U	12.0 U	6,600 U	14.4 U	12.4 U
Vinyl Chloride	ŲG/M3	0.84 U	7.92 UJ	4,350 U	9.46 U	8.18 U
Chloroethane	UG/M3	0.87 U	8.18 U	4,490 U	9.76 U	8.44 U
Methylene Chloride	UG/M3	1.81 J	27.1 UJ	15,300 U	32.0 U	27.8 U
Acetone	UG/M3	19.5 UJ	185 UJ	105,000 U	85.5 J	190 U
Carbon Disulfide	UG/M3	1.03 U	9.97 J	5,300 U	8.41 J	15.3
1,1-Dichloroethene	UG/M3	1.31 U	12.3 UJ	6,740 U	14.7 U	12.7 U
1,1-Dichloroethane	UG/M3	1.34 U	12.6 U	6,880 U	15.0 U	13.0 U
2-Butanone	UG/M3	3.54	12.0 U	12,300 U	27.1 U	23.6 U
Chloroform	UG/M3	1.61 U	215	8,300 U	68.4	15.6 U
1,2-Dichloroethane	UG/M3	1.34 U	12.6 U	6,890 U	15.0 U	13.0 U
1,1,1-Trichloroethane	UG/M3	1.80 U	16.9 U	9,280 U	20.2 U	17.5 U
Carbon Tetrachloride	UG/M3	2.08 U	19.5 U	10,700 U	23.3 U	20.1 U
Bromodichloromethane	UG/M3	2.21 U	20.8 U	11,400 U	24.8 U	21.4 U
1,2-Dichloropropane	UG/M3	1.53 U	14.3 U	7,860 U	17.1 U	14.8 U
cis-1.3-Dichloropropene	UG/M3	1.50 U	14.1 U	7,720 U	16.8 U	14.5 U
Trichloroethene	UG/M3	1.77 U	16.7 U	9,140 U	19.9 U	17.2 U
Benzene	UG/M3	0.51 J	9.90 U	22,700	11.8 U	5.11 J
Dibromochloromethane	UG/M3	2.81 U	26.4 U	14,500 U	31.5 U	27.3 U
trans-1,3-Dichloropropene	UG/M3	1.50 UJ	14.1 U	7,720 UJ	16.8 U	14.5 U
1,1,2-Trichloroethane	UG/M3	1.80 U	16.9 U	9,280 U	20.2 U	17.5 U
Bromoform	UG/M3	3.41 UJ	32.0 U	17,600 UJ	38.3 U	33.1 U

Flags assigned during chemistry validation are shown.

U - The analyte was analyzed for, but not detected. The associated numerical value is at or below the method detection limit

Location ID Sample ID		SG-01	SG-02	SG-03	SG-04	SG-05
		SG-1	SG-2	SG-3	SG-4	SG-5
Matrix	t	Soil Gas				
Depth Interval (ft)		•	•	-	-	-
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units					
Volatile Organic Compounds						
I-Methyl-2-Pentanone	UG/M3	3.36 U	32.0 U	18,000 U	37.7 U	32.8 U
2-Hexanone	UG/M3	3.36 U	32.0 U	18,000 U	37.7 U	32.8 U
etrachloroethene	UG/M3	2.24 U	21.0 U	11,600 U	25.1 U	21.7 U
1.2.2-Tetrachloroethane	UG/M3	2.27 U	21.3 U	11,700 U	25.4 U	22.0 U
Foluene	UG/M3	3.54	23.4	6,410 U	20.7	21.1
Chlorobenzene	UG/M3	1.52 U	14.3 U	7,830 U	17.0 U	14.7 U
thylbenzene	UG/M3	1.43 U	6.51 J	7,380 U	7,82 J	5.64 J
Styrene	UG/M3	1.41 UJ	13.2 U	7,240 UJ	8.52 J	13.6 U
n,p-Xylene	UG/M3	1.39 J	40.0	6,430 J	67.7	40.0
-Xylene	UG/M3	1.43 U	6.51 J	7,380 U	13.5 J	6.95 J
cis-1,2-Dichloroethene	UG/M3	1.31 U	12.3 U	6,740 U	14.7 U	12.7 U
rans-1,2-Dichloroethene	UG/M3	1.31 U	12.3 U	6,740 U	14.7 U	12.7 U

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Location ID		SG-06	SG-06	SG-07	SG-08	SG-09
Sample ID		FD-2	SG-6	SG-7	SG-8	SG-9
Matrix		Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas
Depth Interval (ft)		•	-	•	-	-
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units	Field Duplicate (1-1)				
Volatile Organic Compounds						
Chioromethane	UG/M3	18.8 U	1.84 U	1.98 U	18.6 U	17.6 U
Bromomethane	UG/M3	14.0 U	1.40 U	1.51 U	14.0 U	13.2 U
Vinyl Chloride	UG/M3	9.20 U	0.92 U	1.00 U	9.20 U	8.69 U
Chloroethane	UG/M3	9.50 U	0.95 U	1.03 U	9.50 U	8.97 U
Methylene Chloride	UG/M3	31.6 U	3.09	3.33 U	31.3 U	29.5 U
Acetone	UG/M3	216 U	20.2 J	22.8 U	214 U	202 U
Carbon Disulfide	UG/M3	11.2 U	1.34	1.21 U	68.5	4.67 J
1,1-Dichloroethene	UG/M3	14.3 U	1.43 U	1.55 U	14.3 U	13.5 U
1,1-Dichloroethane	UG/M3	14.6 U	1.46 U	1.58 U	14.6 U	13.8 U
2-Butanone	UG/M3	26.8 U	2.57 J	2.83 U	26.5 U	25.1 U
Chloroform	UG/M3	17.6 U	1.76 U	1.90 U	17.6 U	39.1
1.2-Dichloroethane	UG/M3	14.6 U	1.46 U	1.58 U	14.6 U	13.8 U
1,1,1-Trichloroethane	UG/M3	19.6 U	1.96 U	2.13 U	19.6 U	18.6 U
Carbon Tetrachloride	UG/M3	22.7 U	2.27 U	2.45 U	22.7 U	21.4 U
Bromodichloromethane	UG/M3	24.1 U	2.41 U	2.61 U	24.1 U	22.8 U
1,2-Dichloropropane	UG/M3	16.6 U	1.66 U	1.80 U	16.6 U	15.7 U
cis-1,3-Dichloropropene	UG/M3	16.3 U	1.63 U	1.77 U	16.3 U	15.4 U
Trichloroethene	UG/M3	19.4 U	1.93 U	2.10 U	19.3 U	12.4 J
Benzene	UG/M3	11.5 U	0.89 J	1.25 U	24.0	10.9 U
Dibromochloromethane	UG/M3	30.7 U	3.07 U	3.32 U	30.7 U	29.0 U
trans-1,3-Dichloropropene	UG/M3	16.3 U	1.63 U	1.77 U	16.3 U	15.4 U
1,1,2-Trichloroethane	UG/M3	19.6 U	1.96 U	2.13 U	19.6 U	18.6 U
Bromoform	UG/M3	37.2 U	3.72 U	4.03 U	37.2 U	35.1 U

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Location ID Sample ID		SG-06	SG-06	SG-07	SG-08	SG-09
		FD-2	SG-6	SG-7	SG-8	SG-9
Matrix		Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas
Depth Interval (ft)		· ·		-		•
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units	Field Duplicate (1-1)				
Volatile Organic Compounds						<u></u>
4-Methyl-2-Pentanone	UG/M3	37.3 U	2.29 J	3.93 U	36.9 U	34.8 U
2-Hexanone	UG/M3	37.3 U	3.65 U	3.93 U	36.9 U	34.8 U
Tetrachloroethene	UG/M3	24.4 U	3.26	3.80 J	24.4 U	35.3
1,1,2,2-Tetrachioroethane	UG/M3	24.7 U	2.47 U	2.68 U	24.7 U	23.3 U
Toluene	UG/M3	6.41 J	6.03	6.03 J	136	12.8
Chlorobenzene	UG/M3	16.6 U	1.66 U	1.80 U	16.6 U	15.7 U
Ethylbenzene	UG/M3	15.6 U	2.00	1.65 J	14.3 J	14.8 U
Styrene	UG/M3	15.3 U	1.15 J	1.66 U	15.3 U	14.5 U
m,p-Xylene	UG/M3	31.3 U	13.0	10.4 J	80.8	15.6 J
o-Xylene	UG/M3	15.6 U	2.78	1.95 J	13.9 J	14.8 U
cis-1,2-Dichloroethene	UG/M3	14.3 U	1.43 U	1.55 U	14.3 U	13.5 U
trans-1,2-Dichloroethene	UG/M3	14.3 U	1.43 U	1.55 U	14.3 U	13.5 U

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Location ID		SG-10	SG-10	SG-11	SG-12	SG-13
Sample ID		FD-1	SG-10	\$G-11	SG-12	SG-13
Matrix		Soil Gas	Soil Gas	Soil Gas	Soll Gas	Soil Gas
Depth Interval (ft)		•	•	-	•	•
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units	Field Duplicate (1-1)				
Volatile Organic Compounds	-					
Chloromethane	UG/M3	18.6 U	19.8 U	18.2 U	0.58 J	18.2 UJ
Bromomethane	UG/M3	14.0 U	14.8 U	13.6 U	1.40 U	13.6 U
Vinyl Chloride	UG/M3	9.20 U	9.71 U	8.95 U	0.92 U	8.95 UJ
Chloroethane	UG/M3	9.50 U	10.0 U	9.24 U	0.95 U	9.24 U
Methylene Chloride	UG/M3	31.3 U	17.0 J	25.4 J	3.13 U	30.6 UJ
Acetone	UG/M3	214 U	228 U	209 U	28.5 U	209 UJ
Carbon Disulfide	UG/M3	15.6	11.2 J	9.65 J	15.0	7.47 J
1,1-Dichloroethene	UG/M3	14.3 U	15.1 U	13.9 U	1.43 U	13.9 UJ
1,1-Dichloroethane	UG/M3	14.6 U	15.4 U	14.2 U	1.46 U	14.2 U
2-Butanone	UG/M3	26.5 U	28.3 U	26.0 U	2.86	26.0 U
Chloroform	UG/M3	14.2 J	8.79 J	17.1 U	7.81	17.1 U
1,2-Dichloroethane	UG/M3	14.6 U	15.4 U	14.2 U	1.46 U	14.2 U
1,1,1-Trichloroethane	UG/M3	19.6 U	20.7 U	19.1 U	0.60 J	19.1 U
Carbon Tetrachloride	UG/M3	22.7 U	23.9 U	22.0 U	2.27 U	22.0 U
Bromodichloromethane	UG/M3	24.1 U	25.5 U	23.4 U	2.41 U	23.5 U
1,2-Dichloropropane	UG/M3	16.7 U	17.6 U	16.2 U	1.66 U	16.2 U
cis-1,3-Dichloropropene	UG/M3	16.3 U	17.2 U	15.9 U	1.63 U	15.9 U
Trichloroethene	UG/M3	19.4 U	20.4 U	18.8 U	1.93 U	18.8 U
Benzene	UG/M3	11.5 U	12.1 U	4.79 J	1.89	11.2 U
Dibromochloromethane	UG/M3	30.7 U	32.4 U	29.8 U	3.07 U	29.8 U
trans-1,3-Dichloropropene	UG/M3	16.3 U	17.3 U	15.9 U	1.63 UJ	15.9 U
1,1,2-Trichloroethane	UG/M3	19.6 U	20.7 U	19.1 U	1.96 U	19.1 U
Bromoform	UG/M3	37.2 U	39.3 U	36.2 U	3.72 UJ	36.2 U

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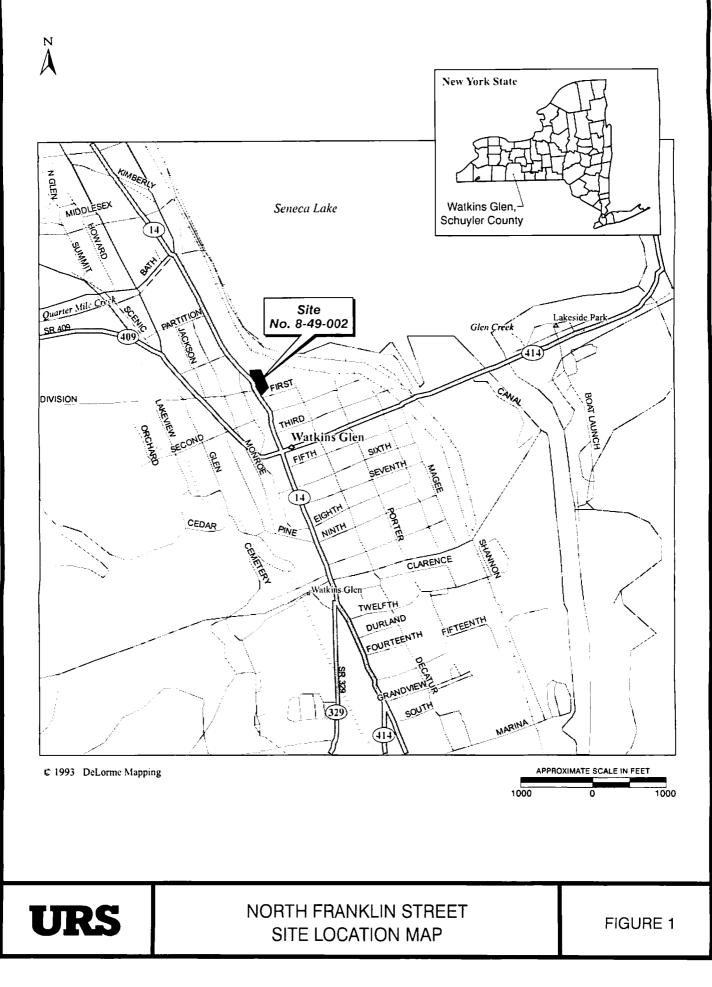
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Location ID	· · · ·	SG-10	SG-10	SG-11	SG-12	SG-13
Sample ID		FD-1	SG-10	SG-11	SG-12	SG-13
Matrix		Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas
Depth Interval (ft)			•	•		
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units	Field Duplicate (1-1)				
Volatile Organic Compounds						
4-Methyl-2-Pentanone	UG/M3	36.9 U	39.3 U	36.1 U	1.97 J	36.1 U
2-Hexanone	UG/M3	36.9 U	39.3 U	36.1 U	3.69 U	36.1 U
Tetrachloroethene	UG/M3	27.8	25.8 U	23.7 U	8.14	23.7 U
1,1,2.2-Tetrachloroethane	UG/M3	24.7 U	26.1 U	24.0 U	2.47 U	24.0 U
Toluene	UG/M3	10.9 J	7.54 J	18.5	15.8	15.1
Chlorobenzene	UG/M3	16.6 U	17.5 U	16.1 U	1.66 U	16.1 U
Ethylbenzene	UG/M3	15.6 U	16.5 U	7.38 J	1.74	15.2 U
Styrene	UG/M3	15.3 U	- 16.2 U	14.9 U	1.53 UJ	14.9 U
m,p-Xylene	UG/M3	37.3	18.2 J	51.2	6.77	15.6 J
o-Xylene	UG/M3	6.95 J	16.5 U	8.25 J	1.26 J	15.2 U
cis-1,2-Dichloroethene	UG/M3	14.3 U	15.1 U	13.9 U	1.43 U	13.9 U
trans-1,2-Dichloroethene	UG/M3	14.3 U	15.1 U	13.9 U	1.43 U	13.9 U

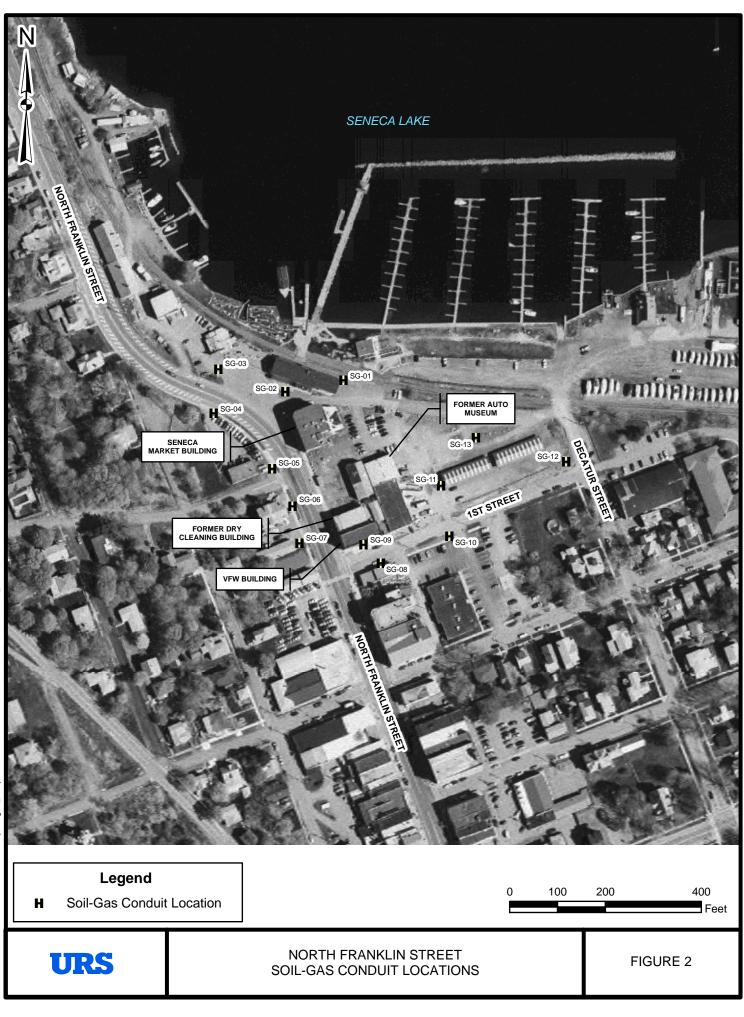
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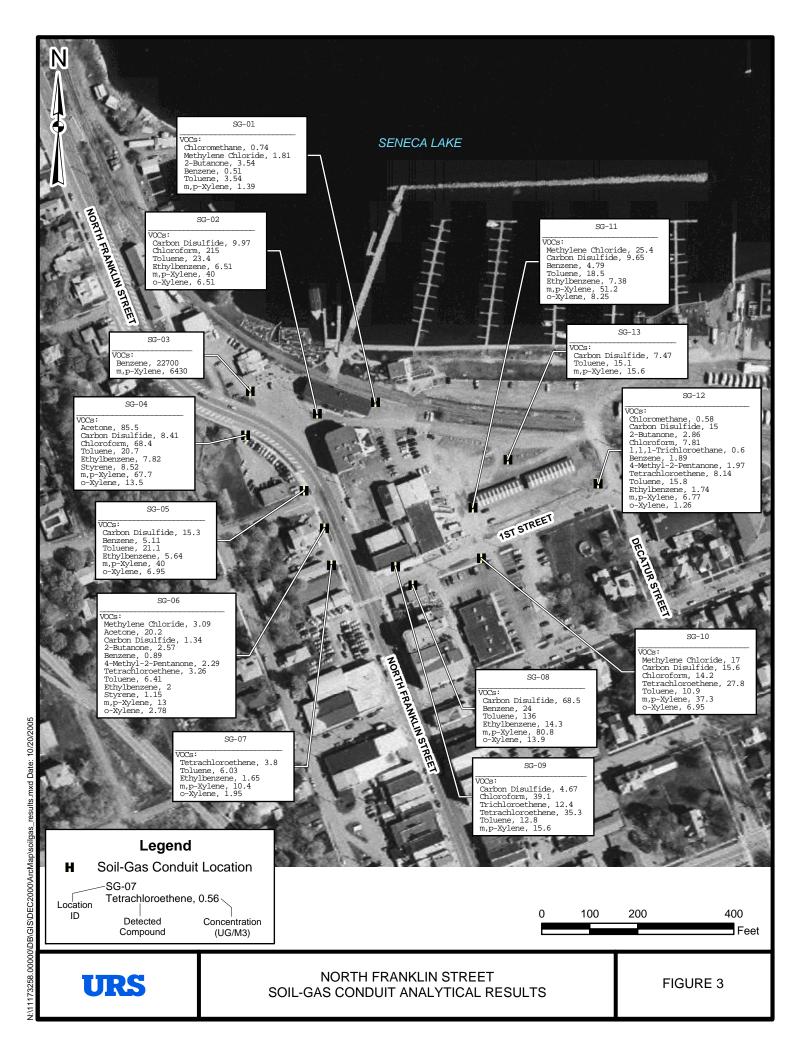
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FIGURES



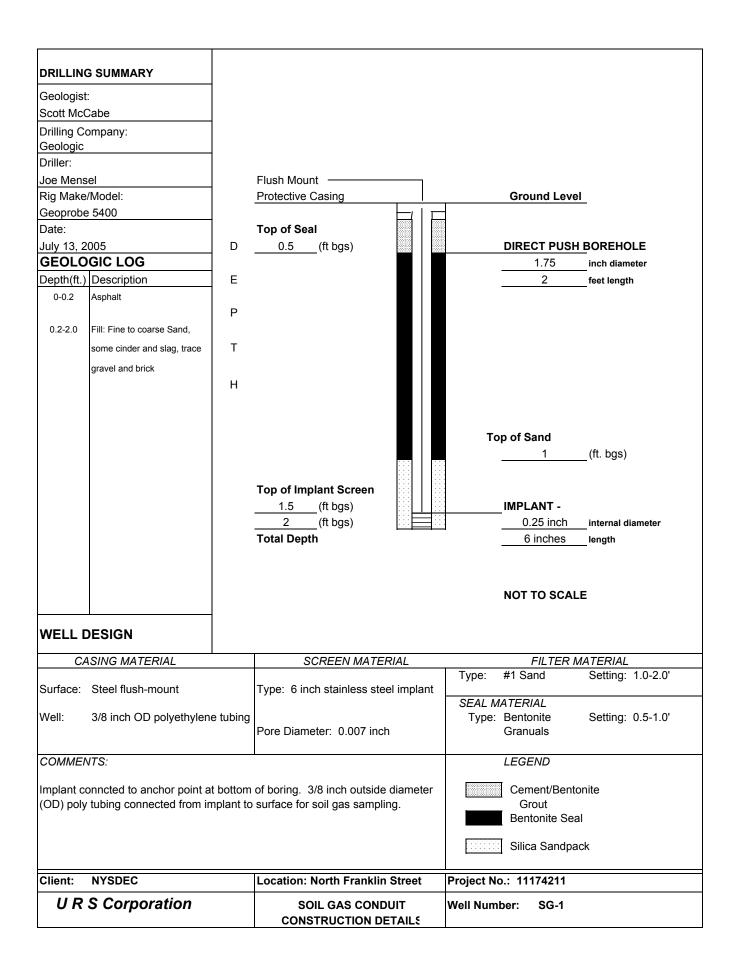
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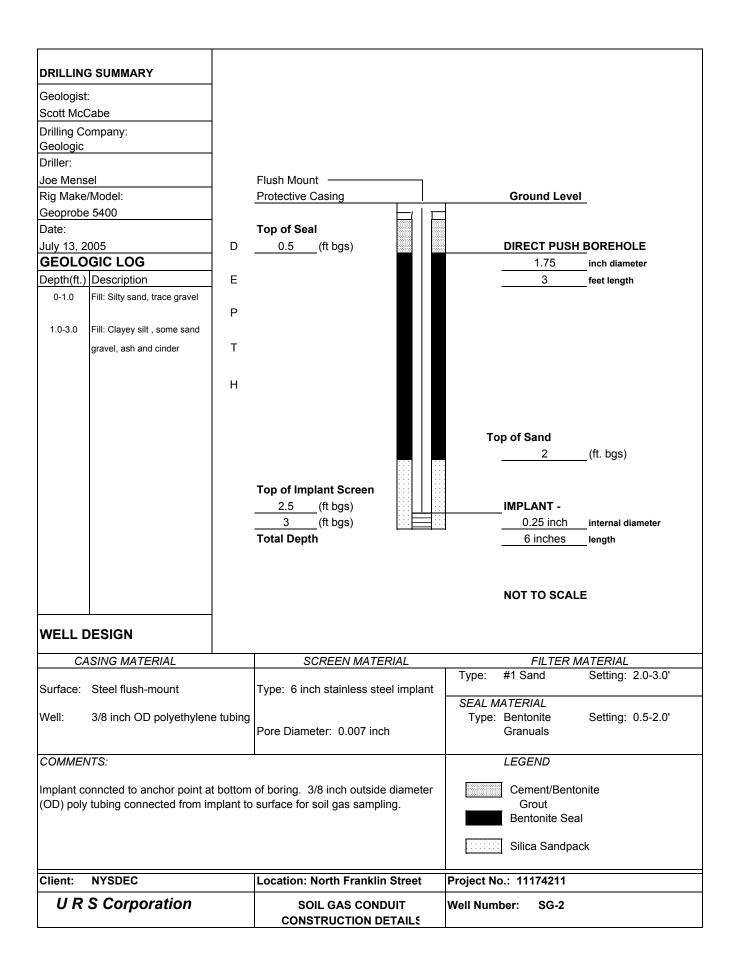


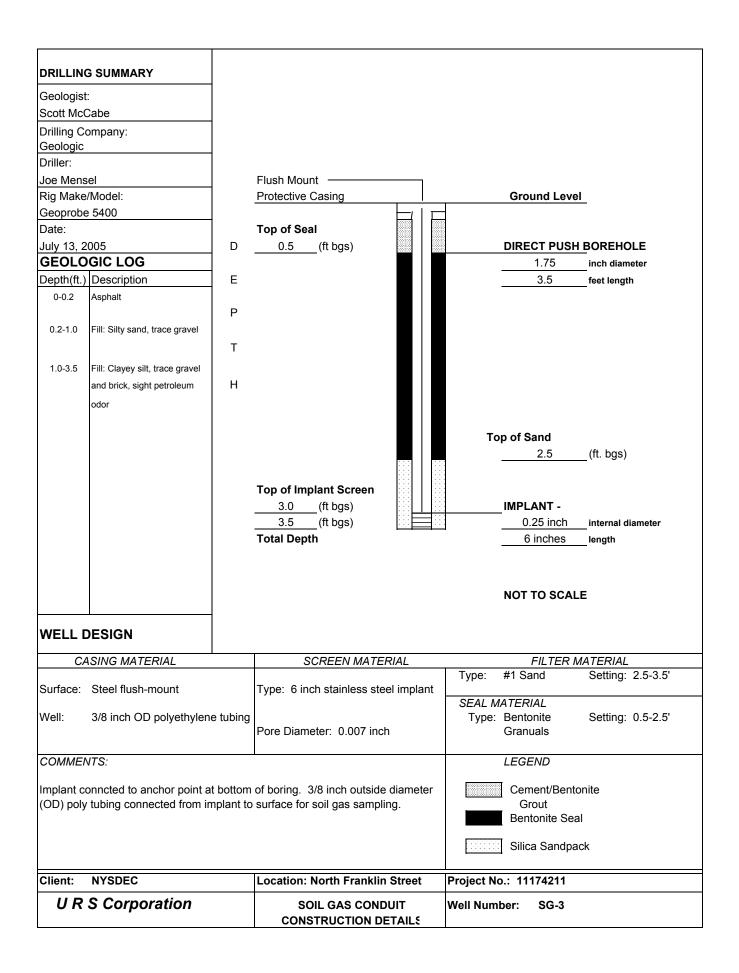


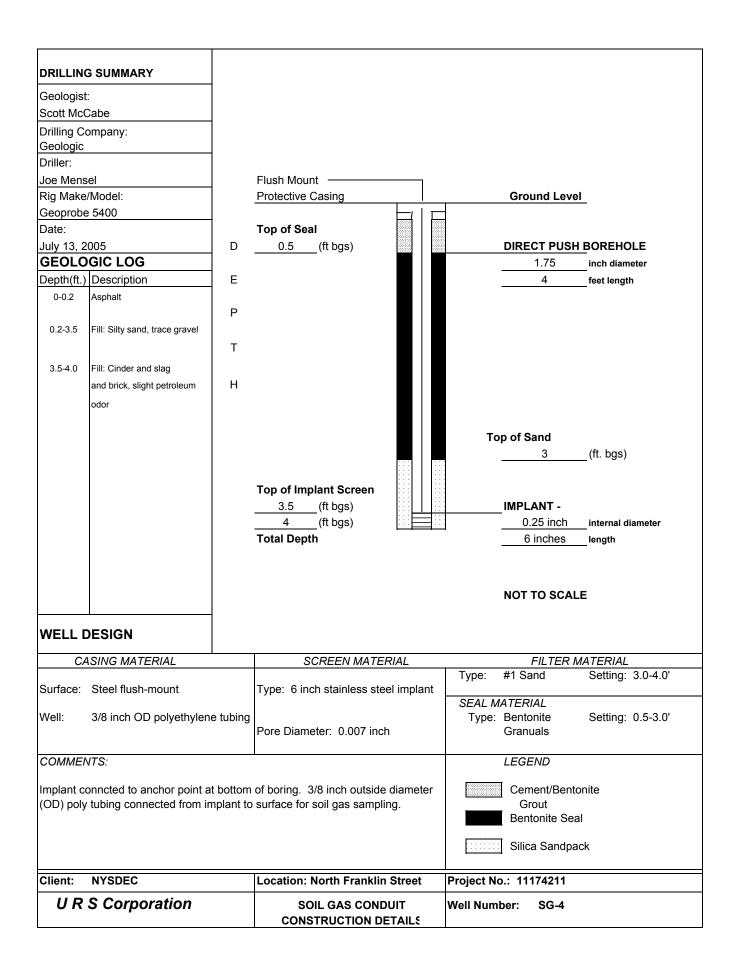
ATTACHMENT 1

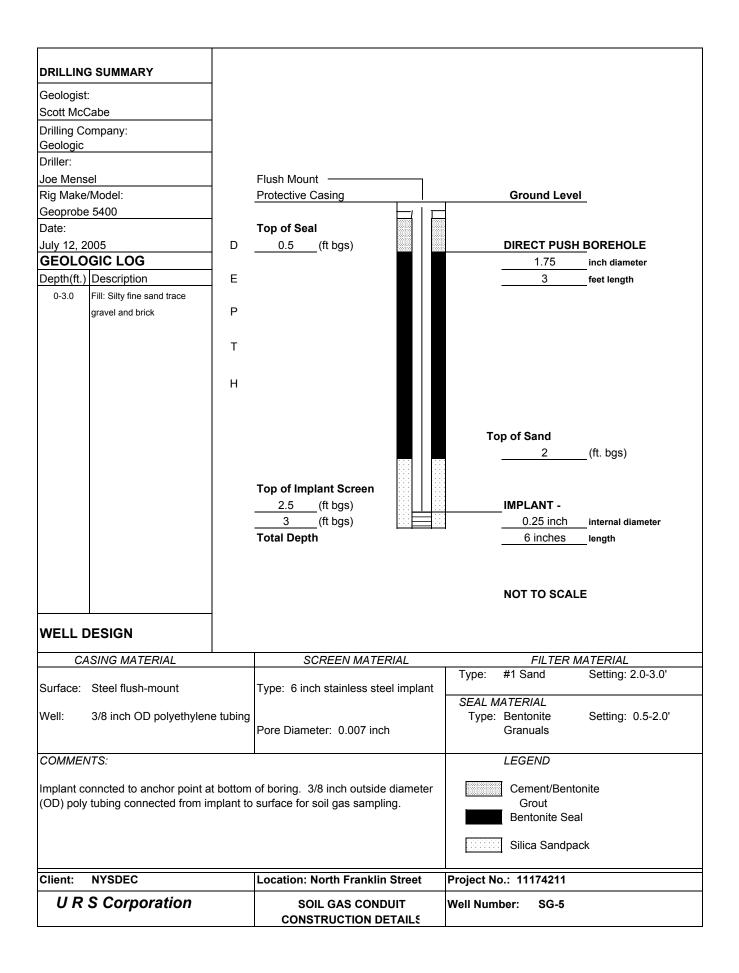
SOIL-GAS CONDUIT CONSTRUCTION DETAILS

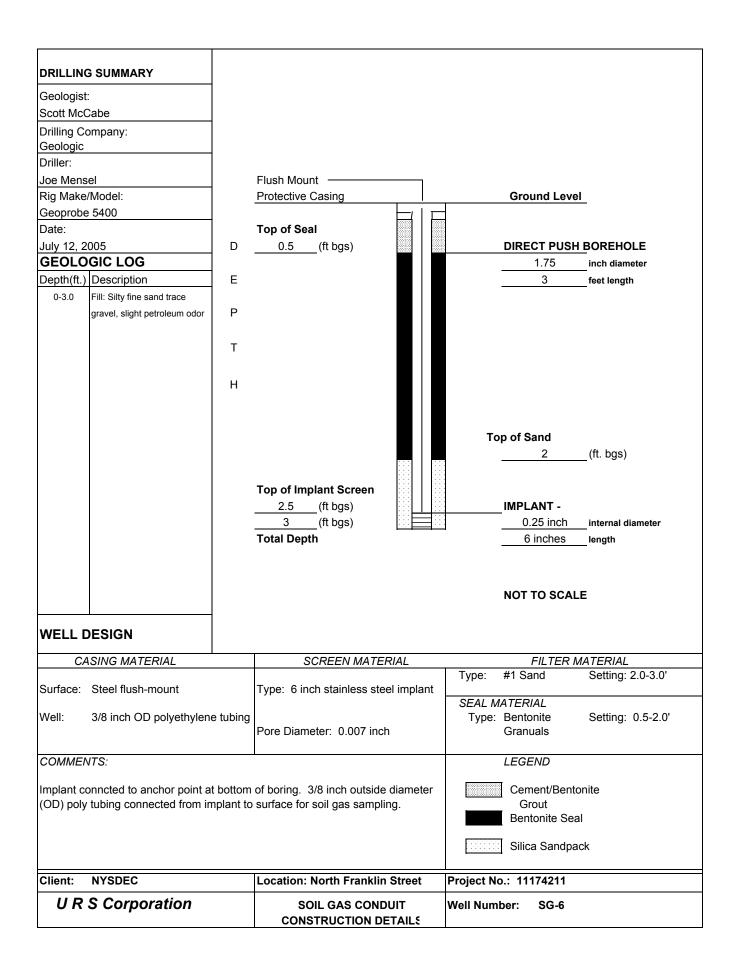


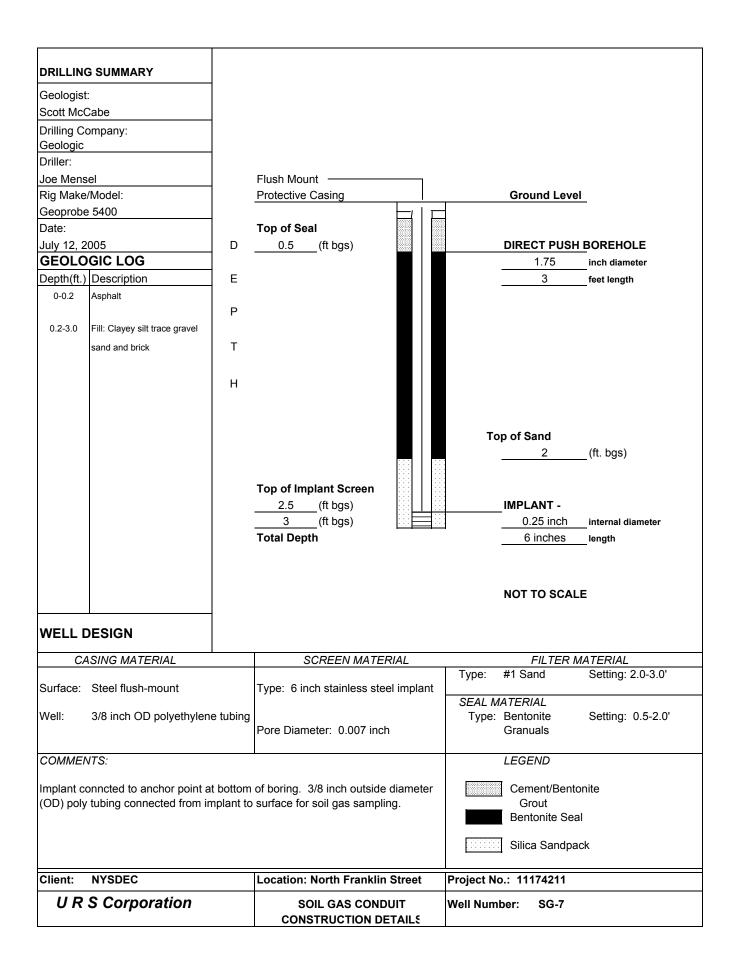


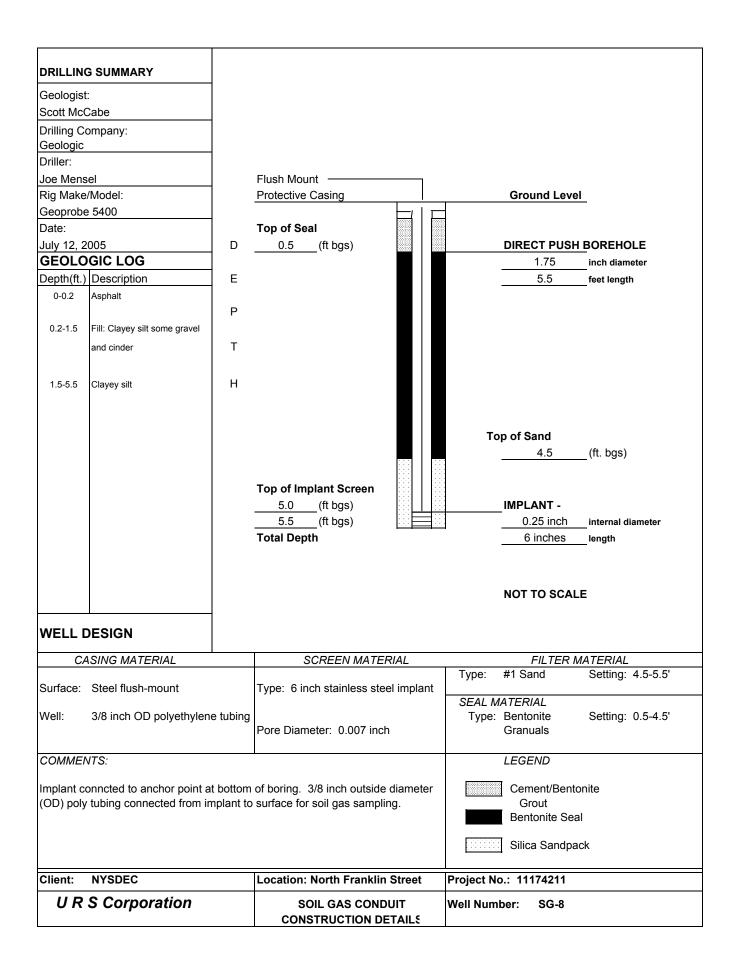


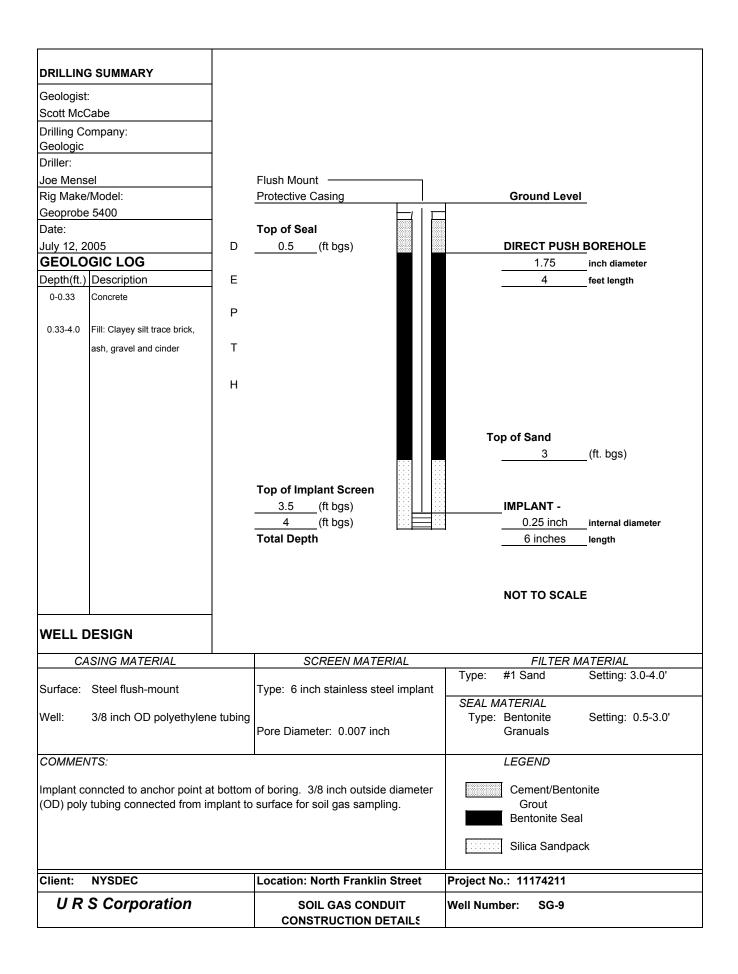


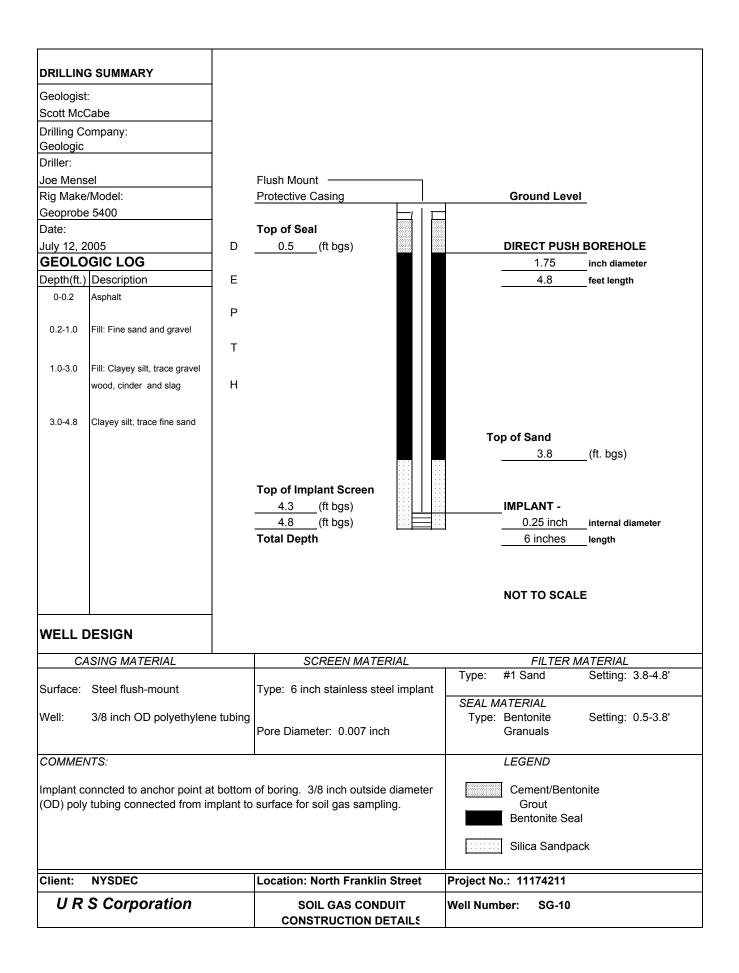


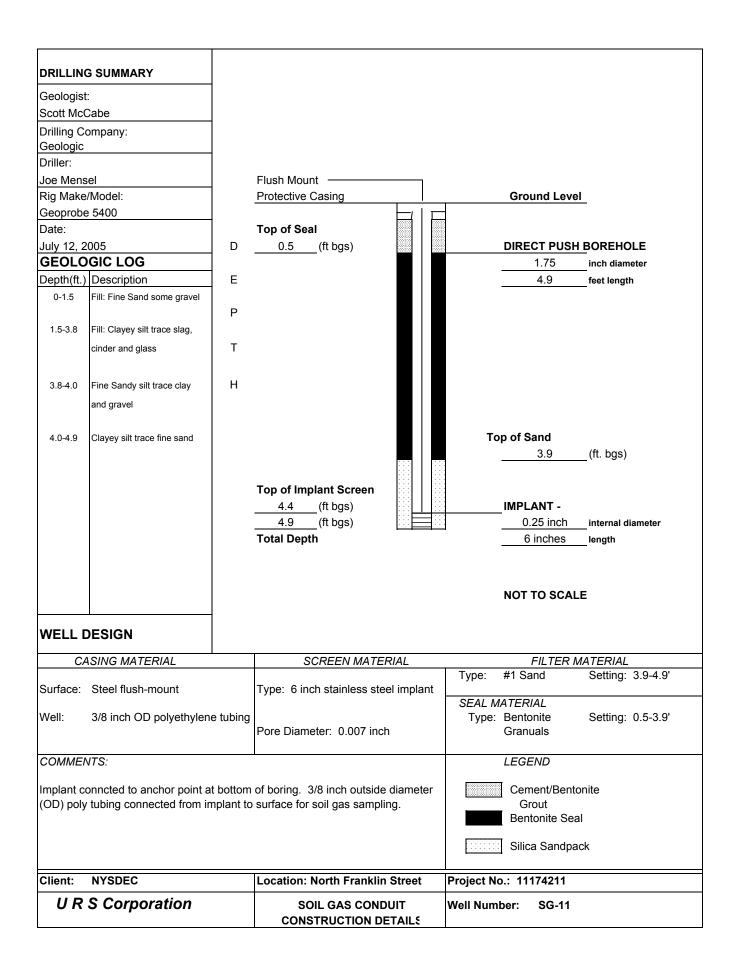


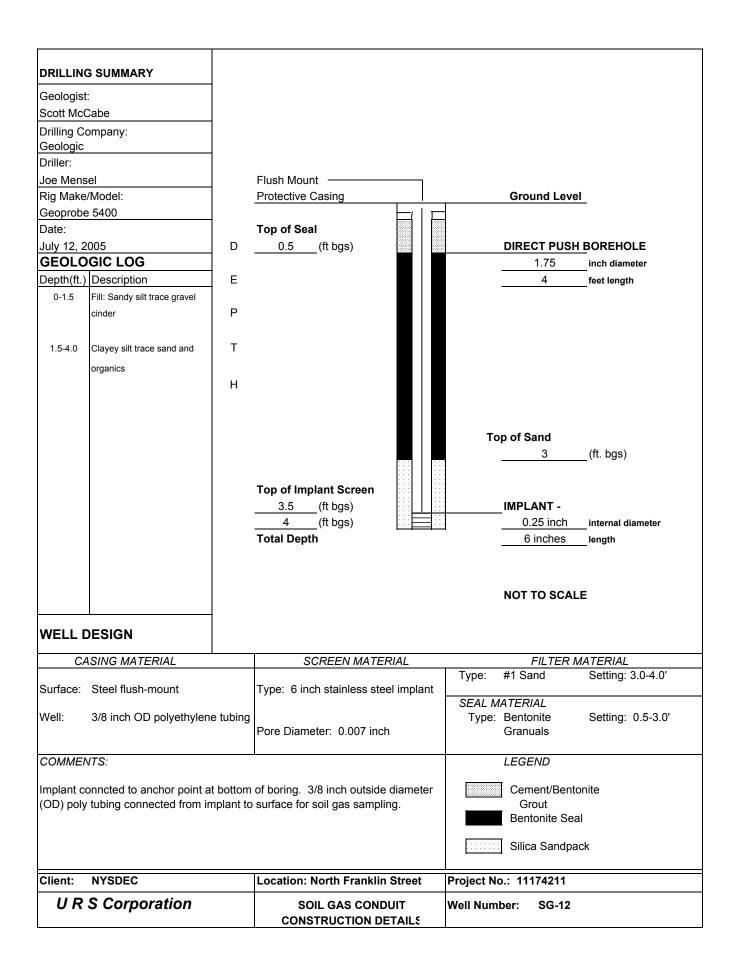


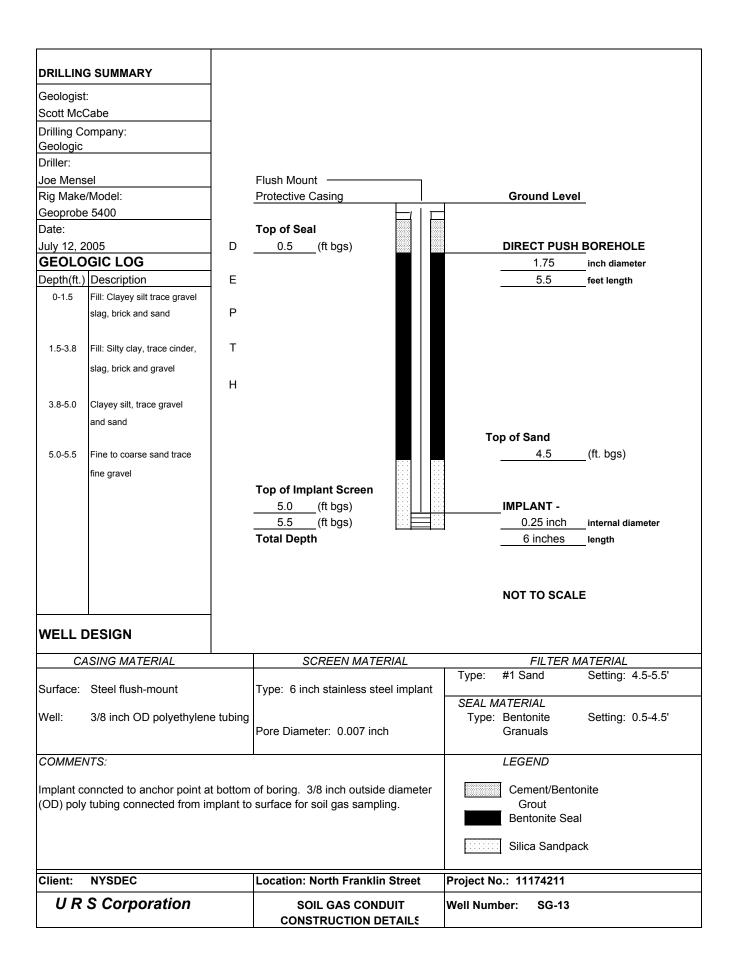












ATTACHMENT 2

PHOTOGRAPHIC LOG



Photo 1: Photograph of a soil gas implant.



Photo 2: Geoprobe operators attaching polyethylene tubing to the soil gas implant.





Photo 3: Photograph of installed soil gas implant prior to installation of bentonite seal and protective casing.



Photo 4: Photograph looking north at SG-1.





Photo 5: Photograph looking north at SG-2.





Photo 6: Photograph looking north at SG-3.

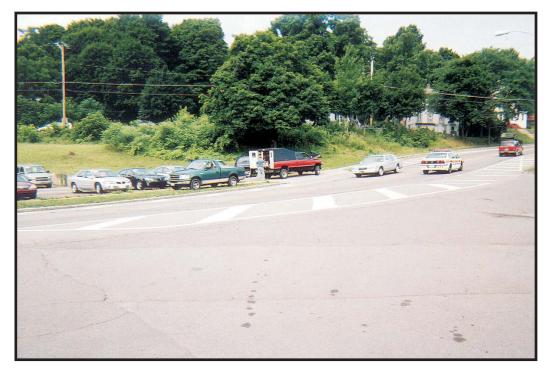


Photo 7: Photograph looking north at SG-4.





Photo 8: Photograph looking west at SG-4.





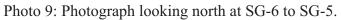






Photo 10: Photograph looking west at SG-6.





Photo 11: Photograph looking west at SG-7.





Photo 12: Photograph looking south at SG-8.



Photo 13: Photograph looking north at SG-9.





Photo 14: Photograph looking south at SG-10.





Photo 15: Photograph looking north at SG-11.





Photo 16: Photograph looking south at SG-12.





Photo 17: Photograph looking south at SG-17.





Photo 18: Photograph looking northwest at tracer gas and summa canister set-up at SG-1 location.



Photo 19: Photograph looking northwest at tracer gas and summa canister set-up at SG-2 location.





Photo 20: Photograph looking west at tracer gas and summa canister set-up at SG-3 location.



Photo 21: Photograph looking west at tracer gas and summa canister set-up at SG-4 location.





Photo 22: Photograph looking west at tracer gas and summa canister set-up at SG-5 location.



Photo 23: Photograph looking south at tracer gas and summa canister set-up at SG-6 location.





Photo 24: Photograph looking southwest at tracer gas and summa canister set-up at SG-7 location.



Photo 25: Photograph looking northwest at tracer gas and summa canister set-up at SG-8 location.





Photo 26: Photograph looking north at tracer gas and summa canister set-up at SG-9 location.



Photo 27: Photograph looking south at tracer gas and summa canister set-up at SG-10 location.





Photo 28: Photograph looking north at tracer gas and summa canister set-up at SG-11 location.





Photo 29: Photograph looking north at tracer gas and summa canister set-up at SG-12 location.



Photo 30: Photograph looking south at tracer gas and summa canister set-up at SG-13 location.



ATTACHMENT 3

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SAMPLING LOGS AND CHAIN-OF-CUSTODY

Site:	North	FRANKLIN STRIFFET
Samp	1	S. MULABIZ
Date:	7/1	

Sample #	56-3 3,0-3,5	56-2 25-30	59-11.5-2.0	20050718-198-1	
Location	56-3	56-2	56-1	AB-1	
Summa Canister ID (Lab ID, if provided)	1495	04311	12738	9 30 83	
Additional Tubing Added		NO/		YES - How much	NO/ YES - How muc
	2	2	2'	-	
Purge Time (Start)	1018	1638	1658		
Purge Time (Stop)	1023	1643	1103		
Total Purge Time (min)	5 min	5 min	5 min	_	
Pressure Gauge - before sampling	- 29	-30	- 29	-29.5	
Sample Time (Start)	1023	1643	1103	1109	
Sample Time (Stop)	1120	1138	1158	1204	
otal Sample Time (min)	- 57	55	55	55	
ressure Gauge - after ampling	-10,0	-8.0	-8.0	-8.0	
anister Pressure Went o Ambient Pressure?	YES /	YES INO	YES NO	YES NO	YES / NO
eneral Comments:	- 2		······		
<u>56-7 30-3,5 (</u> <u>56-2 2,5-3,0 (</u>	DE HE DETECTOR	157pp	nin pine	Of HE AFTE	& sample
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	NE WICK VETCI	Jee pp	D In porco	· UD °L it iz Air	En Supl.

Site: North Frankin STREET. Samplers: S. Microbie Date: 7/18/05

Sample #	56-12 3.5-4.0	56-13 50-5.5	556-11 4.4-4.9	56-10 4.3-4.8	20050718-FD-1
Location	56-12	56-13	56-11	56-10	FD-1 (56-10)
Summa Canister ID (Lab ID, if provided)	04396	3402	9 3074	5/501	1787
Additional Tubing Added	NO/ YES- How much Z ¹	NO/ ES How much Z'	NO/ How much	NO/ TES How much	r
Purge Time (Start)	1222	1239	1254	1309	2'
Purge Time (Stop)	1227	1244	1259	1314	
Total Purge Time (min)	5 min	5 min	5 min	5 mins	1314 5 min
Pressure Gauge - before sampling	-30	-30	-30	~ 36	- 30
Sample Time (Start)	1227	1244	1259	1314	1314
Sample Time (Stop)	1322	1339	1354	1409	1409
otal Sample Time (min)	55	55	55	55	55
ressure Gauge - after ampling	_9	-12	- 10	-10	-/6
anister Pressure Went o Ambient Pressure?	YES INO	YES /	YES /	YESNO	YES (NO)
eneral Comments:		· · · · · · · · · · · · · · · · · · ·			
56-12 0% HE 56-13 0% HE	Detected	174Z pp	0% H	12 ATTER SAMP	0
56 -11 01/0 Hiz	Detected	162 2 ppb 1048 ppb	0% H	L AFRY Sample	P
56-10 0º/6/42	letated	132 (peb	<u> </u>	A CONTRACTOR OF THE OWNER	A REAL PROPERTY AND A REAL PROPERTY A DATACASE IN
FD-1 0% Hoz	Detected	1321 100	0% 14	the second se	1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
	•			J	

Samplers: S. m Date: 7/18/05	CCABIE -				
Sample #	56-4 35-4.0	56-5 2.5.3.0	36-6 2.5.30	20050718-FP-2	
Location	56-4	565	56-6	FD-Z (56-6)	
Summa Canister ID (Lab ID, if provided)	1127	1530	93265	51529	
Additional Tubing Added	NO/ (TES)- How much 2'	NO/ How much	NO/ Tes How much	NO/ YES- How much Z /	NO/ YES - How muc
Purge Time (Start)	1555	1540	14/55	1455	
Purge Time (Stop)	1600	1545	1500	1500	
Total Purge Time (min)	5 mini	5 min	5 min	5 min	
Pressure Gauge - before sampling	-29	- 30	- 22	- 28.5	
Sample Time (Start)	1600	1545	1500	1500	· · ·
Sample Time (Stop)	1655	1640	1555	1555	· · · · · · · · · · · · · · · · · · ·
otal Sample Time (min)	55	55	55	55	
ressure Gauge - after ampling	-9	-6	-9	~11	
anister Pressure Went o Ambient Pressure?	YES / 10	YES / 10	YES/NO	YES/NO	YES / NO
eneral Comments:	······				
56-4 0% HB 5	Viecteo	361,000	00/ 11	as a mar	
6-5 042 149 0	ketecked	563 pp b	0%14	E AFTEN SAPE	volc.
	Detected	<u></u>	- O'LI	Contraction of the local division of the loc	ne
D-Z DYUHB	Derecter)	791 006	0%,14	R AFTER SH	

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Samplers: S. m Date: 7/18	LCABE 105				
Sample #	SG-8 5.0-5.5	567-9 3.5-4.0	56-7 2.5-3.0		
Location	56-8	56-9	56-7		
Summa Canister ID (Lab ID, if provided)	2995	51488	51494		
Additional Tubing Added	NO/ How much	NO/ TES How much	NO/ TES How much	NO/ YES - How much	NO/ YES - How much
Purge Time (Start)	1415	1425	1435		
Purge Time (Stop)	1420	1430	1440		
Total Purge Time (min)	5 min	5 min	5 min		
Pressure Gauge - before sampling	-30	-29.5	- 29		
Sample Time (Start)	1420	1430	1440		
Sample Time (Stop)	1515	1525	1535		
otal Sample Time (min)	55	55	55		
ressure Gauge - after ampling	-11	-12	-9		
anister Pressure Went o Ambient Pressure?	YES / NO	YES / 10	YES NO	YES / NO	YES / NO
eneral Comments: - 4 - 8 0% H7 6 - 9 0% H7 6 - 9 0% H7 6 - 7 0% H7		1427, 1038, 695,	226 · 0 ⁴ b 0 ⁵ b 0 ⁵	6 HE AFTE	Sample

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6-2		1138			2.5-3.0	65							<u>+</u>			1495	- 11	$\left\{ - \right\}$	<u> </u>	
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-13		1339	1		5.0-5.5	65		1	+		<u> </u>					04396	n/			
n-11		1354				65		$\frac{1}{1}$		Sec.						3402	n			
5-10		1409		56-10	4.3-4.8	65		$\frac{1}{2}$								93674	r1		-	54554
-8		1515	IhR	56-8	5.0-5.5	65	 									51501	01			
5-9		1525	the	56-9	3,5-4,0	GS	1	,								2995	n			<u> </u>
-7		1535		56-7	2.5-3.0	GS	1	,								51488	NI			·
-4		1655		56-4	3.5-4.0	65	 	-1	ar Esta	H.	·	•				51494	WI			
-5	V	1640	thR	56-5	25-3.0	10 March 10	1	1			*	84				1127	NI			
MATRIX CODES	AA - AMBIE SE - SEDIN SH - HAZA		ASTE	SL - SLUDGE WP - DRINKIN WW - WASTE	IG WATER S	VG - GROUND O - SOIL OC - DRILL CU	-	(ml - Lea GS - Soil WC - Dri	GAS		. N	NO - Oce NS - Sur NQ - Wat	FACE W	ATER	LH - HAZARDOUS LI LF - FLOATING/FREE	IQUID WAS	TE TON GW		
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56-6	7/18/05	1555	1. 1	S6-6	2,5-3.0	65	· 1	<u> ŀ</u>			ļ	ļ	<u> </u>	 		93265	M			
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MATRIX CODES		IENT RDOUS SOLID W		sl - Sludge WP - Drinkin WW - Waste	G WATER SO	i - Ground - Soil - Drill Cu		G	S - SO	ACHATE IL GAS RILLING W	/Ater	1	WO - OCE WS - SUF WQ - WAT	FACE W	ATER	LH - HAZARDOUS LIQ LF - FLOATING/FREE F	UID WAS PRODUCT	TE ON GV		L SLE
SAMPLE TYPE CODES	TB# - TRIP SD# - MAT	BLANK RIX SPIKE DUPL		RB# - RINSE E FR# - FIELD F		- NORMAL # - MATRIX	ENVIRONN SPIKE	IENTAL S	AMPLE	(# -	SEQUE		JMBER (F	ROM 1 1	O 9) TO /	ACCOMMODATE MULTIPLE	SAMPLES			-
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RELINQUISH		NATURE)	DATE		RECEIVED FO	DR LAB	BY (sig	NATURI	E)	DATE	TIN	ИЕ								

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ATTACHMENT 4

DATA USABILITY SUMMARY REPORT

DATA USABILITY SUMMARY REPORT

NORTH FRANKLIN STREET SITE SITE NO. 84-90-002 WORK ASSIGNMENT D003825-09.3

Analyses Performed by:

SEVERN-TRENT LABORATORIES (STL) – KNOXVILLE, TN

Prepared by:

URS CORPORATION 77 GOODELL STREET BUFFALO, NY 14203

SEPTEMBER 2005

N:\11173258.00000\WORD\North Franklin Street Air July 2005 DUSR.doc

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I.	INTRODUCTION	1
II.	ANALYTICAL METHODOLOGIES	1
III.	DATA DELIVERABLE COMPLETENESS	1
IV.	HOLDING TIMES/SAMPLE RECEIPT	1
V.	NON CONFORMANCES	2
VII.	SUMMARY	4

TABLES (Following Text)

Table 1	Summary of Data Qualifications
Table 2	Validated Soil Gas and Ambient Air Analytical Results

APPENDICES

Appendix A	Support Documentation
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Appendix B Validated Form I's

I. INTRODUCTION

This Data Usability Summary Report (DUSR) has been prepared following the guidelines provided in New York State Department of Environmental Conservation (NYSDEC) Division of Environmental Remediation *Guidance for the Development of Data Usability Summary Reports (DUSR)*, dated June 1999. This DUSR discusses the soil gas and ambient air samples collected on July 18, 2005.

II. ANALYTICAL METHODOLOGIES

The data being evaluated is from the July 18, 2005 sampling of 13 soil gas samples, 1 ambient air sample, and 2 field duplicate samples. The analytical laboratory that performed the analyses is Severn-Trent Laboratories, Inc. (STL) located in Knoxville, TN (STL-Knoxville). The samples were analyzed for volatile organic compounds (VOCs) following USEPA Compendium Method TO-15, *Determination of VOCs in Air Collected in Specially Prepared Canisters and Analyzed By Gas Chromatography/Mass Spectrometry (GC/MS)*.

Table 1 summarizes the data qualifications applied to the sample results. The validated analytical results are presented on Table 2.

A limited data validation was performed following the guidelines in USEPA Region II *Validating Canisters of Volatile Organics in Ambient Air, Rev. 0*, April 1994. Qualifications applied to the data include 'J/UJ' (estimated concentration/estimated quantitation limit) and 'U' (not detected). Documentation supporting the qualification of data is presented in Appendix A. Copies of the validated laboratory results (i.e., Form I's) are presented in Appendix B. Only problems affecting data usability are discussed in this report.

III. DATA DELIVERABLE COMPLETENESS

The laboratory deliverable data package was equivalent to NYSDEC Analytical Services Protocol (ASP) Category B requirements.

IV. HOLDING TIMES/SAMPLE RECEIPT

All samples were received by the laboratory intact and analyzed within the required holding time.

V. NON CONFORMANCES

Quality Control (QC) Blanks

The ambient air sample is considered to be representative of ambient air conditions. Since the soil gas samples are collected approximately 8 feet below ground surface through soil gas conduits, and the ambient air sample represents the air in the breathing zone, the results of the ambient air sample have not been used to qualify the soil gas sample results.

The laboratory method blanks associated with the soil gas samples had detections for acetone. In accordance with the USEPA Region II validation guidelines, those soil gas samples exhibiting detected results for acetone less than five times the concentration detected in the associated laboratory method blank have been qualified as follows: If the value detected in the sample was below the reporting limit (RL) then the compound was raised to the reporting limit and qualified as 'U' (undetected). If the value detected in the sample was greater than the RL, but less than fives times the blank value, the compound was qualified 'U' at the value detected. The affected samples have been listed on Table 1.

Documentation supporting the qualification of data (e.g. method blank Form Is) is presented in Appendix A.

Initial and Continuing Calibrations

The percent difference (%D) between the initial calibration (ICAL) average relative response factor (RRF) and the RRF in one of the continuing calibration (CCAL) standards exceeded the QC limit (i.e., >25%D) for bromoform, trans-1,3-dichloropropene, and styrene. The results for these compounds in the associated samples listed on Table 1 have been qualified 'J' or 'UJ.'

The %D between the ICAL average RRF and the RRF in one of the CCAL standards exceeded the QC limit (i.e., >25%D) for acetone, carbon disulfide, chloromethane, 1,1-dichloroethene, methylene chloride, and/or vinyl chloride. The results for these compounds in the associated samples listed on Table 1 have been qualified 'J' or 'UJ.'

Documentation supporting the qualification of data (e.g., Autosampler Run log, Continuing Calibration Form) is presented in Appendix A.

Surrogates/Internal Standards

The recovery of surrogate 1,2-dichloroethane-d4 in the initial analysis of sample SG-7 was greater than the upper QC limit. The sample was re-analyzed at a dilution and exhibited acceptable recoveries for all compounds. However because the re-analysis was performed at a dilution the RLs were elevated for all compounds. Using professional judgment the results of the initial analysis have been reported on Table 2. The detected results for ethylbenzene, tetrachloroethene, toluene, o-xylene, and m/p-xylene in this sample have been qualified 'J' due to the surrogate outlier.

Target Compounds/Instrument Performance

The following samples were analyzed utilizing a dilution due to elevated concentrations of target compounds and/or matrix interference: SG-2, SG-3, SG-4, SG-5, FD-2 (SG-6), SG-8, SG-9, SG-10, FD-1 (SG-10), SG-11, SG-13, and AB-1. The reporting limits for the non-detect compounds represent the lowest achievable at the diluted level. Samples AB-1, SG-13, and FD-1 (SG-10) were initially analyzed undiluted. These samples exhibited no recovery of the surrogates and internal standards and very poor chromatography. Since two different instruments were used for the analysis of these samples and the chromatography problems were observed on both instruments, the laboratory believes that sample matrix negatively affected instrument performance. The laboratory indicated in the report narrative that the chromatography problems were caused by the freezing of the instrument concentrator traps which restricted the flow of sample to the GC/MS. This could have resulted from water and/or carbon dioxide in the samples causing the trap to freeze-up. The samples were re-analyzed at dilutions and all surrogate and internal standard recoveries were acceptable. The laboratory determined that due to the potential for column and instrument damage by analyzing the remainder of the samples listed above un-diluted, these remaining samples were also analyzed at diluted levels.

VII. SUMMARY

All sample analyses were found to be compliant with the method criteria, except where previously noted. Those results qualified 'J/UJ' (estimated/estimated reporting limit) are considered conditionally usable. Those results qualified 'U' are considered non-detect. All other sample results are usable as reported. URS does not recommend the re-collection of any samples at this time.

Prepared By: Ann Marie Kropovitch, Chemist **Date:** 9|13|05**Date:** 9|13|05Reviewed By: James J. Lehnen, Senior Chemist

DEFINITIONS OF USEPA REGION II DATA QUALIFIERS

- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- J The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
- UJ The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
- R The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.
- D The sample results are reported from a separate secondary dilution analysis.
- NJ Presumptive evidence of a compound at an estimated value.

SAMPLE ID	FRACTION	ANALYTICAL DEVIATION	QUALIFICATION
AB-1, FD-1, FD-2, SG-2, SG-5, SG-7, SG-9, SG-10, SG-11, SG-13	VOCs	Results for acetone < 5X value detected in laboratory method blanks.	Qualify as non-detect ('U') at the detected value or RL, whichever is higher.
SG-1, SG-3, SG-12	VOCs	CCAL %D > 25% for bromoform, styrene, and trans- 1,3-dichloropropene.	Qualify non-detect results 'UJ.'
AB-1, SG-2, SG-13	VOCs	CCAL %D > 25% for chloromethane, vinyl chloride, 1,1-dichloroethene, acetone, carbon disulfide, and methylene chloride.	Qualify detects 'J' and non-detect results 'UJ.'
SG-7	VOCs	The %R of surrogate 1,2- dichloroethane-d4 > QC limit.	Qualify detected results 'J.'

TABLE 1SUMMARY OF DATA QUALIFICATIONS

Location ID		AB-01	SG-01	SG-02	SG-03	SG-04
Sample ID		AB-1	SG-1	SG-2	SG-3	SG-4
Matrix		Ambient Air	Soil Gas	Soil Gas	Soil Gas	Soil Gas
Depth Interval (ft)		•	-	-		-
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units					
Volatile Organic Compounds						
Chloromethane	PPBV	7.8 UJ	0.36 J	7.8 UJ	4,400 U	9.2 U
Bromomethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Vinyl Chloride	PPBV	3.1 UJ	0.33 U	3.1 UJ	1,700 U	3.7 U
Chloroethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Methylene Chloride	PPBV	7.8 UJ	0.52 J	7.8 UJ	4,400 U	9.2 U
Acetone	PPBV	78 UJ	8.2 UJ	78 UJ	44,000 U	36 J
Carbon Disulfide	PPBV	3.1 UJ	0.33 U	3.2 J	1,700 U	2.7 J
1,1-Dichloroethene	PPBV	3.1 UJ	0.33 U	3.1 UJ	1,700 U	3.7 U
1,1-Dichloroethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
2-Butanone	PPBV	7.8 U	1.2	7.8 U	4,400 U	9.2 U
Chloroform	PPBV	3.1 U	0.33 U	44	1,700 U	14
1,2-Dichloroethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
1,1,1-Trichloroethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Carbon Tetrachloride	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Bromodichloromethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
1,2-Dichloropropane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
cis-1,3-Dichloropropene	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Trichloroethene	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Benzene	PPBV	3.1 U	0.16 J	3.1 U	7,100	3.7 U
Dibromochloromethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
trans-1,3-Dichloropropene	PPBV	3.1 U	0.33 UJ	3.1 U	1,700 UJ	3.7 U
1,1,2-Trichloroethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Bromoform	PPBV	3.1 U	0.33 UJ	3.1 U	1,700 UJ	3.7 U

Made by

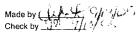
Location ID		AB-01	SG-01	SG-02	SG-03	SG-04
Sample ID		AB-1	SG-1	SG-2	SG-3	SG-4
Matrix		Ambient Air - 07/18/05	Soil Gas - 07/18/05	Soil Gas - 07/18/05	Soil Gas	Soil Gas
Depth Interval (ft)					- 07/18/05	- 07/18/05
Date Sampled						
Parameter	Units					
Volatile Organic Compounds						
1-Methyl-2-Pentanone	PPBV	7.8 U	0.82 U	7.8 U	4,400 U	9.2 U
2-Hexanone	PPBV	7.8 U	0.82 U	7.8 U	4,400 U	9.2 U
Tetrachloroethene	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
1,1,2,2-Tetrachloroethane	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Toluene	PPBV	3.1 U	0.94	6.2	1,700 U	5.5
Chlorobenzene	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
Ethylbenzene	PPBV	3.1 U	0.33 U	1.5 J	1,700 U	1.8 J
Styrene	PPBV	3.1 U	0.33 UJ	3.1 U	1,700 UJ	2.0 J
n,p-Xylene	PPBV	3.1 U	0.16 J	4.6	740 J	7.8
o-Xylene	PPBV	3.1 U	0.33 U	1.5 J	1,700 U	3.1 J
cis-1,2-Dichloroethene	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U
rans-1,2-Dichloroethene	PPBV	3.1 U	0.33 U	3.1 U	1,700 U	3.7 U

Flags assigned during chemistry validation are shown. Made by Check by The Market Strength Check by The

Location ID		SG-05	SG-06	SG-06	SG-07	SG-08
Sample ID		SG-5	FD-2	SG-6	SG-7	SG-8
Matrix Depth Interval (ft)		Soil Gas -	Soil Gas	Soil Gas -	Soil Gas	Soil Gas -
			-			
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units		Field Duplicate (1-1)			
Volatile Organic Compounds						
Chloromethane	PPBV	8.0 U	9.1 U	0.89 U	0.96 U	9.0 U
Bromomethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Vinyl Chloride	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Chloroethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Methylene Chloride	PPBV	8.0 U	9.1 U	0.89	0.96 U	9.0 U
Acetone	PPBV	80 U	91 U	8.5 J	9.6 U	90 U
Carbon Disulfide	PPBV	4.9	3.6 U	0.43	0.39 U	22
1,1-Dichloroethene	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
1,1-Dichloroethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
2-Butanone	PPBV	8.0 U	9.1 U	0.87 J	0.96 U	9.0 U
Chloroform	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
1,2-Dichloroethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 Ų	3.6 U
1,1,1-Trichloroethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Carbon Tetrachloride	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Bromodichloromethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
1,2-Dichloropropane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
cis-1,3-Dichloropropene	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Trichloroethene	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Benzene	PPBV	1.6 J	3.6 U	0.28 J	0.39 U	7.5
Dibromochloromethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
trans-1,3-Dichloropropene	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
1,1,2-Trichloroethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Bromoform	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U

Made by

Location ID Sample ID Matrix Depth Interval (ft)		SG-05 SG-5 Soil Gas -	SG-06	SG-06 SG-6 Soil Gas -	SG-07 SG-7 Soil Gas	SG-08 SG-8
			FD-2			
			Soil Gas			Soil Gas
			-		-	-
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units		Field Duplicate (1-1)			
Volatile Organic Compounds						
4-Methyl-2-Pentanone	PPBV	8.0 U	9.1 U	0.56 J	0.96 U	9.0 U
2-Hexanone	PPBV	8.0 U	9.1 U	0.89 U	0.96 U	9.0 U
Tetrachloroethene	PPBV	3.2 U	3.6 U	0.48	0.56 J	3.6 U
1,1,2,2-Tetrachloroethane	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Toluene	PPBV	5.6	1.7 J	1.6	1.6 J	36
Chlorobenzene	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
Ethylbenzene	PPBV	1.3 J	3.6 U	0.46	0.38 J	3.3 J
Styrene	PPBV	3.2 U	3.6 U	0.27 J	0.39 U	3.6 U
m,p-Xylene	РРВУ	4.6	3.6 U	1.5	1.2 J	9.3
o-Xylene	PPBV	1.6 J	3.6 U	0.64	0.45 J	3.2 J
cis-1,2-Dichloroethene	ррви	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U
trans-1,2-Dichloroethene	PPBV	3.2 U	3.6 U	0.36 U	0.39 U	3.6 U



Location ID		SG-09	SG-10	SG-10	SG-11	SG-12
Sample ID Matrix Depth Interval (ft)		SG-9 Soil Gas	FD-1 Soil Gas	SG-10 Soil Gas	SG-11 Soil Gas	SG-12 Soil Gas
					Date Sampled	
Parameter	Units		Field Duplicate (1-1)			
Volatile Organic Compounds					· · · · · · · · · · · · · · · · · · ·	
Chloromethane	PPBV	8.5 U	9.0 U	9.6 U	8.8 U	0.28 J
Bromomethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
Vinyl Chloride	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
Chloroethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
Methylene Chloride	PPBV	8.5 U	9.0 U	4.9 J	7.3 J	0.90 U
Acetone	PPBV	85 U	90 U	96 U	88 U	12 U
Carbon Disulfide	PPBV	1.5 J	5.0	3.6 J	3.1 J	4.8
1,1-Dichloroethene	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
1,1-Dichloroethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
2-Butanone	PPBV	8.5 U	9.0 U	9.6 U	8.8 U	0.97
Chloroform	PPBV	8.0	2.9 J	1.8 J	3.5 U	1.6
1,2-Dichloroethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
1,1,1-Trichloroethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.11 J
Carbon Tetrachloride	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
Bromodichloromethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
1,2-Dichloropropane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
cis-1,3-Dichloropropene	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
Trichloroethene	PPBV	2.3 J	3.6 U	3.8 U	3.5 U	0.36 U
Benzene	PPBV	3.4 U	3.6 U	3.8 U	1.5 J	0.59
Dibromochloromethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
trans-1,3-Dichloropropene	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 UJ
1,1,2-Trichloroethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
Bromoform	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 UJ

Made by ALA 9/13/05 Check by

Location ID Sample ID Matrix Depth Interval (ft)		SG-09 SG-9 Soil Gas -	SG-10	SG-10 SG-10 Soil Gas -	SG-11 SG-11 Soil Gas	SG-12 SG-12
			FD-1			
			Soil Gas			Soil Gas
			-			-
Date Sampled		07/18/05	07/18/05	07/18/05	07/18/05	07/18/05
Parameter	Units		Field Duplicate (1-1)			
Volatile Organic Compounds						
4-Methyl-2-Pentanone	PPBV	8.5 U	9.0 U	9.6 U	8.8 U	0.48 J
2-Hexanone	PPBV	8.5 U	9.0 U	9.6 U	8.8 U	0.90 U
Tetrachloroethene	PPBV	5.2	4.1	3.8 U	3.5 U	1.2
1,1,2,2-Tetrachloroethane	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
Toluene	PPBV	3.4	2.9 J	2.0 J	4.9	4.2
Chlorobenzene	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
Ethylbenzene	PPBV	3.4 U	3.6 U	3.8 U	1.7 J	0.40
Styrene	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 UJ
m,p-Xylene	PPBV	1.8 J	4.3	2.1 J	5.9	0.78
o-Xylene	PPBV	3.4 U	1.6 J	3.8 U	1.9 J	0.29 J
cis-1,2-Dichloroethene	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U
trans-1,2-Dichloroethene	PPBV	3.4 U	3.6 U	3.8 U	3.5 U	0.36 U

Flags assigned during chemistry validation are shown.

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TABLE 2VALIDATED SOIL GAS AND AMBIENT AIR ANALYTICAL RESULTSNORTH FRANKLIN ST. SITE

Location ID	I	SG-13
Sample ID		SG-13
Matrix		Soil Gas
Depth Interval (ft)		-
Date Sampled		07/18/05
Parameter	Units	
Volatile Organic Compounds		
Chloromethane	PPBV	8.8 UJ
Bromomethane	PPBV	3.5 U
Vinyl Chloride	PPBV	3.5 UJ
Chloroethane	PPBV	3.5 U
Methylene Chloride	PPBV	8.8 UJ
Acetone	PPBV	88 UJ
Carbon Disulfide	PPBV	2.4 J
1,1-Dichloroethene	PPBV	3.5 UJ
1,1-Dichloroethane	PPBV	3.5 U
2-Butanone	PPBV	8.8 U
Chloroform	PPBV	3.5 U
1,2-Dichloroethane	PPBV	3.5 U
1,1,1-Trichloroethane	PPBV	3.5 U
Carbon Tetrachloride	PPBV	3.5 U
Bromodichloromethane	PPBV	3.5 U
1,2-Dichloropropane	PPBV	3.5 U
cis-1,3-Dichloropropene	PPBV	3.5 U
Trichloroethene	PPBV	3.5 U
Benzene	PPBV	3.5 U
Dibromochloromethane	PPBV	3.5 U
trans-1,3-Dichloropropene	PPBV	3.5 U
1,1,2-Trichloroethane	PPBV	3.5 U
Bromoform	PPBV	3.5 U

Flags assigned during chemistry validation are shown.

Made by the state 1/13/05 Check by 12 1/05

TABLE 2 VALIDATED SOIL GAS AND AMBIENT AIR ANALYTICAL RESULTS NORTH FRANKLIN ST. SITE

Location ID	<u> </u>	SG-13
Sample ID		SG-13
Matrix		Soil Gas
Depth Interval (ft)		-
Date Sampled		07/18/05
Parameter	Units	
Volatile Organic Compounds		
4-Methyl-2-Pentanone	PPBV	8.8 U
2-Hexanone	PPBV	8.8 U
Tetrachloroethene	PPBV	3.5 U
1,1,2,2-Tetrachloroethane	PPBV	3.5 U
Toluene	PPBV	4.0
Chlorobenzene	PPBV	3.5 U
Ethylbenzene	PPBV	3.5 U
Styrene	PPBV	3.5 U
m,p-Xylene	PPBV	1.8 J
o-Xylene	PPBV	3.5 U
cis-1,2-Dichloroethene	PPBV	3.5 U
trans-1,2-Dichloroethene	PPBV	3.5 U

Flags assigned during chemistry validation are shown.

Made by

APPENDIX A

SUPPORT DOCUMENTATION

H5G200167

	IN		F CI	IST	ODY	REC	ORI	D	1	1	T	ESTS			U	R	5		
PROJECT N).	4000			SITE NAME	ukun Ste			10-15						LAB_STC-K	of	3		
SAMPLERS				Ē.					В	OTTLE	TYPE /	AND PRE	ESERVATI	VE	PAGE	of	2		-
DELIVERY S					AIRBILL NO	D.:		TOTAL NO.# OF CONTAINERS	- sumus 1						REMARKS	SAMPLE TYPE	BEGINNING DEPTH (IN FEET)	ENDING DEPTH (IN FEET)	(ERPINS)
LOCATION IDENTIFIER	DA	TE	TIME	COMP/ GRAB	SAM	IPLE ID	MATRIX	8 2	é						SUMMA #	0 N			
56-3	7/18	105	1120	the	56-3-3	0-35	65	i	1						1495				
56-2		i	1138	the	SG-Z 2	5-3.0	65	<u> </u>	1						04311				
56-1			1158	IhR	56-1 1.	5-2.0	65	L							1 2738	M			
AB-1			1204	The	20050718	-AB-1	AA	1_	1						93083	M			
SG-12		1	1322	1hr	56-12	35-4.0	65	1	1						04396	w1			
56-13			1339		56-13 -		65	1	1						3402				
56-11	1		1354			4.4.4.9	65	(1						93674	ri	<u> </u>	$\left - \right $	
56-10	+	1	1409		56-10	43-4.8	65	1	(ļ		51501	101			
56-8 66-8	+	1	1515	Ibr		5.0-5.5	65	t	1						2995	n			
56-9		1	1525	the		3.5-4.0	65	ι	1						51488	M			
56-7		+	1535	Ihr	+	25-30	65	1	1					_	51494	WI			
56-4		1	1655	The		3.5-4.0	6,5	1	1				<u> </u>		1127	M			
567-5		丁	1640	IhR	56-5	25-3.0	65	1	1						1530 LH - HAZARDOUS	NU NOLID W	ASTE		
MATRIX		SE · SED	BIENT AIR		SL - SLUDGE WP - DRINKIN WW - WASTE		WG - GROUI SO - SOIL DC - DRILL (WL LEA GS - SOIL WC - DRI	GAS	ATER	WO - OCEAN WS - SURFAC WQ - WATER	E WATER	LF - FLOATING/FR	EPRODU	ICT ON	GW TABI	μE · · ·
SAMPLE TYPE COD		TB# - TR	IP BLANK ATRIX SPIKE DU		RB# - RINSE FR# - FIELD I		N# - NORMA MS# - MATE		NMENTA	L SAMPLE	(# - S					LE SAMPL	.es in a	SINGLE	DAY)
REKINQUI		BY (s	GIGNATURE)	DA	TE TIME	RECEIVED	BY (SIG	NATURE	I		DATE	TIME		/	RUCTIONS Seals Intac	t			
Kitte	ple	le		7/18	65 1900								n a a l	N 14		+			
RELINQUI	SHED) BY (\$	SIGNATURE)	Ď.	TE TIME	RECEIVED	FOR LA			· /-	DATE 07200	TIME 0900	2 B	oxes	FEDEX 4798 1855 1 1833				
Distributio	n: Ori	ginal ac	ccompanies	shipmer	nt, copy to c	oordinator fiel							4	L A	V 1844 DF 072005				

URSF-075C/1 OF 1/CofCR/GCM

HSG200167

СНА	IN O	FC	JST	ODY	REC	OR	D				TESTS			U	R	5		
PROJECT N 11742 SAMPLERS	0. 2 <i>11. 890</i> (print/signa	00 TURE)		site name Nota Fia				<u>で、</u>	OTTLE		AND PR	ESERVAT	TVE	LAB <u>STL K</u> COOLER <u>I</u> PAGE <u>Z</u>	of	3		-
DELIVERY S	uccabe ervice: Pe			, AIRBILL NO			TOTAL NO.# OF CONTAINERS	L Summer						REMARKS	SAMPLE TYPE	BEGINNING DEPTH (IN FEET)	ENDING DEPTH (IN FEET)	FIGLD LOT NO. # (ERPIMS)
IDENTIFIER SG-G FD-1 FD-2	DATE	TIME 1555	GRAB 1 h.e 1 hr 1 hr	SAMF SG-6 2, 20050718 20050718	-FD-1	MATRIX 65 65 65 67 5	1 1 1	9 1 1						93265 1787 51529.	KJ FRI FRZ			
				FD- FD-	1 =	5		00										
MATRIX CODES	SE - SEC SH - HAZ	BENT AIR IMENT ARDOUS SOLID	WASTE	SL - SLUDGE WP - DRINKING WW - WASTE V RB# - RINSE BI	WATER	WG - GROUN SO - SOIL DC - DRILL (N# - NORMA	OUTTINGS	<u></u>	GS - SC WC - D	RILLING W		WO - OCEAN WS - SURFA WQ - WATEF	CE WATER		E PRODU			
RELINQUI	SHED BY (SHED BY (SHE	ATRIX SPIKE DU SIGNATURE) SIGNATURE)	DA <mark>7/18,</mark> DA	FR# - FIELD RE	RECEIVED	FOR LA	NATURE)		JRE)	DATE				RUCTIONS				

URSF-075C/1 OF 1/CotCR/GCM

PROJECT NARRATIVE H5G200167 REVISED

This report has been revised. The narrative has been updated and expanded chromatograms have been included.

The results reported herein are applicable to the samples submitted for analysis only.

This report shall not be reproduced except in full, without the written approval of the laboratory.

The original chain of custody documentation is included with this report.

Sample Receipt

There were no problems with the condition of the samples received.

Quality Control

Unless otherwise noted, all holding times and QC criteria were met and the test results shown in this report meet all applicable NELAC requirements.

Samples AB-1, and SG-13 analyzed on instrument MG and samples SG-7, and FD-1 analyzed on instrument MT, were originally analyzed at 500 mL and exhibited poor chromatography and poor or no internal standard and/or surrogate recoveries. The chromatography observed on these two separate instruments is consistent with matrix freezing of the instrument concentrator traps, restricting the flow of sample to the GCMS. This could have resulted from water and/or carbon dioxide in the sample causing the trap freeze-up. These samples impacted the recoveries of internal standards in subsequent samples run on the instrument. The samples were re-analyzed at 50 mL. Both sets of data for these four samples are reported. Due to the detrimental effects of the sample matrix on the instrument, the remaining samples (SG-2, SG-11, SG-10, SG-8, SG-9, SG-4, SG-5, and FD-2) were analyzed at 50 mL.

STL Knoxville maintains the following certifications, approvals and accreditations: Arkansas DEQ Cert. # 03-049-0, California DHS ELAP Cert. #2423, Colorado DPHE, Connecticut DPH Cert. #PH-0223, Florida DOH Cert. #E87177, Georgia DNR Cert. #906, Hawaii DOH, Illinois EPA Cert. # 000687, Indiana DOH Cert. #C-TN-02, Kansas DHE Cert. # E-10349, Kentucky DEP Lab ID #90101, Louisiana DEQ Cert. #03079, Louisiana DOHH Cert. #LA030024, Maryland DHMH Cert. #277, Massachusetts DEP Cert. #M-TN009, Michigan DEQ Lab ID #9933, New Jersey DEP Cert. #TN001, New York DOH Lab #10781, North Carolina DPH Lab ID #21705, North Carolina DEHNR Cert. #64, Oklahoma DEQ ID #9415, Pennsylvania DEP Cert. # 68-576, South Carolina DHEC Lab ID #84001001, Tennessee DOH Lab ID #02014, Utah DOH Cert. #QUAN3, Virginia DGS Lab ID #00165, Washington DOE Lab #C120, Wisconsin DNR Lab ID #998044300, US Army Corps of Engineers, Naval Facilities Engineering Service Center, US EPA Perchlorate Approval and USDA Soil Permit #S-46424. This list of approvals is subject to change and does not imply that laboratory certification is available for all parameters reported in this environmental sample data report.

PROJECT NARRATIVE H5G200167 REVISED

Sample SG-3 was reported with elevated reporting limits for all analytes due to the presence of non-target compounds. A dilution was necessary prior to analysis, and the reporting limits were adjusted accordingly.

STL Knoxville maintains the following certifications, approvals and accreditations: Arkansas DEQ Cert. # 03-049-0, California DHS ELAP Cert. #2423, Colorado DPHE, Connecticut DPH Cert. #PH-0223, Florida DOH Cert. #E87177, Georgia DNR Cert. #906, Hawaii DOH, Illinois EPA Cert. # 000687, Indiana DOH Cert. #C-TN-02, Kansas DHE Cert. # E-10349, Kentucky DEP Lab ID #90101, Louisiana DEQ Cert. #03079, Louisiana DOHH Cert. #LA030024, Maryland DHMH Cert. #277, Massachusetts DEP Cert. #M-TN009, Michigan DEQ Lab ID #9933, New Jersey DEP Cert. #TN001, New York DOH Lab #10781, North Carolina DPH Lab ID #21705, North Carolina DEHNR Cert. #64, Oklahoma DEQ ID #9415, Pennsylvania DEP Cert. # 68-576, South Carolina DHEC Lab ID #84001001, Tennessee DOH Lab ID #02014, Utah DOH Cert. #QUAN3, Virginia DGS Lab ID #00165, Washington DOE Lab #C120, Wisconsin DNR Lab ID #998044300, US Army Corps of Engineers, Naval Facilities Engineering Service Center, US EPA Perchlorate Approval and USDA Soil Permit #S-46424. This list of approvals is subject to change and does not imply that laboratory certification is available for all parameters reported in this environmental sample data report.

Client Sample ID: INTRA-LAB BLANK

GC/MS Volatiles

Lot-Sample # H5G220000 - 06	ib W	ork Order # HF2WT	IAA Ma	trix: AIR
Prep Date: 7/21/05 Prep Batch #: 5203061		ate Received: 7/21/05 nalysis Date 7/21/05		
Dilution Factor.: 1	Ν	Iethod TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	0.20	ND	0.91
Acetone	0.47	5.0	1.1 J	12
Ethylbenzene	ND	0.20	ND	0.87
2-Hexanone	ND	0.50	ND	2.0
Methylene chloride	ND	0.50	ND	1.7
Benzene	ND	0.20	ND	0.64
Styrene	ND	0.20	ND	0.85
1,1,2,2-Tetrachloroethane	ND	0.20	ND	1.4
Tetrachloroethene	ND	0.20	ND	1.4
Toluene	ND	0.20	ND	0.75
1,1,1-Trichloroethane	ND	0.20	ND	1.1
1,1,2-Trichloroethane	ND	0.20	ND	1.1
Trichloroethene	ND	0.20	ND	1.1
Vinyl chloride	ND	0.20	ND	0.51
o-Xylene	ND	0.20	ND	0.87
m-Xylene & p-Xylene	ND	0.20	ND	0.87
Bromodichloromethane	ND	0.20	ND	1.3
2-Butanone (MEK)	ND	0.50	ND	1.5
4-Methyl-2-pentanone (MIBK)	ND	0.50	ND	2.0
Bromoform	ND	0.20	ND	2.1
Bromomethane	ND	0.20	ND	0.78
Carbon disulfide	ND	0.20	ND	0.62
Carbon tetrachloride	ND	0.20	ND	1.3
Chlorobenzene	ND	0.20	ND	0.92
Dibromochloromethane	ND	0.20	ND	1.7
Chloroethane	ND	0.20	ND	0.53
Chloroform	ND	0.20	ND	0.98
Chloromethane	ND	0.50	ND	1.0
1,1-Dichloroethane	ND	0.20	ND	0.81
1,2-Dichloroethane	ND	0.20	ND	0.81
1,1-Dichloroethene	ND	0.20	ND	0.79
cis-1,2-Dichloroethene	ND	0.20	ND	0.79
trans-1,2-Dichloroethene	ND	0.20	ND	0.79
	ND ND	0.20	ND	0.92
1,2-Dichloropropane		0.20	ND	0.91
cis-1,3-Dichloropropene	ND	0.20		0.71

TO-14 Conv Rev. 5

Client Sample ID: INTRA-LAB BLANK

GC/MS Volatiles

Lot-Sample # H5G250000 - 28	8B	Work Order # HF7E71	AA Ma	trix: AIR
Prep Date: 7/25/05 Prep Batch #: 5206288		Date Received: 7/20/05 Analysis Date 7/25/05		
Dilution Factor.: 1		Method: TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	0.20	ND	0.91
Acetone	0.32	5.0	0.75 J	12
Ethylbenzene	ND	0.20	ND	0.87
2-Hexanone	ND	0.50	ND	2.0
Methylene chloride	ND	0.50	ND	1.7
Benzene	ND	0.20	ND	0.64
Styrene	ND	0.20	ND	0.85
1,1,2,2-Tetrachloroethane	ND	0.20	ND	1.4
Tetrachloroethene	ND	0.20	ND	1.4
Toluene	ND	0.20	ND	0.75
1,1,1-Trichloroethane	ND	0.20	ND	1.1
1,1,2-Trichloroethane	ND	0.20	ND	1,1
Trichloroethene	ND	0.20	ND	1.1
Vinyl chloride	ND	0.20	ND	0.51
o-Xylene	ND	0.20	ND	0.87
m-Xylene & p-Xylene	ND	0.20	ND	0.87
Bromodichloromethane	ND	0.20	ND	1.3
2-Butanone (MEK)	ND	0.50	ND	1.5
4-Methyl-2-pentanone (MIBK)	ND	0.50	ND	2.0
Bromoform	ND	0.20	ND	2.1
Bromomethane	ND	0.20	ND	0.78
Carbon disulfide	ND	0.20	ND	0.62
Carbon tetrachloride	ND	0.20	ND	1.3
Chlorobenzene	ND	0.20	ND	0.92
Dibromochloromethane	ND	0.20	ND	1.7
Chloroethane	ND	0.20	ND	0.53
Chloroform	ND	0.20	ND	0.98
Chloromethane	ND	0.50	ND	1.0
1,1-Dichloroethane	ND	0.20	ND	0.81
1,2-Dichloroethane	ND	0.20	ND	0.81
1,1-Dichloroethene	ND	0.20	ND	0.79
cis-1,2-Dichloroethene	ND	0.20	ND	0.79
trans-1,2-Dichloroethene	ND	0.20	ND	0.79
1,2-Dichloropropane	ND	0.20	ND	0.92
cis-1,3-Dichloropropene	ND	0.20	ND	0.91

TO-14 Conv Rev. 5

Client Sample ID: INTRA-LAB BLANK

GC/MS Volatiles

		ate Received: 7/20/05 nalvsis Date 7/22/05		
Prep Date: 7/22/05 Prep Batch #: 5206164	А	nalysis Date 7/22/05		
Dilution Factor.: 1	Μ	lethod: TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3)	REPORTING LIMIT (ug/m3)
rans-1,3-Dichloropropene	ND	0.20	ND	0.91
Acetone	0.26	5.0	0.61 J	12
Ethylbenzene	ND	0.20	ND	0.87
2-Hexanone	ND	0.50	ND	2.0
Methylene chloride	ND	0.50	ND	1.7
Benzene	ND	0.20	ND	0.64
Styrene	ND	0.20	ND	0.85
1,1,2,2-Tetrachloroethane	ND	0.20	ND	1.4
Tetrachloroethene	ND	0.20	ND	1.4
Toluene	ND	0.20	ND	0.75
1,1,1-Trichloroethane	ND	0.20	ND	1.1
1,1,2-Trichloroethane	ND	0.20	ND	1.1
Trichloroethene	ND	0.20	ND	1.1
Vinyl chloride	ND	0.20	ND	0.51
o-Xylene	ND	0.20	ND	0.87
m-Xylene & p-Xylene	ND	0.20	ND	0.87
Bromodichloromethane	ND	0.20	ND	1.3
2-Butanone (MEK)	ND	0.50	ND	1.5
4-Methyl-2-pentanone (MIBK)	ND	0.50	ND	2.0
Bromoform	ND	0.20	ND	2.1
Bromomethane	ND	0.20	ND	0.78
Carbon disulfid e	ND	0.20	ND	0.62
Carbon tetrachloride	ND	0.20	ND	1.3
Chlorobenzene	ND	0.20	ND	0.92
Dibromochloromethane	ND	0.20	ND	1.7
Chloroethane	ND	0.20	ND	0.53
Chloroform	ND	0.20	ND	0.98
Chloromethane	ND	0.50	ND	1.0
1,1-Dichloroethane	ND	0.20	ND	0.81
1,2-Dichloroethane	ND	0.20	ND	0.81
1,1-Dichloroethene	ND	0.20	ND	0.79
cis-1,2-Dichloroethene	ND	0.20	ND	0.79
trans-1,2-Dichloroethene	ND	0.20	ND	0.79
1,2-Dichloropropane	ND	0.20	ND	0.92
cis-1,3-Dichloropropene	ND	0.20	ND	0.91

Client Sample ID: INTRA-LAB BLANK

GC/MS Volatiles

Lot-Sample # H5G220000 - 12	VIB V	Vork Order # HF2141	AA Ma	trix: AIR
Prep Date: 7/21/05 Prep Batch #: 5203121		Date Received: 7/20/05 Analysis Date 7/21/05		
Dilution Factor.: 1	Ν	Aethod TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	0.20	ND	0.91
Acetone	0.27	5.0	0.63 J	12
Ethylbenzene	ND	0.20	ND	0.87
2-Hexanone	ND	0.50	ND	2.0
Methylene chloride	ND	0.50	ND	1.7
Benzene	ND	0.20	ND	0.64
Styrene	ND	0.20	ND	0.85
1,1,2,2-Tetrachloroethane	ND	0.20	ND	1.4
Tetrachloroethene	ND	0.20	ND	1.4
Toluene	ND	0.20	ND	0.75
1,1,1-Trichloroethane	ND	0.20	ND	1.1
1,1,2-Trichloroethane	ND	0.20	ND	1.1
Trichloroethene	ND	0.20	ND	1.1
Vinyl chloride	ND	0.20	ND	0.51
o-Xylene	ND	0.20	ND	0.87
m-Xylene & p-Xylene	ND	0.20	ND	0.87
Bromodichloromethane	ND	0.20	ND	1.3
2-Butanone (MEK)	ND	0.50	ND	1.5
4-Methyl-2-pentanone (MIBK)	ND	0.50	ND	2.0
Bromoform	ND	0.20	ND	2.1
Bromomethane	ND	0.20	ND	0.78
Carbon disulfide	ND	0.20	ND	0.62
Carbon tetrachloride	ND	0.20	ND	1.3
Chlorobenzene	ND	0.20	ND	0.92
Dibromochloromethane	ND	0.20	ND	1.7
Chloroethane	ND	0.20	ND	0.53
Chloroform	ND	0.20	ND	0.98
Chloromethane	ND	0.50	ND	1.0
1,1-Dichloroethane	ND	0.20	ND	0.81
1,2-Dichloroethane	ND	0.20	ND	0.81
1,1-Dichloroethene	ND	0.20	ND	0.79
cis-1,2-Dichloroethene	ND	0.20	ND	0.79
trans-1,2-Dichloroethene	ND	0.20	ND	0.79
1,2-Dichloropropane	ND	0.20	ND	0.92
cis-1,3-Dichloropropene	ND	0.20	ND	0.91

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TO-14 Conv Rev. 5

		STL-Knoxv TO-14 Autosam 'ME'		
Sample	Volume	Position	Date	Time ====================================
CCV LCSA BLK BLKA HFVXP HFVX4 HFV05D HFVXW HFVX1 HFVX6 HFVX1R HFVX1RR HFVX1RR HFVX7	103m] 101m] 501m] 501m] 253m] 501m] 253m] 502m] 502m] 493m] 27m] 28m] 51m]	15 15 16 16 1 2 3 5 6 7 6 6 8	7/21/200 7/21/200 7/21/200 7/21/200 7/21/200 7/21/200 7/21/200 7/21/200 7/21/200 7/21/200 7/21/200 7/22/200 7/22/200	12:26:48 PM 2:04:26 PM 4:29:13 PM 5:07:24 PM 5:44:23 PM 6:22:55 PM 6:59:39 PM 7:37:45 PM 8:15:58 PM 9:38:51 PM 10:35:41 PM 7:10:08 AM 8:22:12 AM

050721 -Knoxville CTI

STL Knoxville

CONTINUING CALIBRATION COMPOUNDS

Instrument ID: mg.i Injection Date: 21-JUL-2005 12:26 Jab File ID: gccvg21.d Init. Cal. Date(s): 01-JUL-2005 01-JUL-2005 Analysis Type: AIR Init. Cal. Times: 12:54 16:32 Jab Sample ID: GCCV Quant Type: ISTD Aethod: /chem/gcms/mg.i/G072105.b/TO14.m

			MIN		MAX	
COMPOUND	RRF / AMOUNT	RF10			\$D / \$DRIFT	
39 Chloroform	2.78591					
40 1,1,1-Trichloroethane	2.48983	2.66647	0.000			•
41 Cyclohexane	0.76498				•	
41 Cyclonexand 42 Carbon Tetrachloride	2.76630					
43 Benzene	0.99049				,	
44 1,2-Dichloroethane	0.43213	0.46033	0.000	-6.52560		
44 1,2-Dichiolocchance 45 2,2,4-trimethylpentane	2.31797	2.30177	0.000	0.69923	•	
	0.90299		0.000	-0.23187	•	•
46 Heptane 47 Trichloroethene	0.46535	0.47624	0.000	-2.34172		
	0.16823	0.18182	10.000	-8.0817		
48 1-Butanol 49 1,2-Dichloropropane	0.38015		0.000	-5.4355		•
	0.32811		0.000	-7.9499	30.0000	
50 Dibromomethane 51 Bromodichloromethane	0.58594	0.6508	5 0.000	-11.0792	30.0000	
	0.51706		0.000		-	•
52 cis-1,3-Dichloropropene	0.84075		3 0 . 000	-1.8776		•
53 4-Methyl-2-pentanone	1.22579		7 0.000	-10.0404		
54 Toluene	0.41504	•			30.0000	
55 Octane 56 trans-1,3-Dichloropropene	0.33654		5 0.00	o∣ <u>€26.5667</u>	8) 30.0000	•
56 trans-1, 3-Dichiorophopene	0.38698		0 0.00	-17.2909	8 30.0000	•
57 1,1,2-Trichloroethane	0.45442		0 0.00	0 -6.2886	7 30.0000	
58 Tetrachloroethene	0.44979		5 0.00	0 -4.5502	9 30.0000	
59 2-Hexanone	0.5908		7 0.00	0 -22.9663	9 30.0000	
60 Dibromochloromethane	0.55724	•	4 0.00	0 -22.3967	6 30.0000	
61 1,2-Dibromoethane	0.96941		8 0.00			
62 Chlorobenzene	1.3656	- 1	6 0.00		30.0000	00 Averaged
63 Ethylbenzene	1.0086	- 1	1 0.00			
64 m-Xylene (For p-)	0.9248	- 1	.9 0.00		.5 30.0000	
65 Nonane	1.0233	•	6 0.00		30.0000	
66 o-Xylene	0.7046		10.00		30.0000	•
67 Styrene	0.5340			0 (-31.500		00 Average
68 Bromoform	0.8518		24 0.00			00 Average
69 a-Pinene	1.6592	•	95 0.00		38 30.000	00 Average
70 Cumene	0.5935		4 0.00		75 30.000	00 Average
71 Camphene	0.8485	-	0.00	- 1		00 Average
72 1,1,2,2-Tetrachloroethane			61 0.00			00 Average
73 1,2,3-Trichloropropane	0.2582	0.279	1			

		TO-14 Autosam 'ME'	pler Log	
Sample	Volume	Position	Date	Time
CCV LCS BLK HFV05 HFVXV HFVX1 HFVX6 HFVX7 HFVX8 HFV0A HFV0D HFV0D	101ml 101ml 503ml 503ml 51ml 51ml 51ml 51ml 51ml 51ml 51ml 51	15 15 16 16 3 4 6 7 8 9 10 11 11	7/22/200 7/22/200 7/22/200 7/22/200 7/22/200 7/22/200 7/22/200 7/22/200 7/22/200 7/22/200 7/22/200 7/22/200 7/22/200	9:33:23 AM 10:09:24 AM 10:51:53 AM 11:30:19 AM 12:07:29 PM 12:43:22 PM 1:19:20 PM 1:55:11 PM 2:31:10 PM 5:02:58 PM 5:39:23 PM 6:15:59 PM 6:52:27 PM

050722 STL-Knoxville TO-14 Autosampler Log 'MF'

ata File: /chem/gcms/mg.i/G072205.b/gccvg22.d eport Date: 22-Jul-2005 09:57

STL Knoxville

CONTINUING CALIBRATION COMPOUNDS

Instrument ID: mg.i Injection Date: 22-JUL-2005 09:33 Lab File ID: gccvg22.d Init. Cal. Date(s): 01-JUL-2005 01-JUL-2005 Analysis Type: AIR Init. Cal. Times: 12:54 16:32 Lab Sample ID: GCCV Quant Type: ISTD Method: /chem/gcms/mg.i/G072205.b/T014.m

		1	MIN	I	ļ	MAX	
COMPOUND	RRF / AMOUNT	RF10				%D / %DRIFT	
				====		*********	=============================
\$ 4 1,2-Dichloroethane-d4	0.16189	0.15951			1.46963	30.0000	
\$ 5 Toluene-d8	0.96626	0.98255	0.000	·	-1.68589	30.00000	Averaged
\$ 6 4-Bromofluorobenzene	0.64632	0.55755	0.000	:	13.73420	30.00000	Averaged
7 Chlorodifluoromethane	3.47060	2.47879	0.000	$1 \in$	28.57746	30.00000	Averaged
8 Dichlorodifluoromethane	3.73704	2.74008	0.000	1	26.67783	30.00000	Averaged
9 1,2-Dichlorotetrafluoroetha	2.86670	2.17921	0.000	:	23.98184	30.00000	·
10 Chloromethane	0.59116	0.43548	0.000	14. 1	26.33379	30.00000	Averaged
11 Vinyl Chloride	1.87479	1.33903	0.000		28.57686	30.0000	
12 n-Butane	4.16059	2.87217	0.000	1	30.96722	* ··	
13 1,3-Butadiene	1.56206	1.21784	0.000		22.03669	30.0000	-
14 Methanol	0.84045	0.65022	0.000		22.63398	30.0000	· .
15 Bromomethane	1.54138	1.18145	10.000	1	23.35152	30.0000	
16 Chloroethane	0.96859	0.75423	0.000	1	22.13069	30.0000	
17 Vinyl Bromide	3.31117	2.51923	0.000	1	23.91706	30.00000	Averaged
18 Trichlorofluoromethane	3.84505	2.94536	0.000		23.39873	30.00000	
19 Pentane	0.66348	0.45393	0.000	$ \leq$	31.58394	30.00000	Averaged <
20 Ethyl Ether	1.64256	1.42715	0.000)	13.11410	30.0000	Averaged
21 Acrolein	0.59698	0.48102	0.000		19.42476	30.00000	Averaged
22 1,1-Dichloroethene	1.76544	1.27211	0.000	NC	27.94386	30.0000) Averaged
23 1,1,2-Trichlorotrifluoroeth	3.48059	2.70066	0.000)	22.40778		
24 Acetone	0.89685	0.65260	0.000	ИĆ	27.23440	1	
25 Carbon Disulfide	5.14606	3.63433	10.000	$ \langle \rangle $	29.37649	30.0000	
26 Acetonitrile	1.27860	1.07193	0.000	0	16.16396	30.0000	•
27 3-Chloropropene	1.98672	1.60275	5 0 . 000	0	19.32661	30.0000	· .
28 Methylene Chloride	1.57247	1.14947	10.00	o⊧⊂	26.90017	30.0000	
29 2,3-Dimethyl butane	5.73089	4.30142	L 0.00	0	24.94349	30.0000	· .
30 tert-butanol	. 2.99272	2.5536	9 0.00	0	14.66989	30.0000	0 Averaged
31 Acrylonitrile	1.22058	1.0934	5 0.00	0	10.41448	30.0000	0 Averaged
32 trans-1,2-Dichloroethene	1.69810	1.3067	1 0.00	0	23.0486	5 30.0000	0 Averaged
33 Methyl-t-Butyl Ether	2.65088	2.3215	1 0.00	0	12.4247	30.0000	0 Averaged
	1.87298	1.3834	7 0.00	01 (26.1354	3) 30.0000	0 Averaged
34 Hexane 35 1,1-Dichloroethane	3.02922	2.4528	210.00	0	19.0280	B 30.0000	0 Averaged
	3.40733	3.1740	6 0.00	0	6.8462	0 30.0000	0 Averaged
<pre> 36 Vinyl Acetate 37 cis 1,2-Dichloroethene</pre>	1.65646	1.3004	3 0.00	0	21.4934	1 30.0000	0 Averaged
	0.60471	0.5442	4 0.00	0	9.9999	7 30.0000	0 Averaged
38 2-Butanone			1	L			

APPENDIX B

VALIDATED FORM I's

N:\11173258.00000\WORD\North Franklin Street Air July 2005 DUSR.doc 09/13/05 11:30 AM

Client Sample ID: SG-1

GC/MS Volatiles

Prep Batch #: 5203121 Dilution Factor: 1.64 Method: TO-15 PARAMETER (ppb(v/v)) LIMIT (ppb(v/v)) RESULTS (ug/m3) RBPORT LIMIT (p trans-1,3-Dichloropropene ND 0.33 ND 1.5 Acetone 5.0 8.2 -12 J 19 Ethylbenzene ND 0.33 ND 1.4 2-Hexanone ND 0.32 ND 1.4 2-Hexanone ND 0.33 ND 1.4 2-Hexanone ND 0.33 ND 1.4 2-Hexanone ND 0.33 ND 2.8 Benzene 0.16 0.33 0.52 J 1.0 Styrene ND 0.33 ND 2.2 1.0 Styrene ND 0.33 ND 2.2 1.0 Styrene ND 0.33 ND 1.8 1.1,2 1.1 1.1 1.2 1.2 1.2 1.2	
PARAMETER (ppb(v/v)) LIMIT (ppb(v/v)) RESULTS (ug/m3) LIMIT (ug/m3) trans-1,3-Dichloropropene ND 0.33 ND 1.5 Acetone 5.0 6.5 8.2 D-13 JF 19 Ethylbenzene ND 0.33 ND 1.4 2.4 2.4 1.4 2-Hexanone ND 0.82 ND 3.4 Methylene chloride 0.52 0.82 1.8 J 2.8 Benzene 0.16 0.33 ND 1.4 1.1,2,2-2 1.0 1.4 1,1,2,2-Tetrachloroethane ND 0.33 ND 2.3 1.0 Styrene ND 0.33 ND 2.2 2 2 Toluene 0.94 0.33 3.5 1.2 1.1 1.1 1.4 1,1,2-Trichloroethane ND 0.33 ND 1.8 1.1,1-Trichloroethane ND 0.33 ND 1.4 m-Xylene & p-Xylene 0.16 0.33 <th></th>	
Acteone 5.0 5.0 8.2 D-12 JF 19 Ethylbenzene ND 0.33 ND 1.4 2-Hexanone ND 0.82 ND 3.4 Methylene chloride 0.52 0.82 1.8 J 2.8 Benzene 0.16 0.33 0.52 J 1.0 Styrene ND 0.33 ND 2.3 1.4 1,1,2,2-Tetrachloroethane ND 0.33 ND 2.2 Toluene 0.94 0.33 3.5 1.2 1,1,1-Trichloroethane ND 0.33 ND 1.8 Trichloroethane ND 0.33 ND 1.8 Trichloroethane ND 0.33 ND 1.4 m-Xylene & p-Xylene 0.16 0.33 ND 1.4 Bromodichloromethane ND 0.33 ND 1.4 Bromodichloromethane ND 0.33 ND 3.4 Bromodichloromethane	
Activity Instrume	
Lith ND 0.82 ND 3.4 2-Hexanone ND 0.52 0.82 1.8 J 2.8 Benzene 0.16 0.33 0.52 J 1.0 Styrene ND 0.33 ND 2.3 Tetrachloroethane ND 0.33 ND 2.2 Totuene 0.94 0.33 3.5 1.2 1,1,1-Trichloroethane ND 0.33 ND 2.2 Toluene 0.94 0.33 3.5 1.2 1,1,1-Trichloroethane ND 0.33 ND 1.8 1,1,2-Trichloroethane ND 0.33 ND 1.8 1,1,2-Trichloroethane ND 0.33 ND 1.8 1,2-Trichloroethane ND 0.33 ND 1.4 m-Xylene & p-Xylene ND 0.33 ND 2.2 2-Butanone (MEK) 1.2 0.82 3.4 2.4 4-Methyl-2-pentanone (MIBK) ND 0.33 ND 3.4 Bromoform ND 0.33 ND	
Description Description Description Methylene chloride 0.52 0.82 1.8 J 2.8 Benzene 0.16 0.33 0.52 J 1.0 Styrene ND 0.33 ND 2.3 Tetrachloroethane ND 0.33 ND 2.2 Toluene 0.94 0.33 3.5 1.2 1,1,1-Trichloroethane ND 0.33 ND 1.8 1,1,2-Trichloroethane ND 0.33 ND 1.8 Trichloroethane ND 0.33 ND 1.8 Vinyl chloride ND 0.33 ND 1.4 m-Xylene & p-Xylene 0.16 0.33 ND 1.4 Bromodichloromethane ND 0.33 ND 1.4 Bromodichloromethane ND 0.33 ND 2.2 2-Butanone (MEK) 1.2 0.82 ND 3.4 Bromoform ND 0.33 ND 1.0 </td <td></td>	
Net Hyper Childre 0.16 0.33 0.52 J 1.0 Styrene ND 0.33 ND 1.4 1,1,2,2-Tetrachloroethane ND 0.33 ND 2.3 Tetrachloroethane ND 0.33 ND 2.3 Tetrachloroethane ND 0.33 ND 2.2 Toluene 0.94 0.33 3.5 1.2 1,1,1-Trichloroethane ND 0.33 ND 1.8 1,1,2-Trichloroethane ND 0.33 ND 1.8 Vinyl chloride ND 0.33 ND 1.4 m-Xylene & p-Xylene 0.16 0.33 ND 1.4 Bromodichloromethane ND 0.33 ND 2.2 2-Butanone (MEK) 1.2 0.82 ND 3.4 Bromoform ND 0.33 ND 1.3 Carbon disulfide ND 0.33 ND 1.4 Dibromochloromethane ND 0.33 ND	
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trans-1,2-Dichloroethene ND 0.33 ND 1.3	
1,2-Dichloropropane ND 0.33 ND 1.5	

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Client Sample ID: SG-1

GC/MS Volatiles

Lot-Sample # H5G200167 - 003	Work Order #	HFVXW1AD	Matrix:	AIR
SURROGATE	PERCENT RECOVERY		LABORATOR CONTROL LIMITS (%)	Y
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	109 101 85		70 - 130 70 - 130 70 - 130	

Oualifiers

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.
 J Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

Client Sample ID: SG-2

GC/MS Volatiles

Lot-Sample # H5G200167 - 002	2	Work Order #	HFVXVIAD	Ma	trix:	AIR
Date Sampled: 7/18/05 Prep Date: 7/22/05 Date Date 5206164		Date Received: Analysis Date				
Prep Batch #: 5206164 Dilution Factor.: 15.5		Method	: TO-15			
					REPORTI	NG
PARAMETER	RESULTS (ppb(v/v))	LIMIT (p	5 15 07 7	LTS (ug/m3)	LIMIT (up	
trans-1,3-Dichloropropene	ND	3.1	ND		14	
Acetone	-H ND	78	+ 13-26-	YBL	15180	
Ethylbenzene	1.5	3.1	6.6	J	13	
2-Hexanone	ND	7.8	ND		32	
Methylene chloride	ND	7.8	ND	55	27	
Benzene	ND	3.1	ND		9.9	
Styrene	ND	3.1	ND		13	
1,1,2,2-Tetrachloroethane	ND	3.1	ND		21	
Tetrachloroethene	ND	3.1	ND		21	
Toluene	6.2	3.1	23		12	
1,1,1-Trichloroethane	ND	3.1	ND		17	
1,1,2-Trichloroethane	ND	3.1	ND		17	
Trichloroethene	ND	3.1	ND	/	17	
Vinyl chloride	ND	3.1	ND	\mathcal{T}	7.9	
o-Xylene	1.5	3.1	6.5	J	13	
m-Xylene & p-Xylene	4.6	3.1	20		13	
Bromodichloromethane	ND	3.1	ND		21	
2-Butanone (MEK)	ND	7.8	ND		23	
4-Methyl-2-pentanone (MIBK)	ND	7.8	ND		32	
Bromoform	ND	3.1	ND		32	
Bromomethane	ND	3.1	ND	-	12	
Carbon disulfide	3.2	3.1	10	-5	9.7	
Carbon tetrachloride	ND	3.1	ND		20	
Chlorobenzene	ND	3.1	ND		14	
Dibromochloromethane	ND	3.1	ND		26	
Chloroethane	ND	3.1	ND		8.2	
Chloroform	44	3.1	220	/	15	
Chloromethane	ND	7.8	ND	\odot	16	
1,1-Dichloroethane	ND	3.1	ND		13	
1,2-Dichloroethane	ND	3.1	ND		13	
1,1-Dichloroethene	ND	3.1	ND	125	12	
cis-1,2-Dichloroethene	ND	3.1	ND		12	
trans-1,2-Dichloroethene	ND	3.1	ND		12	
1,2-Dichloropropane	ND	3.1	ND		14	•
cis-1,3-Dichloropropene	ND	3.1	ND		14	, (

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Client Sample ID: SG-2

GC/MS Volatiles

Lot-Sample # H5G200167 - 002	Work Order # HFVXV1AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	120 92 83	70 - 130 70 - 130 70 - 130

Qualifiers

BMethod blank contamination. The associated method blank contains the target analyte at a reportable level.JEstimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

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Client Sample ID: SG-3

GC/MS Volatiles

Lot-Sample # H5G200167 - 00	v 1 v	Vork Order # HFVXP	1AD Ma	trix: AIR
Date Sampled: 7/18/05 Prep Date: 7/21/05 Prep Batch #: 5203121		Pate Received 7/20/05 nalysis Date 7/21/05		
Dilution Factor.: 8739.4	N	fethod TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	1700	ND OF	7900
Acetone	ND	44000	ND	100000
Ethylbenzene	ND	1700	ND	7600
2-Hexanone	ND	4400	ND	18000
Methylene chloride	ND	4400	ND .	15000
Benzene	7100	1700	23000	5600
Styrene	ND	1700	ND OF	7400
1,1,2,2-Tetrachloroethane	ND	1700	ND	12000
Tetrachloroethene	ND	1700	ND	12000
Toluene	ND	1700	ND	6600
1,1,1-Trichloroethane	ND	1700	ND	9500
1,1,2-Trichloroethane	ND	1700	ND	9500
Trichloroethene	ND	1700	ND	9400
Vinyl chloride	ND	1700	ND	4500
o-Xylene	ND	1700	ND	7600
m-Xylene & p-Xylene	740	1700	3200 J	7600
Bromodichloromethane	ND	1700	ND	12000
2-Butanone (MEK)	ND	4400	ND	13000
4-Methyl-2-pentanone (MIBK)	ND	4400	ND	18000
Bromoform	ND	1700	ND OS	18000
Bromomethane	ND	1700	ND	6800
Carbon disulfide	ND	1700	ND	5400
Carbon tetrachloride	ND	1700	ND	11000
Chlorobenzene	ND	1700	ND	8000
Dibromochloromethane	ND	1700	ND	15000
Chloroethane	ND	1700	ND	4600
Chloroform	ND	1700	ND	8500
Chloromethane	ND	4400	ND	9000
1,1-Dichloroethane	ND	1700	ND	7100
1,2-Dichloroethane	ND	1700	ND	7100
1,1-Dichloroethene	ND	1700	ND	6900
cis-1,2-Dichloroethene	ND	1700	ND	6900
trans-1,2-Dichloroethene	ND	1700	ND	6900
1,2-Dichloropropane	ND	1700	ND	8100
cis-1,3-Dichloropropene	ND	1700	ND	7900

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Client Sample ID: SG-3

GC/MS Volatiles

Lot-Sample # H5G200167 - 001	Work Order # HFVXP1AD	Matrix AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	98 103 80	70 - 130 70 - 130 70 - 130

Qualifiers

J

Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

Client Sample ID: SG-4

GC/MS Volatiles

Lot-Sample # H5G200167 - 01	2	Work Order # HFV0F1	AD	Matrix: AIR
Date Sampled: 7/18/05 Prep Date: 7/21/05 Prep Batch #: 5203061		Date Received: 7/20/05 Analysis Date 7/22/05		
Prep Batch #: 5203061 Dilution Factor.: 18.4		Method ; TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3	REPORTING) LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	3.7	ND (17
Acetone	36	92	85 J)5 8.0 J	220
Ethylbenzene	1.8	3.7	8.0 J	16
2-Hexanone	ND	9.2	ND	38
Methylene chloride	ND	9.2	ND	32
Benzene	ND	3.7	ND	12
Styrene	2.0	3.7	8.3 J	16
1,1,2,2-Tetrachloroethane	ND	3.7	ND	25
Tetrachloroethene	ND	3.7	ND	25
Toluene	5.5	3.7	21	14
1,1,1-Trichloroethane	ND	3.7	ND	20
1,1,2-Trichloroethane	ND	3.7	ND	20
Trichloroethene	ND	3.7	ND	20
Vinyl chloride	ND	3.7	ND	9.4
o-Xylene	3.1	3.7	13 J	16
m-Xylene & p-Xylene	7.8	3.7	34	16
Bromodichloromethane	ND	3.7	ND	25
2-Butanone (MEK)	ND	9.2	ND	27
4-Methyl-2-pentanone (MIBK)	ND	9.2	ND	38
Bromoform	ND	3.7	ND	38
Bromomethane	ND	3.7	ND	14
Carbon disulfide	2.7	3.7	8.4 J	11
Carbon tetrachloride	ND	3.7	ND	23
Chlorobenzene	ND	3.7	ND	17
Dibromochloromethane	ND	3.7	ND	31
Chloroethane	ND	3.7	ND	9.7
Chloroform	14	3.7	71	18
Chloromethane	ND	9.2	ND	19
1,1-Dichloroethane	ND	3.7	ND	15
1,2-Dichloroethane	ND	3.7	ND	15
1,1-Dichloroethene	ND	3.7	ND	15
cis-1,2-Dichloroethene	ND	3.7	ND	15
trans-1,2-Dichloroethene	ND	3.7	ND	15
1,2-Dichloropropane	ND	3.7	ND	17
cis-1,3-Dichloropropene	ND	3.7	ND	17

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Client Sample ID: SG-4

GC/MS Volatiles

Lot-Sample #	H5G200167 - 012	Work Order #	HFV0F1AD	Matrix:	AIR
SURROGATE		PERCENT RECOVERY		LABORATORY CONTROL LIMITS (%)	
1,2-Dichloroeth Toluene-d8	ane-d4	125 104		70 - 130 70 - 130	
4-Bromofluorol	benzene	116		70 - 130	

Qualifiers

BMethod blank contamination. The associated method blank contains the target analyte at a reportable level.JEstimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

Client Sample ID: SG-5

GC/MS Volatiles

Lot-Sample # H5G200167 - 013	3	Work Order #	HFV0H1	AD	Matr	ix:	AIR
Date Sampled: 7/18/05 Prep Date: 7/21/05 Prep Batch #: 5203061		Date Received: Analysis Date	7/22/05				
Dilution Factor.: 15.9		Method	10-15				
PARAMETER	RESULTS (ppb(v/v))	REPORTI LIMIT (pr		RESUL	TS (ug/m3)	REPORTI LIMIT (ug	
trans-1,3-Dichloropropene	ND	3.2		ND		14	
Acetone	23-1.0	80	NA)	-54	🔾 کارنج	190	
Ethylbenzene	1.3	3.2		5.6	Ĵ	14	
2-Hexanone	ND	8.0		ND		33	
Methylene chloride	ND	8.0		ND		28	
Benzene	1.6	3.2		5.1	J	10	
Styrene	ND	3.2		ND		14	
1,1,2,2-Tetrachloroethane	ND	3.2		ND		22	
Tetrachloroethene	ND	3.2		ND		22	
Toluene	5.6	3.2		21		12	
1,1,1-Trichloroethane	ND	3.2		ND		17	
1,1,2-Trichloroethane	ND	3.2		ND		17	
Trichloroethene	ND	3.2		ND		17	
Vinyl chloride	ND	3.2		ND		8.1	
o-Xylene	1.6	3.2		6.9	J	14	
m-Xylene & p-Xylene	4.6	3.2		20		14	
Bromodichloromethane	ND	3.2		ND		21	
2-Butanone (MEK)	ND	8.0		ND		23	
4-Methyl-2-pentanone (MIBK)	ND	8.0		ND		33	
Bromoform	ND	3.2		ND		33	
Bromomethane	ND	3.2		ND		12	
Carbon disulfide	4.9	3.2		15		9.9	
Carbon tetrachloride	ND	3.2		ND		20	
Chlorobenzene	ND	3.2		ND		15	
Dibromochloromethane	ND	3.2		ND		27	
Chloroethane	ND	3.2		ND		8.4	
Chloroform	ND	3.2		ND		16	
Chloromethane	ND	8.0		ND		16	
1,1-Dichloroethane	ND	3.2		ND		13	
1,2-Dichloroethane	ND	3.2		ND		13	
1,1-Dichloroethene	ND	3.2		ND		13	
cis-1,2-Dichloroethene	ND	3.2		ND		13	
trans-1,2-Dichloroethene	ND	3.2		ND		13	Nº m
1,2-Dichloropropane	ND	3.2		ND		15	\$5. CV
cis-1,3-Dichloropropene	ND	3.2		ND		14	
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Client Sample ID: SG-5

GC/MS Volatiles

Lot-Sample # H5G200167 - 013	Work Order # HFV0	H1AD Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	124 105 112	70 - 130 70 - 130 70 - 130

Qualifiers

BMethod blank contamination. The associated method blank contains the target analyte at a reportable level.JEstimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

Client Sample ID: SG-6

GC/MS Volatiles

Lot-Sample # H5G200167 - 014		Work Order # HFV	V0Q1AD	Mat	trix:	AIR
Date Sampled: 7/18/05		Date Received: 7/20)/05			
Prep Date: 7/21/05		Analysis Date 7/21				
Prep Batch #: 5203061		-				
Dilution Factor.: 1.78		Method TO-	-15			
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v)) RESULT	S (ug/m3)	REPORTING LIMIT (ug/n	
trans-1,3-Dichloropropene	ND	0.36	ND		1.6	
Acetone	8.5	8.9	20	J ந	21	
Ethylbenzene	0.46	0.36	2.0	1	1.5	
2-Hexanone	ND	0.89	ND		3.6	
Methylene chloride	0.89	0.89	3.1		3.1	
Benzene	0.28	0.36	0.91	J	1.1	
Styrene	0.27	0.36	1.1	J	1.5	
1,1,2,2-Tetrachloroethane	ND	0.36	ND		2.4	
Tetrachloroethene	0.48	0.36	3.2		2.4	
Toluene	1.6	0.36	6.2		1.3	
1,1,1-Trichloroethane	ND	0.36	ND		1.9	
1,1,2-Trichloroethane	ND	0.36	ND		1.9	
Trichloroethene	ND	0.36	ND		1.9	
Vinyl chloride	ND	0.36	ND		0.91	
o-Xylene	0.64	0.36	2.8		1.5	
m-Xylene & p-Xylene	1.5	0.36	6.4		1.5	
Bromodichloromethane	ND	0.36	ND		2.4	
2-Butanone (MEK)	0.87	0.89	2.6	J	2.6	
4-Methyl-2-pentanone (MIBK)	0.56	0.89	2.3	J	3.6	
Bromoform	ND	0.36	ND		3.7	
Bromomethane	ND	0.36	ND		1.4	
Carbon disulfide	0.43	0.36	1.3		1.1	
Carbon tetrachloride	ND	0.36	ND		2.2	
Chlorobenzene	ND	0.36	ND		1.6	
Dibromochloromethane	ND	0.36	ND		3.0	
Chloroethane	ND	0.36	ND		0.94	
Chloroform	ND	0.36	ND		1.7	
Chloromethane	ND	0.89	ND		1.8	
1,1-Dichloroethane	ND	0.36	ND		1.4	
1,2-Dichloroethane	ND	0.36	ND		1.4	
1,1-Dichloroethene	ND	0.36	ND		1.4	
cis-1,2-Dichloroethene	ND	0.36	ND		1.4	
trans-1,2-Dichloroethene	ND	0.36	ND		1.4	
1,2-Dichloropropane	ND	0.36	ND		1.6	12
cis-1,3-Dichloropropene	ND	0.36	ND		1.6	- (Y

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Client Sample ID: SG-6

GC/MS Volatiles

Lot-Sample # H5G200	167 - 014 Work Order	# HFV0Q1AD	Matrix: AIR
SURROGATE	PERCENT RECOVER	Y	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8	123 104		70 - 130 70 - 130
4-Bromofluorobenzene	117		70 - 130

В

J

Method blank contamination. The associated method blank contains the target analyte at a reportable level. Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

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Client Sample ID: FD-2

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GC/MS Volatiles

Lot-Sample # H5G200167 - 01	6	Work Order # HFV051A	7.2	1720411	•ix Al	
Date Sampled: 7/18/05 Prep Date: 7/25/05 Prep Batch #: 5206288		Date Received: 7/20/05 Analysis Date 7/25/05				
Prep Batch #: 5206288 Dilution Factor.: 18.2		Method: TO-15				
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS	(ug/m3)	REPORTING LIMIT (ug/m3))
trans-1,3-Dichloropropene	ND	3.6	ND		17	
Acetone	.6.3 ND	91 15) 15-	JB 🔾	220	
Ethylbenzene	ND	3.6	ND	. ,	16	
2-Hexanone	ND	9.1	ND		37	
Methylene chloride	ND	9.1	ND		32	
Benzene	ND	3.6	ND		12	
Styrene	ND	3.6	ND		16	
1,1,2,2-Tetrachloroethane	ND	3.6	ND		25	
Fetrachloroethene	ND	3.6	ND		25	
Foluene	1.7	3.6	6.2	J	14	
,1,1-Trichloroethane	ND	3.6	ND		20	
1,1,2-Trichloroethane	ND	3.6	ND		20	
Frichloroethene	ND	3.6	ND		20	
Vinyl chloride	ND	3.6	ND		9.3	
o-Xylene	ND	3.6	ND		16	
m-Xylene & p-Xylene	ND	3.6	ND		16	
Bromodichloromethane	ND	3.6	ND		24	
2-Butanone (MEK)	ND	9.1	ND		27	
4-Methyl-2-pentanone (MIBK)	ND	9.1	ND		37	
Bromoform	ND	3.6	ND		38	
Bromomethane	ND	3.6	ND		14	
Carbon disulfide	ND	3.6	ND		11	
Carbon tetrachloride	ND	3.6	ND		23	
Chlorobenzene	ND	3.6	ND		17	
Dibromochloromethane	ND	3.6	ND		31	
Chloroethane	ND	3.6	ND		9.6	
Chloroform	ND	3.6	ND		18	
Chloromethane	ND	9.1	ND		19	
1,1-Dichloroethane	ND	3.6	ND		15	
1,2-Dichloroethane	ND	3.6	ND		15	
1,1-Dichloroethene	ND	3.6	ND		14	
cis-1,2-Dichloroethene	ND	3.6	ND		14	
trans-1,2-Dichloroethene	ND	3.6	ND		14	
1,2-Dichloropropane	ND	3.6	ND		17	
1. 1. 2. D'ablemennenen	NT	3.6	ND		17	

3.6

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ND

cis-1,3-Dichloropropene

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ND

Client Sample ID: FD-2

GC/MS Volatiles

Lot-Sample # H5G200167 - 016	Work Order #	HFV051AD	Matrix: AIR	
SURROGATE	PERCENT RECOVERY		LABORATORY CONTROL LIMITS (%)	
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	113 94 81		70 - 130 70 - 130 70 - 130	

Qualifiers

В

J

Method blank contamination. The associated method blank contains the target analyte at a reportable level. Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

Client Sample ID: SG-7

GC/MS Volatiles

Lot-Sample # H5G200167 - 01	1	Work Order # HFV0E1.	AD Mat	rix: AIR
Date Sampled: 7/18/05		Date Received: 7/20/05		
Prep Date: 7/21/05		Analysis Date 7/22/05		
Prep Batch #: 5203061 Dilution Factor.: 1.93		Method TO-15		
Different Pactorn. 1990				
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	0.39	ND	1.8
Acetone	ND	9.6	ND	23
Ethylbenzene	0.38	0.39	1.6 J)	1.7
2-Hexanone	ND	0.96	ND	4.0
Methylene chloride	ND	0.96	ND	3.4
Benzene	ND	0.39	ND	1.2
Styrene	ND	0.39	ND	1.6
1,1,2,2-Tetrachloroethane	ND	0.39	ND	2.6
Tetrachloroethene	0.56	0.39	3.8 5	2.6
Toluene	1.6	0.39	5.9 🖒	1.5
1,1,1-Trichloroethane	ND	0.39	ND	2.1
1,1,2-Trichloroethane	ND	0.39	ND	2.1
Trichloroethene	ND	0.39	ND	2.1
Vinyl chloride	ND	0.39	ND	0.99
o-Xylene	0.45	0.39	2.0	1.7
m-Xylene & p-Xylene	1.2	0.39	5.2	1.7
Bromodichloromethane	ND	0.39	ND	2.6
2-Butanone (MEK)	ND	0.96	ND	2.8
4-Methyl-2-pentanone (MIBK)	ND	0.96	ND	4.0
Bromoform	ND	0.39	ND	4.0
Bromomethane	ND	0.39	ND	1.5
Carbon disulfide	ND	0.39	ND	1.2
Carbon tetrachloride	ND	0.39	ND	2.4
Chlorobenzene	ND	0.39	ND	1.8
Dibromochloromethane	ND	0.39	ND	3.3
Chloroethane	ND	0.39	ND	1.0
Chloroform	ND	0.39	ND	1.9
Chloromethane	ND	0.96	ND	2.0
1,1-Dichloroethane	ND	0.39	ND	1.6
1,2-Dichloroethane	ND	0.39	ND	1.6
1,1-Dichloroethene	ND	0.39	ND	1.5
cis-1,2-Dichloroethene	ND	0.39	ND	1.5
trans-1,2-Dichloroethene	ND	0.39	ND	1.5 s.
1,2-Dichloropropane	ND	0.39	ND	1.8
cis-1,3-Dichloropropene	ND	0.39	ND	- 1.8

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Client Sample ID: SG-7

GC/MS Volatiles

Lot-Sample # H5G200167 - 011	Work Order #	HFV0E1AD	Matrix:	AIR
SURROGATE	PERCENT RECOVERY		LABORATOR CONTROL LIMITS (%)	Ŷ
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	148 * 113 108		70 - 130 70 - 130 70 - 130	

Qualifiers

* Surrogate recovery is outside stated control limits.

J Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

Client Sample ID: SG-7

GC/MS Volatiles

Date Sampled: 7/18/05 Prep Date: 7/21/05		ate Received: 7/20/05 nalysis Date 7/22/05		
Prep Batch #: 5203061 Dilution Factor.: 19.3		ethod: TO-15		/
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3)	REPORTING LIMIT (ug/m3)
rans-1,3-Dichloropropene	ND	3.9	ND /	18
Acetone	13	96	31 XB	230
Ethylbenzene	ND	3.9	ND	17
2-Hexanone	ND	9.6	ND /	40
Methylene chloride	ND	9.6	ND /	34
Benzene	ND	3.9	ND /	12
Styrene	ND	3.9	ND /	16
,1,2,2-Tetrachloroethane	ND	3.9	ND	26
Fetrachloroethene	ND	3.9	D	26
Foluene	2.3	3.9	/8.6 J	15
1,1,1-Trichloroethane	ND	3.9	ND	21
,1,2-Trichloroethane	ND	3.9	ND	21
[richloroethene	ND	3.9	ND	21
vinyl chloride	ND	3.9	ND	9.9
-Xylene	ND	3.9	ND	17
n-Xylene & p-Xylene	1.5	3.9	6.7 J	17
Bromodichloromethane	ND	3.9	ND	26
2-Butanone (MEK)	ND	9.6 /	ND	28
4-Methyl-2-pentanone (MIBK)	ND	9,6	ND	40
Bromoform	ND	3.9	ND	40
Bromomethane	ND	/ 3.9	ND	15
Carbon disulfide	2.6	3.9	8.1 J	12
Carbon tetrachloride	ND	3.9	ND	24
Chlorobenzene	ND	3.9	ND	18
Dibromochloromethane	ND /	3.9	ND	33
Chloroethane	ND	3.9	ND	10
Chloroform	ND /	3.9	ND	19
Chloromethane	ND	9.6	ND	20
,1-Dichloroethane	ND	3.9	ND	16
,2-Dichloroethane	ŅĎ	3.9	ND	16
,1-Dichloroethene	ND	3.9	ND	15
is-1,2-Dichloroethene	ND	3.9	ND	15
rans-1,2-Dichloroethene	ND	3.9	ND	15
,2-Dichloropropane	ND	3.9	ND	18
cis-1,3-Dichloropropene	ND	3.9	ND	18

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Client Sample ID: SG-7

GC/MS Volatiles

Lot-Sample # H5G200167 - 011	Work Order # HFV0E2AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	125 102 116	70 - 130 70 - 130 70 - 130

Oualifiers

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.
 J Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

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Client Sample ID: SG-8

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GC/MS Volatiles

Lot-Sample # H5G200167 - 00	9	Work Order #	HFV0A1AD	Mat	rix: AIR
Date Sampled: 7/18/05		Date Received:	7/20/05		
Prep Date 7/25/05		Analysis Date			
Prep Batch #: 5206288					
Dilution Factor.: 18		Method:	TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTII LIMIT (pp		°S (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	3.6	ND		16
Acetone	- 31 + X)	90	1.1.2 73	JB	210
Ethylbenzene	3.3	3.6	14	J	16
2-Hexanone	ND	9.0	ND		37
Methylene chloride	ND	9.0	ND		31
Benzene	7.5	3.6	24		12
Styrene	ND	3.6	ND		15
1,1,2,2-Tetrachloroethane	ND	3.6	ND		25
Tetrachloroethene	ND	3.6	ND		24
Toluene	36	3.6	140		14
1,1,1-Trichloroethane	ND	3.6	ND		20
1,1,2-Trichloroethane	ND	3.6	ND		20
Trichloroethene	ND	3.6	ND		19
Vinyl chloride	ND	3.6	ND		9.2
o-Xylene	3.2	3.6	14	J	16
m-Xylene & p-Xylene	9.3	3.6	41		16
Bromodichloromethane	ND	3.6	ND		24
2-Butanone (MEK)	ND	9.0	ND		27
4-Methyl-2-pentanone (MIBK)	ND	9.0	ND		37
Bromoform	ND	3.6	ND		37
Bromomethane	ND	3.6	ND		14
Carbon disulfide	22	3.6	68		11
Carbon tetrachloride	ND	3.6	ND		23
Chlorobenzene	ND	3.6	ND		17
Dibromochloromethane	ND	3.6	ND		31
Chloroethane	ND	3.6	ND		9.5
Chloroform	ND	3.6	ND		18
Chloromethane	ND	9.0	ND		19
1,1-Dichloroethane	ND	3.6	ND		15
1,2-Dichloroethane	ND	3.6	ND		15
1,1-Dichloroethene	ND	3.6	ND		14
cis-1,2-Dichloroethene	ND	3.6	ND		14
trans-1,2-Dichloroethene	ND	3.6	ND		14
1,2-Dichloropropane	ND	3.6	ND		17
cis-1,3-Dichloropropene	ND	3.6	ND		16

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Client Sample ID: SG-8

GC/MS Volatiles

Lot-Sample # H5G200167 - 009	Work Order # HFV0A1AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	115 93 80	70 - 130 70 - 130 70 - 130

Qualifiers

BMethod blank contamination. The associated method blank contains the target analyte at a reportable level.JEstimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

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Client Sample ID: SG-9

GC/MS Volatiles

Lot-Sample # H5G200167 - 01	0	Work Order #	HFV0D1A	D	Matr	ix AIR	
Date Sampled: 7/18/05 Prep Date: 7/25/05 Prep Batch #: 5206288		Date Received: Analysis Date	7/20/05 7/25/05				
Prep Batch #: 5206288 Dilution Factor.: 17		Method:	TO-15				
PARAMETER	RESULTS (ppb(v/v))	REPORTI LIMIT (pr		RESULT	ſS (ug/m3)	REPORTING LIMIT (ug/m3)	
trans-1,3-Dichloropropene	ND	3.4		ND		15	
Acetone	6.1 N.X.	85	CK-1	14	J¥ U	200	
Ethylbenzene	ND	3.4		ND	· •	15	
2-Hexanone	ND	8.5		ND		35	
Methylene chloride	ND	8.5		ND		30	
Benzene	ND	3.4		ND		11	
Styrene	ND	3.4		ND		14	
1,1,2,2-Tetrachloroethane	ND	3.4		ND		23	
Tetrachloroethene	5.2	3.4		35		23	
Toluene	3.4	3.4		13		13	
1,1,1-Trichloroethane	ND	3.4		ND		19	
1,1,2-Trichloroethane	ND	3.4		ND		19	
Trichloroethene	2.3	3.4		13	J	18	
Vinyl chloride	ND	3.4		ND		8.7	
o-Xylene	ND	3.4		ND		15	
m-Xylene & p-Xylene	1.8	3.4		8.0	J	15	
Bromodichloromethane	ND	3.4		ND		23	
2-Butanone (MEK)	ND	8.5		ND		25	
4-Methyl-2-pentanone (MIBK)	ND	8.5		ND		35	
Bromoform	ND	3.4		ND		35	
Bromomethane	ND	3.4		ND		13	
Carbon disulfide	1.5	3.4		4.6	J	11	
Carbon tetrachloride	ND	3.4		ND		21	
Chlorobenzene	ND	3.4		ND		16	
Dibromochloromethane	ND	3.4		ND		29	
Chloroethane	ND	3.4		ND		9.0	
Chloroform	8.0	3.4		39		17	
Chloromethane	ND	8.5		ND		18	
1,1-Dichloroethane	ND	3.4		ND		14	
1,2-Dichloroethane	ND	3.4		ND		14	
1,1-Dichloroethene	ND	3.4		ND		13	
cis-1,2-Dichloroethene	ND	3.4		ND		13	
trans-1,2-Dichloroethene	ND	3.4		ND		13	
1,2-Dichloropropane	ND	3.4		ND		16	
cis-1,3-Dichloropropene	ND	3.4		ND		15	ĺ,

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Client Sample ID: SG-9

GC/MS Volatiles

Lot-Sample # H5G200167 - 010	Work Order #	HFV0D1AD	Matrix AIR
SURROGATE	PERCENT RECOVERY		LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	116 95 82		70 - 130 70 - 130 70 - 130

BMethod blank contamination. The associated method blank contains the target analyte at a reportable level.JEstimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

Client Sample ID: SG-10

GC/MS Volatiles

Lot-Sample # H5G200167 - 003	3	Work Order # HF	VX81AD	Mat	rix: AIR
Date Sampled: 7/18/05 Prep Date 7/25/05		Date Received: 7/2 Analysis Date 7/2	0/05 5/05		
Prep Batch #: 5206288 Dilution Factor.: 19.1		Method TC	0-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/	(v)) <u>RE</u>	SULTS (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	3.8	NI)	17
Acetone	7.5 (X)	96	- N 18	- X X -	230
Ethylbenzene	ND	3.8	NI		17
2-Hexanone	ND	9.6	NI)	39
Methylene chloride	4.9	9.6	17	J	33
Benzene	ND	3.8	NI)	12
Styrene	ND	3.8	NI	2	16
1,1,2,2-Tetrachloroethane	ND	3.8	NI	D	26
Tetrachloroethene	ND	3.8	N	2	26
Toluene	2.0	3.8	7.4	4 J	14
1,1,1-Trichloroethane	ND	3.8	N	D	21
1,1,2-Trichloroethane	ND	3.8	N	D	21
Trichloroethene	ND	3.8	N	D	21
Vinyl chloride	ND	3.8	N	D	9.8
o-Xylene	ND	3.8	N		17
m-Xylene & p-Xylene	2.1	3.8	9.		17
Bromodichloromethane	ND	3.8	N		26
2-Butanone (MEK)	ND	9.6	N		28
4-Methyl-2-pentanone (MIBK)	ND	9.6	N		39
Bromoform	ND	3.8	N		39
Bromomethane	ND	3.8	N		15
Carbon disulfide	3.6	3.8	11		12
Carbon tetrachloride	ND	3.8	N		24
Chlorobenzene	ND	3.8	N		18
Dibromochloromethane	ND	3.8		D	33
Chloroethane	ND	3.8	N		10
Chloroform	1.8	3.8	9.		19
Chloromethane	ND	9.6		D	20
1,1-Dichloroethane	ND	3.8		D	15
1,2-Dichloroethane	ND	3.8		D	15
1,1-Dichloroethene	ND	3.8		ID	15
cis-1,2-Dichloroethene	ND	3.8		ID	15
trans-1,2-Dichloroethene	ND	3.8		ID	15
1,2-Dichloropropane	ND	3.8		ID	18
cis-1,3-Dichloropropene	ND	3.8	Ν	ID	17

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Client Sample ID: SG-10

GC/MS Volatiles

Lot-Sample # H5G200167 - 008	Work Order # HFVX81AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	111 94 82	70 - 130 70 - 130 70 - 130

Qualifiers

BMethod blank contamination. The associated method blank contains the target analyte at a reportable level.JEstimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

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Client Sample ID: FD-1

GC/MS Volatiles

		GC/MS Vol	atiles			1	1
Lot-Sample # H5G200167 - 015		Work Order #	HFV001A	D	Matrix.		AIR
Date Sampled: 7/18/05 Prep Date: 7/21/05 Prep Batch #: 5203061		Date Received: Analysis Date					
Dilution Factor.: 1.81		Method	TO-15				
PARAMETER	RESULTS (ppb(v/v))	REPORTI LIMIT (pr		RESULTS (ug/m		EPORTI IMIT (ug	
trans-1,3-Dichloropropene	NR			NR			
Acetone	NR			NR	, Alexandre de la construcción d		
Ethylbenzene	NR			NR	1		
2-Hexanone	NR			NR	1		
Methylene chloride	NR			NR /			
Benzene	NR			NR			
Styrene	NR			NR /			
1,1,2,2-Tetrachloroethane	NR			NR /			
Tetrachloroethene	NR			NR			
Toluene	NR			ŊŔ			
1,1,1-Trichloroethane	NR			NR			
1,1,2-Trichloroethane	NR		/	NR			
Trichloroethene	NR			NR			
Vinyl chloride	NR			NR			
o-Xylene	NR			NR			
m-Xylene & p-Xylene	NR	,	/	NR			
Bromodichloromethane	NR	/		NR			
2-Butanone (MEK)	NR			NR			
4-Methyl-2-pentanone (MIBK)	NR	1		NR			
Bromoform	NR	1		NR			
Bromomethane	NR	1		NR			
Carbon disulfide	NR			NR			
Carbon tetrachloride	NR			NR			
Chlorobenzene	NR			NR			
Dibromochloromethane	NR			NR			
Chloroethane	NR			NR			
Chloroform	NR			NR			
Chloromethane	NR			NR			
1,1-Dichloroethane	NR			NR			
1,2-Dichloroethane	NR			NR			
1,1-Dichloroethene	NR			NR			
cis-1,2-Dichloroethene	NR			NR			
trans-1,2-Dicbloroethene	NR			NR			
1,2-Dichloropropane	NR			NR			
cis-1,3-Dichloropropene	NR			NR			
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Client Sample ID: FD-1

GC/MS Volatiles

Lot-Sample # H5G200167 - 015	Work Order # HFV001AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	0.0 * 58 * 102	70 - 130 70 - 130 70 - 130
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Oualifiers

* Surrogate recovery is outside stated control limits.NR Not reportable.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

Client Sample ID: FD-1

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GC/MS Volatiles

Lot-Sample # H5G200167 - 01	5	Work Order # HFV002	AD	Matr	ix:	AIR
Date Sampled: 7/18/05 Prep Date: 7/21/05 Prep Batch #: 5203061		Date Received 7/20/05 Analysis Date 7/22/05				
Dilution Factor.: 18.1		Method TO-15				
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULT	°S (ug/m3)	REPORTIN LIMIT (ug/	
trans-1,3-Dichloropropene	ND	3.6	ND		16	
Acetone	-31 NA	90 5	74		210	
Ethylbenzene	ND	3.6	ND	-	16	
2-Hexanone	ND	9.0	ND		37	
Methylene chloride	ND	9.0	ND		31	
Benzene	ND	3.6	ND		12	
Styrene	ND	3.6	ND		15	
1,1,2,2-Tetrachloroethane	ND	3.6	ND		25	
Tetrachloroethene	4.1	3.6	28		25	
Toluene	2.9	3.6	11	J	14	
1,1,1-Trichloroethane	ND	3.6	ND		20	
1,1,2-Trichloroethane	ND	3.6	ND		20	
Trichloroethene	ND	3.6	ND		19	
Vinyl chloride	ND	3.6	ND		9.3	
o-Xylene	1.6	3.6	6.8	J	16	
m-Xylene & p-Xylene	4.3	3.6	19		16	
Bromodichloromethane	ND	3.6	ND		24	
2-Butanone (MEK)	ND	9.0	ND		27	
4-Methyl-2-pentanone (MIBK)	ND	9.0	ND		37	
Bromoform	ND	3.6	ND		37	
Bromomethane	ND	3.6	ND		14	
Carbon disulfide	5.0	3.6	16		11 23	
Carbon tetrachloride	ND	3.6	ND			
Chlorobenzene	ND	3.6	ND		17 31	
Dibromochloromethane	ND	3.6	ND		9.6	
Chloroethane	ND	3.6	ND	Ŧ		
Chloroform	2.9	3.6	14 ND	J	18 19	
Chloromethane	ND	9.0			15	
1,1-Dichloroethane	ND	3.6	ND ND		15	
1,2-Dichloroethane	ND	3.6	ND ND		13 14	
1,1-Dichloroethene	ND	3.6			14	
cis-1,2-Dichloroethene	ND	3.6	ND ND		14 14	
trans-1,2-Dichloroethene	ND	3.6			14	
1,2-Dichloropropane	ND	3.6	ND		16	\sim
cis-1,3-Dichloropropene	ND	3.6	ND		10	1.

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Client Sample ID: FD-1

GC/MS Volatiles

Lot-Sample # H5G2001	67 - 015 Work Order #	HFV002AD	Matrix:	AIR
SURROGATE	PERCENT RECOVERY		LABORATORY CONTROL LIMITS (%)	{
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	122 105 116		70 - 130 70 - 130 70 - 130	

Oualifiers

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.
 J Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

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Client Sample ID: SG-11

GC/MS Volatiles

Lot-Sample # H5G200167 - 007	7	Work Order #	HFVX71A	D	Matr	ix:	AIR
Date Sampled: 7/18/05 Prep Date 7/25/05		Date Received: Analysis Date	7/20/05 7/25/05				
Prep Batch #: 5206288 Dilution Factor.: 17.7		Method	: TO-15				
PARAMETER	RESULTS (ppb(v/v))	REPORT LIMIT (p		RESUL	TS (ug/m3)	REPORTIN LIMIT (ug	
trans-1,3-Dichloropropene	ND	3.5	1 1	ND		16	
Acetone	-13 ND	88	12	-32	XB ()	210	
Ethylbenzene	1.7	3.5		7.6	J	15	
2-Hexanone	ND	8.8		ND	_	36	
Methylene chloride	7.3	8.8		25	J	31	
Benzene	1.5	3.5		4.9	J	11	
Styrene	ND	3.5		ND		15	
1,1,2,2-Tetrachloroethane	ND	3.5		ND		24	
Tetrachloroethene	ND	3.5		ND		24	
Toluene	4.9	3.5		19		13	
1,1,1-Trichloroethane	ND	3.5		ND		19	
1,1,2-Trichloroethane	ND	3.5		ND		19	
Trichloroethene	ND	3.5		ND		19	
Vinyl chloride	ND	3.5		ND		9.0	
o-Xylene	1.9	3.5		8.2	1	15	
m-Xylene & p-Xylene	5.9	3.5		26		15	
Bromodichloromethane	ND	3.5		ND		24	
2-Butanone (MEK)	ND	8.8		ND		26	
4-Methyl-2-pentanone (MIBK)	ND	8.8		ND		36	
Bromoform	ND	3.5		ND		37	
Bromomethane	ND	3.5		ND		14	
Carbon disulfide	3.1	3.5		9.5	J	11	
Carbon tetrachloride	ND	3.5		ND		22	
Chlorobenzene	ND	3.5		ND		16	
Dibromochloromethane	ND	3.5		ND		30	
Chloroethane	ND	3.5		ND		9.3	
Chloroform	ND	3.5		ND		17	
Chloromethane	ND	8.8		ND		18	
1,1-Dichloroethane	ND	3.5		ND		14	
1,2-Dichloroethane	ND	3.5		ND		14	
1,1-Dichloroethene	ND	3.5		ND		14	
cis-1,2-Dichloroethene	ND	3.5		ND		14	
trans-1,2-Dichloroethene	ND	3.5		ND		14	
1,2-Dichloropropane	ND	3.5		ND		16	
cis-1,3-Dichloropropene	ND	3.5		ND		16	Č, Š

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Client Sample ID: SG-11

GC/MS Volatiles

Lot-Sample # H5G200167 - 007	Work Order # HFVX7	1AD Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	112 96 83	70 - 130 70 - 130 70 - 130

Qualifiers

BMethod blank contamination. The associated method blank contains the target analyte at a reportable level.JEstimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

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Client Sample ID: SG-12

GC/MS Volatiles

Lot-Sample # H5G200167 - 005	5	Work Order # HFVX41	AD Ma	atrix: AIR
Date Sampled: 7/18/05 Prep Date: 7/21/05		Date Received: 7/20/05 Analysis Date 7/21/05		
Prep Batch #: 5203121 Dilution Factor.: 1.8		Method: TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v/v))	RESULTS (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	0.36	ND U.J	1.6
Acetone	12	9.0	28 J	21
Ethylbenzene	0.40	0.36	1.8	1.6
2-Hexanone	ND	0.90	ND	3.7
Methylene chloride	ND	0.90	ND	3.1
Benzene	0.59	0.36	1.9	1.2
Styrene	ND	0.36	ND CAS	1.5
1,1,2,2-Tetrachloroethane	ND	0.36	ND	2.5
Tetrachloroethene	1.2	0.36	8.0	2.4
Toluene	4.2	0.36	16	1.4
1,1,1-Trichloroethane	0.11	0.36	0.60 J	2.0
1,1,2-Trichloroethane	ND	0.36	ND	2.0
Trichloroethene	ND	0.36	ND	1.9
Vinyl chloride	ND	0.36	ND	0.92
o-Xylene	0.29	0.36	1.3 J	1.6
m-Xylene & p-Xylene	0.78	0.36	3.4	1.6
Bromodichloromethane	ND	0.36	ND	2.4
2-Butanone (MEK)	0.97	0.90	2.9	2.7
4-Methyl-2-pentanone (MIBK)	0.48	0.90	2.0 J	3.7
Bromoform	ND	0.36	ND 55	3.7
Bromomethane	ND	0.36	ND	1.4
Carbon disulfide	4.8	0.36	15	1.1
Carbon tetrachloride	ND	0.36	ND	2.3
Chlorobenzene	ND	0.36	ND	1.7
Dibromochloromethane	ND	0.36	ND	3.1
Chloroethane	ND	0.36	ND	0.95
Chloroform	1.6	0.36	7.9	1.8
Chloromethane	0.28	0.90	0.57 J	1.9
1,1-Dichloroethane	ND	0.36	ND	1.5
1,2-Dichloroethane	ND	0.36	ND	1.5
1,1-Dichloroethene	ND	0.36	ND	1.4
cis-1,2-Dichloroethene	ND	0.36	ND ·	1.4
trans-1,2-Dichloroethene	ND	0.36	ND	1.4
1,2-Dichloropropane	ND	0.36	ND	1.7
cis-1,3-Dichloropropene	ND	0.36	ND	1.6

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Client Sample ID: SG-12

GC/MS Volatiles

Lot-Sample #	H5G200167 - 005	Work Order #	HFVX41AD	Matrix:	AIR
SURROGATE		PERCENT RECOVERY		LABORATORY CONTROL LIMITS (%)	-
1,2-Dichloroethan Toluene-d8 4-Bromofluorober		112 89 86		70 - 130 70 - 130 70 - 130	

Qualifiers

BMethod blank contamination. The associated method blank contains the target analyte at a reportable level.JEstimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

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Client Sample ID: SG-13

GC/MS Volatiles

Lot-Sample # H5G200167 - 006	5	Work Order #	HFVX62AD	Matr	ix: AIR
Date Sampled: 7/18/05 Prep Date: 7/22/05		Date Received: Analysis Date			
Prep Batch #: 5206164 Dilution Factor.: 17.5		Method:	TO-15		
PARAMETER	RESULTS (ppb(v/v))	REPORTIN LIMIT (pp		TS (ug/m3)	REPORTING LIMIT (ug/m3)
trans-1,3-Dichloropropene	ND	3.5	ND		16
Acetone	· 12 · · · ·	88	K.A.) 29	ت کور –	5210
Ethylbenzene	ND	3.5	ND		15
2-Hexanone	ND	8.8	ND	a. The second	36
Methylene chloride	ND	8.8	ND	153 L	30
Benzene	ND	3.5	ND		11
Styrene	ND	3.5	ND		15
1,1,2,2-Tetrachloroethane	ND	3.5	ND		24
Tetrachloroethene	ND	3.5	ND		24
Toluene	4.0	3.5	15		13
1,1,1-Trichloroethane	ND	3.5	ND		19
1,1,2-Trichloroethane	ND	3.5	ND		19
Trichloroethene	ND	3.5	ND		19
Vinyl chloride	ND	3.5	ND ·	55	8.9
o-Xylene	ND	3.5	ND		15
m-Xylene & p-Xylene	1.8	3.5	7.8	J	15
Bromodichloromethane	ND	3.5	ND		23
2-Butanone (MEK)	ND	8.8	ND		26
4-Methyl-2-pentanone (MIBK)	ND	8.8	ND		36
Bromoform	ND	3.5	ND		36
Bromomethane	ND	3.5	ND	and the second	14
Carbon disulfide	2.4	3.5	7.6	, ≭ ∖ ∖	11
Carbon tetrachloride	ND	3.5	ND		22
Chlorobenzene	ND	3.5	ND		16
Dibromochloromethane	ND	3.5	ND		30
Chloroethane	ND	3.5	ND		9.2
Chloroform	ND	3.5	ND	e	17
Chloromethane	ND	8.8	ND	(S)	18
1,1-Dichloroethane	ND	3.5	ND		14
1,2-Dichloroethane	ND	3.5	ND	·	14
1,1-Dichloroethene	ND	3.5	ND	C12	14
cis-1,2-Dichloroethene	ND	3.5	ND		14
trans-1,2-Dichloroethene	ND	3.5	ND		14
1,2-Dichloropropane	ND	3.5	ND		16
cis-1,3-Dichloropropene	ND	3.5	ND		16 (

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Client Sample ID: SG-13

GC/MS Volatiles

Lot-Sample # H5G200167 - 006	Work Order # HFVX62AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8	123 94	70 - 130 70 - 130
4-Bromofluorobenzene	78	70 - 130

Oualifiers

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Method blank contamination. The associated method blank contains the target analyte at a reportable level. Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

Client Sample ID: SG-13

GC/MS Volatiles

Lot-Sample # H5G200167 - 006	5	Work Order #	HFVX61A	AD Ma	trix: AIR
Date Sampled: 7/18/05		Date Received:	7/20/05		
Prep Date: 7/21/05		Analysis Date	7/21/05		
Prep Batch #: 5203121					
Dilution Factor.: 1.75		Method	: TO-15		
	RESULTS	REPORT	NG		REPORTING
PARAMETER	(ppb(v/v))	LIMIT (p		RESULTS (ug/m3)	LIMIT (ug/p13)
trans-1,3-Dichloropropene	NR			NR	
Acetone	NR			NR	
Ethylbenzene	NR			NR	
2-Hexanone	NR			NR	
Methylene chloride	NR			NR /	/
Benzene	NR			NR	
Styrene	NR			NR	
1,1,2,2-Tetrachloroethane	NR			NR	
Tetrachloroethene	NR			NR	
Toluene	NR			NR	
1,1,1-Trichloroethane	NR			NR	
1,1,2-Trichloroethane	NR			NR	
Trichloroethene	NR			NR	
Vinyl chloride	NR		/	/ NR	
o-Xylene	NR			NR	
m-Xylene & p-Xylene	NR			NR	
Bromodichloromethane	NR			NR	
2-Butanone (MEK)	NR			NR NR	
4-Methyl-2-pentanone (MIBK)	NR	/	/	NR	
Bromoform	NR			NR	
Bromomethane	NR	/		NR	
Carbon disulfide	NR			NR	
Carbon tetrachloride	NR NR	/		NR	
Chlorobenzene	NR			NR	
Dibromochloromethane	NR	1		NR	
Chloroethane Chloroform	NR	1		NR	
Chloromethane	NR /	1. 1		NR	
1,1-Dichloroethane	NR			NR	
1,2-Dichloroethane	NR			NR	
1,1-Dichloroethene	NR			NR	
cis-1,2-Dichloroethene	NŔ			NR	
trans-1,2-Dichloroethene	ŃR			NR	
1,2-Dichloropropane	NR			NR	
cis-1,3-Dichloropropene	NR			NR	

Client Sample ID: SG-13

GC/MS Volatiles

Lot-Sample # H5G200167 - 006	Work Order # HFVX61AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	0.0 * 0.0 * 0.0 *	70 - 130 70 - 130 70 - 130

Qualifiers

Surrogate recovery is outside stated control limits.
 NR Not reportable.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

Client Sample ID: AB-1

GC/MS Volatiles

Lot-Sample # H5G200167 - 00	4	Work Order # Hl	FVX12AD	M	atrix:	AIR
Date Sampled: 7/18/05 Prep Date 7/22/05		Date Received: 7/2 Analysis Date 7/2				
Prep Batch #: 5206164 Dilution Factor.: 15.5	1	Method: To	O-15			
PARAMETER	RESULTS (ppb(v/v))	REPORTING LIMIT (ppb(v		JLTS (ug/m3)	REPORTI LIMIT (ug	
trans-1,3-Dichloropropene	ND	3.1	ND		14	
Acetone	7.6 p.X)	78	<u>()</u> 18	AB C	5 180	
Ethylbenzene	ND '	3.1	ND		13	
2-Hexanone	ND	7.8	ND		32	
Methylene chloride	ND	7.8	ND	85° 2	27	
Benzene	ND	3.1	ND		9.9	
Styrene	ND	3.1	ND		13	
1,1,2,2-Tetrachloroethane	ND	3.1	ND		21	
Tetrachloroethene	ND	3.1	ND		21	
Toluene	ND	3.1	ND		12	
1,1,1-Trichloroethane	ND	3.1	ND		17	
1,1,2-Trichloroethane	ND	3.1	ND		17	
Trichloroethene	ND	3.1	ND		17	
Vinyl chloride	ND	3.1	ND	Cí)	7.9	
o-Xylene	ND	3.1	ND		13	
m-Xylene & p-Xylene	ND	3.1	ND		13	
Bromodichloromethane	ND	3.1	ND		21	
2-Butanone (MEK)	ND	7.8	ND		23	
4-Methyl-2-pentanone (MIBK)	ND	7.8	ND		32	
Bromoform	ND	3.1	ND		32	
Bromomethane	ND	3.1	ND	~~~~	12	
Carbon disulfide	ND	3.1	ND	NT 5	9.7	
Carbon tetrachloride	ND	3.1	ND		20	
Chlorobenzene	ND	3.1	ND		14	
Dibromochloromethane	ND	3.1	ND		26	
Chloroethane	ND	3.1	ND		8.2	
Chloroform	ND	3.1	ND		15	
Chloromethane	ND	7.8	ND	155	16	
1,1-Dichloroethane	ND	3.1	ND		13	
1,2-Dichloroethane	ND	3.1	ND		13	
1,1-Dichloroethene	ND	3.1	ND	55	12	
cis-1,2-Dichloroethene	ND	3.1	ND		12	
trans-1,2-Dichloroethene	ND	3.1	ND		12	
1,2-Dichloropropane	ND	3.1	ND		14	•.
cis-1,3-Dichloropropene	ND	3.1	ND		14	1

9/14/2004

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Client Sample ID: AB-1

GC/MS Volatiles

Lot-Sample # H5G200167 - 00	4 Work Order # HFVX12AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	115 94 77	70 - 130 70 - 130 70 - 130

Oualifiers

B Method blank contamination. The associated method blank contains the target analyte at a reportable level.
 J Estimated result. Result is less than RL.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24.45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)

TO-14 Conv Rev. 5

Client Sample ID: AB-1

GC/MS Volatiles

		Client Sample II	D: AB-1		1
		GC/MS Vol	atiles		
Lot-Sample # H5G200	0167 - 004	Work Order #	HFVX11AD	Matrix:	AIR
Prep Date 7/2	8/05 1/05	Date Received: Analysis Date	7/20/05 7/21/05		
Prep Batch #: 520 Dilution Factor.: 1.5)3121 5	Method:	TO-15		
Dilution Pactor 1.5	5	Methodistation	10-15	/	/
PARAMETER	RESULTS (ppb(v/v))	REPORTI LIMIT (pp		S (ug/m3) REPORT	
trans-1,3-Dichloroprope	ne NR		NR		
Acetone	NR		NR	1	
Ethylbenzene	NR		NR		
2-Hexanone	NR		NR	1	
Methylene chloride	NR		NR	1	
Benzene	NR		NR	1	
Styrene	NR		NR	·	
1,1,2,2-Tetrachloroethan			NR		
Tetrachloroethene	NR		NR		
Toluene	NR		NR		
1,1,1-Trichloroethane	NR		ŃR		
1,1,2-Trichloroethane	NR		/ NR		
Trichloroethene	NR				
Vinyl chloride	NR		NR		
o-Xylene m-Xylene & p-Xylene	NR NR		NR NR		
Bromodichloromethane	NR		NR NR		
2-Butanone (MEK)	NR	į	NR		
4-Methyl-2-pentanone (!		Ĵ	NR		
Bromoform	NR	/	NR		
Bromomethane	NR	1	NR		
Carbon disulfide	NR	1	NR		
Carbon tetrachloride	NR		NR		
Chlorobenzene	NR		NR		,
Dibromochloromethane	NR		NR		
Chloroethane	NR		NR		
Chloroform	NR		NR		
Chloromethane	NR		NR		
1,1-Dichloroethane	NR		NR		
1,2-Dichloroethane	NR		NR		
1,1-Dichloroethene	NR		NR		
cis-1,2-Dichloroethene	NR		NR		
trans-1,2-Dichloroethene			NR		
1,2-Dichloropropane	NR		NR		
cis-1,3-Dichloropropene	NR		NR		

TO-14 Conv Rev. 5

Client Sample ID: AB-1

GC/MS Volatiles

Lot-Sample # H5G200167 - 004	Work Order # HFVX11AD	Matrix: AIR
SURROGATE	PERCENT RECOVERY	LABORATORY CONTROL LIMITS (%)
1,2-Dichloroethane-d4 Toluene-d8 4-Bromofluorobenzene	0.0 * 0.0 * 0.0 *	70 - 130 70 - 130 70 - 130

Qualifiers

* Surrogate recovery is outside stated control limits. NR Not reportable.

The 'Result' in ug/m3 is calculated using the following equation: Amount Found(before rounding)*(Molecular Weight/24,45)

The 'Reporting Limit' in ug/m3 is calculated using the following equation: (Reporting Limit(before rounding) * Dilution Factor) * (Molecular Weight/24.45)