DEPARTMENT OF THE ARMY
CAMP STANLEY STORAGE ACTIVITY, RRAD
25800 RALPH FAIR ROAD, BOERNE, TX 78015-4800

August 28, 2001

Office of the Commander U - 029 - 01

Mr.

Mr.

Camp Stanley Storage Activity (CSSA) collected groundwater samples from your water well on 1 August and 15 August 2001. These samples were submitted to a laboratory contracted by CSSA's environmental contractor for volatile organic compound (VOC) analysis. This letter provides you with the VOC data from the laboratory results and a formal thank you for your assistance in this groundwater monitoring effort.

We provide below a summary of detected VOC compounds compared to maximum contaminant levels (MCLs) allowed in drinking water by the U.S. Environmental Protection Agency (EPA) under the Safe Drinking Water Act. We have attached the laboratory summary sheets from the initial August 1, 2001 sampling event. Results from the August 15, 2001 event are provided in the table also, however the data is still "preliminary" since validation and verification required for these samples have not yet been completed. Validation and verification normally take six weeks to complete. When we receive validation and verification we will provide it to you promptly.

LS-6 Well Adjacent to Church

Date Sampled	*VOC Compound*	*Result (ppb)*	*MCL (ppb)*
First Sample Event at Wellhead			
August 1, 2001	Tetrachloroethene (PCE)	6.0	5
	Trichloroethene (TCE)	0.31	5
Confirmation Sample Event at Wellhead			
August 15, 2001	PCE	6.5	5
	TCE	0.34	5
First Sample Event at Cafeteria Tap			
August 15, 2001	PCE	3.91	5
	TCE	0.24	5
First Sample Post-GAC			
August 28, 2001	PCE	<5.0	5
	TCE	<5.0	5

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LS-5 Up the Street from Church

Date Sampled	*VOC Compound*	*Result (ppb)*	*MCL (ppb)*
First Tap Sample Event at Vacant Residence on Curres Creek Rd.			
August 1, 2001	PCE	0.28	5
	TCE	0.51	5

Analytical results from the initial August 1 sampling event found low levels of the VOCs, PCE and TCE, in both LS-5 and LS-6. Our prime contractor contacted us on 9August to advise us that the samples had a PCE level in well LS-6 that was slightly above the MCL. Consequently on August 15, CSSA took a second set of confirmatory samples from your primary drinking water well (LS-6). One sample came from the wellhead and the second from the cafeteria tap. The purpose of these samples was to confirm the original results and to determine what VOC levels were at the point of use.

Analyses of the follow-up samples also identified PCE and TCE. Of special note is the lower concentration found in the cafeteria tap sample taken during the second sampling event of August 15, 2001. Tap sample concentrations were below the MCL for both TCE and PCE and indicate water at that tap meets federal and state drinking water standards. It is likely that the lower levels were a result of the VOCs volatilizing in the well cistern and during movement through the water distribution system.

On Tuesday, August 15, 2001, CSSA arranged to have a granular activated carbon (GAC) filtration system installed on your well. CSSA initiated the installation of a granular activated carbon treatment unit for your wellhead to reduce VOC contaminant levels to within drinking water standards. A copy of the invoice is attached for your records.

Carbonair Environmental Systems of San Marcos, Texas installed the filtration system on August 15. CSSA contracted with Carbonair to install the system at no cost to you. The system will remain in operation for the foreseeable future or until significant reductions in contamination levels are seen in your well. As we discussed at the time of installation, CSSA will be responsible for all costs associated with operation and maintenance of this system. However, we ask that you take responsibility for exchanging the five-micron pre- and post-filters that are located within the system housing. The spare filters are located inside the housing and should be changed at least monthly. Spent filters can be disposed of with your regular household trash. Please contact the installer at (800) 893-5937 should you have questions regarding the filter change-out.

Carbonair is scheduled to perform maintenance on the system every six months. Their first maintenance visit is scheduled for approximately February 8, 2002. At that time, Carbonair will exchange the first carbon canister and perform other routine maintenance operations. If you experience any problems with the system, please let the installer or CSSA know immediately. Carbonair is very responsive and can make additional maintenance visits if needed.

On August 28, 2001, CSSA collected a confirmation water sample from LS-6, above the GAC filtration system. This sample was submitted for rush VOC analyses using a quick 24-hour turnaround. Because of the rapid turnaround, it will not be possible to perform all of the validation and verification and other quality assurance procedures that our analytical results usually are subjected to. For this reason, these analytical results are considered "screening" level only. The resampling post-GAC had no detections of any contaminants above 5 ppb. On Thursday, August 30, CSSA will collect a second post GAC confirmation sample. This sample will also be submitted for VOC analyses using a seven-day turnaround time. After the results from this second sample have been validated and verified, we will forward the results to you. The results from this second post GAC sample will be considered more definitive.

CSSA will also take periodic samples from your well during the drilling processes of an additional monitoring well. The installation of the new well is located approximately 500 feet from yours and we want to ensure that our activities do not adversely impact your drinking water supply. We will contact you as these sampling events occur to keep you informed and provide analytical data as results become available.

As part of the ongoing CSSA environmental program, we are continuing to investigate and cleanup VOC source areas on the installation and to track these compounds in ground water on- and off-post. As part of this effort, we would like to contact you in the future to schedule another sampling event. When arrangements with contractors are complete, we will contact you with a proposed sampling date and time. Once we have arranged a date with you, CSSA will attempt to provide at least 72 hours notice prior to proposed sampling events. CSSA anticipates sampling your drinking water before and after treatment at least quarterly for the foreseeable future.

Again, we would like to thank you for your patience and cooperation. We regret that your well has been impacted, but remain committed to making sure your water is safe to use and

keeping you informed. If you have any questions concerning this letter, please contact me at 295-7416.

Sincerely,

Jason D. Shirley
Lieutenant Colonel, U.S. Army
Commanding Officer

CC: Mr. Greg Lyssy
EPA Region 6

Mr. Kirk Coulter
TNRCC Central Office

Mr. Tom Haberle
TNRCC Region 13

Ms. Abigail Power
TNRCC Region 13

Ms. Kyle Cunningham, R.S.
San Antonio Metropolitan Health District

Ms. Susan Roberts
Parsons ES

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ORGANIC ANALYSES DATA SHEET 2
RESULTS

Analytical Method: EPA 8260B Preparatory Method: 5030B AAB #: 010804AH-38194
Lab Name: APPL, Inc Contract #: F11623-94-D0024-RL83
Field Sample ID: LS-5 Lab Sample ID: AP20277 Matrix: Water
% Solids: NA Initial Calibration ID: H010804
Date Received: 02-Aug-01 Date Prepared: 04-Aug-01 Date Analyzed: 04-Aug-01
Concentration Units: ug/L

Analyte	MDL	RL	Concentration	Dilution	Confirm	Qualifier
1,1,1,2-Tetrachloroethane	0.14	0.5	0.14	1.0		U
1,1,1-TCA	0.08	0.8	0.08	1.0		U
1,1,2,2-Tetrachloroethane	0.20	0.4	0.20	1.0		U
1,1,2-TCA	0.16	1.0	0.16	1.0		U
1,1-DCA	0.07	0.4	0.07	1.0		U
1,1-DCE	0.16	1.2	0.16	1.0		U
1,1-Dichloropropene	0.12	1.0	0.12	1.0		U
1,2,3-Trichlorobenzene	0.13	0.5	0.13	1.0		U
1,2,3-Trichloropropane	0.23	3.2	0.23	1.0		U
1,2,4-Trichlorobenzene	0.08	0.5	0.08	1.0		U
1,2,4-Trimethylbenzene	0.07	1.3	0.07	1.0		U
1,2-DCA	0.10	0.6	0.10	1.0		U
1,2-DCB	0.08	0.3	0.08	1.0		U
1,2-Dibromo-3-chloropropane	0.72	2.6	0.72	1.0		U
1,2-Dichloropropane	0.14	0.4	0.14	1.0		U
1,2-EDB	0.11	0.6	0.11	1.0		U
1,3,5-Trimethylbenzene	0.06	0.5	0.06	1.0		U
1,3-DCB	0.12	1.2	0.12	1.0		U
1,3-Dichloropropane	0.10	0.4	0.10	1.0		U
1,4-DCB	0.09	0.3	0.09	1.0		U
1-Chlorohexane	0.12	0.6	0.12	1.0		U
2,2-Dichloropropane	0.53	3.5	0.53	1.0		U
2-Chlorotoluene	0.12	0.4	0.12	1.0		U
4-Chlorotoluene	0.09	0.6	0.09	1.0		U
Benzene	0.12	0.4	0.13	1.0		F
Bromobenzene	0.08	0.3	0.08	1.0		U
Bromochloromethane	0.16	0.4	0.16	1.0		U
Bromodichloromethane	0.12	0.8	0.12	1.0		U
Bromoform	0.14	1.2	0.14	1.0		U
Bromomethane	0.36	1.1	0.36	1.0		U
Carbon tetrachloride	0.09	2.1	0.09	1.0		U
Chlorobenzene	0.09	0.4	0.09	1.0		U
Chloroethane	0.26	1.0	0.26	1.0		U
Chloroform	0.06	0.3	0.06	1.0		U
Chloromethane	0.41	1.3	0.41	1.0		U
Cis-1,2-DCE	0.11	1.2	0.11	1.0		U
Cis-1,3-Dichloropropene	0.09	1.0	0.09	1.0		U

Comments:

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ORGANIC ANALYSES DATA SHEET 2
RESULTS

Analytical Method: EPA 8260B Preparatory Method: 5030B AAB #: 010804AH-38194
Lab Name: APPL, Inc Contract #: F11623-94-D0024-RL83
Field Sample ID: LS-5 Lab Sample ID: AP20277 Matrix: Water
% Solids: NA Initial Calibration ID: H010804
Date Received: 02-Aug-01 Date Prepared: 04-Aug-01 Date Analyzed: 04-Aug-01
Concentration Units: ug/L

Analyte	MDL	RL	Concentration	Dilution	Confirm	Qualifier
Dibromochloromethane	0.09	0.5	0.09	1.0		U
Dibromomethane	0.10	2.4	0.10	1.0		U
Dichlorodifluoromethane	0.24	1.0	0.24	1.0		U
Ethylbenzene	0.06	0.6	0.06	1.0		U
Hexachlorobutadiene	0.19	1.1	0.19	1.0		U
Isopropylbenzene	0.08	0.5	0.08	1.0		U
m&p-Xylene	0.14	0.5	0.14	1.0		U
Methylene chloride	0.19	1.0	0.19	1.0		U
n-Butylbenzene	0.11	1.1	0.11	1.0		U
n-Propylbenzene	0.10	0.4	0.10	1.0		U
Naphthalene	0.08	0.8	0.08	1.0		U
o-Xylene	0.07	1.1	0.07	1.0		U
p-Isopropyltoluene	0.06	1.2	0.06	1.0		U
Sec-Butylbenzene	0.05	1.3	0.05	1.0		U
Styrene	0.07	0.4	0.07	1.0		U
TCE	0.14	1.0	0.51	1.0		F
Tert-Butylbenzene	0.05	1.4	0.05	1.0		U
Tetrachloroethene	0.11	1.4	0.28	1.0		F
Toluene	0.11	1.1	0.11	1.0		U
Trans-1,2-DCE	0.14	0.6	0.14	1.0		U
Trans-1,3-Dichloropropene	0.14	1.0	0.14	1.0		U
Trichlorofluoromethane	0.09	0.8	0.09	1.0		U
Vinyl chloride	0.27	1.1	0.27	1.0		U

Surrogate	Recovery	Control Limits	Qualifier
1,2-DCA-D4(S)	101	62-139	
4-Bromofluorobenzene(S)	101	75-125	
Dibromofluoromethane(S)	97.8	75-125	
Toluene-D8(S)	94.9	75-125	

Internal Std	Qualifier
1,4-Dichlorobenzene-D(IS)	
Chlorobenzene-D5(IS)	
Fluorobenzene(IS)	

Comments:

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ORGANIC ANALYSES DATA SHEET 2
RESULTS

Analytical Method: EPA 8260B Preparatory Method: 5030B AAB #: 010804AH-38194
Lab Name: APPL, Inc Contract #: F11623-94-D0024-RL83
Field Sample ID: LS-6 Lab Sample ID: AP20278 Matrix: Water
% Solids: NA Initial Calibration ID: H010804
Date Received: 02-Aug-01 Date Prepared: 04-Aug-01 Date Analyzed: 04-Aug-01
Concentration Units: ug/L

Analyte	MDL	RL	Concentration	Dilution	Confirm	Qualifier
1,1,1,2-Tetrachloroethane	0.14	0.5	0.14	1.0		U
1,1,1-TCA	0.08	0.8	0.08	1.0		U
1,1,2,2-Tetrachloroethane	0.20	0.4	0.20	1.0		U
1,1,2-TCA	0.16	1.0	0.16	1.0		U
1,1-DCA	0.07	0.4	0.07	1.0		U
1,1-DCE	0.16	1.2	0.16	1.0		U
1,1-Dichloropropene	0.12	1.0	0.12	1.0		U
1,2,3-Trichlorobenzene	0.13	0.5	0.13	1.0		U
1,2,3-Trichloropropane	0.23	3.2	0.23	1.0		U
1,2,4-Trichlorobenzene	0.08	0.5	0.08	1.0		U
1,2,4-Trimethylbenzene	0.07	1.3	0.07	1.0		U
1,2-DCA	0.10	0.6	0.10	1.0		U
1,2-DCB	0.08	0.3	0.08	1.0		U
1,2-Dibromo-3-chloropropane	0.72	2.6	0.72	1.0		U
1,2-Dichloropropane	0.14	0.4	0.14	1.0		U
1,2-EDB	0.11	0.6	0.11	1.0		U
1,3,5-Trimethylbenzene	0.06	0.5	0.06	1.0		U
1,3-DCB	0.12	1.2	0.12	1.0		U
1,3-Dichloropropane	0.10	0.4	0.10	1.0		U
1,4-DCB	0.09	0.3	0.09	1.0		U
1-Chlorohexane	0.12	0.6	0.12	1.0		U
2,2-Dichloropropane	0.53	3.5	0.53	1.0		U
2-Chlorotoluene	0.12	0.4	0.12	1.0		U
4-Chlorotoluene	0.09	0.6	0.09	1.0		U
Benzene	0.12	0.4	0.14	1.0		F
Bromobenzene	0.08	0.3	0.08	1.0		U
Bromochloromethane	0.16	0.4	0.16	1.0		U
Bromodichloromethane	0.12	0.8	0.12	1.0		U
Bromoform	0.14	1.2	0.14	1.0		U
Bromomethane	0.36	1.1	0.36	1.0		U
Carbon tetrachloride	0.09	2.1	0.09	1.0		U
Chlorobenzene	0.09	0.4	0.09	1.0		U
Chloroethane	0.26	1.0	0.26	1.0		U
Chloroform	0.06	0.3	0.06	1.0		U
Chloromethane	0.41	1.3	0.41	1.0		U
Cis-1,2-DCE	0.11	1.2	0.11	1.0		U
Cis-1,3-Dichloropropene	0.09	1.0	0.09	1.0		U

Comments:

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ORGANIC ANALYSES DATA SHEET 2
RESULTS

Analytical Method: EPA 8260B Preparatory Method: 5030B AAB #: 010804AH-38194
Lab Name: APPL, Inc Contract #: F11623-94-D0024-RL83
Field Sample ID: LS-6 Lab Sample ID: AP20278 Matrix: Water
% Solids: NA Initial Calibration ID: H010804
Date Received: 02-Aug-01 Date Prepared: 04-Aug-01 Date Analyzed: 04-Aug-01
Concentration Units: ug/L

Analyte	MDL	RL	Concentration	Dilution	Confirm	Qualifier
Dibromochloromethane	0.09	0.5	0.09	1.0		U
Dibromomethane	0.10	2.4	0.10	1.0		U
Dichlorodifluoromethane	0.24	1.0	0.24	1.0		U
Ethylbenzene	0.06	0.6	0.06	1.0		U
Hexachlorobutadiene	0.19	1.1	0.19	1.0		U
Isopropylbenzene	0.08	0.5	0.08	1.0		U
m&p-Xylene	0.14	0.5	0.14	1.0		U
Methylene chloride	0.19	1.0	0.19	1.0		U
n-Butylbenzene	0.11	1.1	0.11	1.0		U
n-Propylbenzene	0.10	0.4	0.10	1.0		U
Naphthalene	0.08	0.8	0.08	1.0		U
o-Xylene	0.07	1.1	0.07	1.0		U
p-Isopropyltoluene	0.06	1.2	0.06	1.0		U
Sec-Butylbenzene	0.05	1.3	0.05	1.0		U
Styrene	0.07	0.4	0.07	1.0		U
TCE	0.14	1.0	0.31	1.0		F
Tert-Butylbenzene	0.05	1.4	0.05	1.0		U
Tetrachloroethene	0.11	1.4	6.00	1.0		
Toluene	0.11	1.1	0.11	1.0		U
Trans-1,2-DCE	0.14	0.6	0.14	1.0		U
Trans-1,3-Dichloropropene	0.14	1.0	0.14	1.0		U
Trichlorofluoromethane	0.09	0.8	0.09	1.0		U
Vinyl chloride	0.27	1.1	0.27	1.0		U

Surrogate	Recovery	Control Limits	Qualifier
1,2-DCA-D4(S)	101	62-139	
4-Bromofluorobenzene(S)	101	75-125	
Dibromofluoromethane(S)	98.1	75-125	
Toluene-D8(S)	95.2	75-125	

Internal Std	Qualifier
1,4-Dichlorobenzene-D(IS)	
Chlorobenzene-D5(IS)	
Fluorobenzene(IS)	

Comments:

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ORGANIC ANALYSES DATA SHEET 2
RESULTS

Analytical Method: EPA 8260B Preparatory Method: 5030B AAB #: 010816AH-39050
Lab Name: APPL, Inc Contract #: F11623-94-D0024-RL83
Field Sample ID: LS-6 WELL Lab Sample ID: AP21005 Matrix: Water
% Solids: NA Initial Calibration ID: H010813
Date Received: 16-Aug-01 Date Prepared: 16-Aug-01 Date Analyzed: 16-Aug-01
Concentration Units: ug/L

Analyte	MDL	RL	Concentration	Dilution	Confirm	Qualifier
1,1,1,2-Tetrachloroethane	0.14	0.5	0.14	1.0		U
1,1,1-TCA	0.08	0.8	0.08	1.0		U
1,1,2,2-Tetrachloroethane	0.20	0.4	0.20	1.0		U
1,1,2-TCA	0.16	1.0	0.16	1.0		U
1,1-DCA	0.07	0.4	0.07	1.0		U
1,1-DCE	0.16	1.2	0.16	1.0		U
1,1-Dichloropropene	0.12	1.0	0.12	1.0		U
1,2,3-Trichlorobenzene	0.13	0.5	0.13	1.0		U
1,2,3-Trichloropropane	0.23	3.2	0.23	1.0		U
1,2,4-Trichlorobenzene	0.08	0.5	0.08	1.0		U
1,2,4-Trimethylbenzene	0.07	1.3	0.07	1.0		U
1,2-DCA	0.10	0.6	0.10	1.0		U
1,2-DCB	0.08	0.3	0.08	1.0		U
1,2-Dibromo-3-chloropropane	0.72	2.6	0.72	1.0		U
1,2-Dichloropropane	0.14	0.4	0.14	1.0		U
1,2-EDB	0.11	0.6	0.11	1.0		U
1,3,5-Trimethylbenzene	0.06	0.5	0.06	1.0		U
1,3-DCB	0.12	1.2	0.12	1.0		U
1,3-Dichloropropane	0.10	0.4	0.10	1.0		U
1,4-DCB	0.09	0.3	0.09	1.0		U
1-Chlorohexane	0.12	0.6	0.12	1.0		U
2,2-Dichloropropane	0.53	3.5	0.53	1.0		U
2-Chlorotoluene	0.12	0.4	0.12	1.0		U
4-Chlorotoluene	0.09	0.6	0.09	1.0		U
Benzene	0.12	0.4	0.13	1.0		F
Bromobenzene	0.08	0.3	0.08	1.0		U
Bromochloromethane	0.16	0.4	0.16	1.0		U
Bromodichloromethane	0.12	0.8	0.12	1.0		U
Bromoform	0.14	1.2	0.14	1.0		U
Bromomethane	0.36	1.1	0.36	1.0		U
Carbon tetrachloride	0.09	2.1	0.09	1.0		U
Chlorobenzene	0.09	0.4	0.09	1.0		U
Chloroethane	0.26	1.0	0.26	1.0		U
Chloroform	0.06	0.3	0.06	1.0		U
Chloromethane	0.41	1.3	0.41	1.0		U
Cis-1,2-DCE	0.11	1.2	0.11	1.0		U
Cis-1,3-Dichloropropene	0.09	1.0	0.09	1.0		U

Comments:

AFCEE
ORGANIC ANALYSES DATA SHEET 2
RESULTS

Analytical Method: EPA 8260B Preparatory Method: 5030B AAB #: 010816AII-39050
Lab Name: APPL, Inc Contract #: F11623-94-D0024-RL83
Field Sample ID: LS-6 WELL Lab Sample ID: AP21005 Matrix: Water
% Solids: NA Initial Calibration ID: H010813
Date Received: 16-Aug-01 Date Prepared: 16-Aug-01 Date Analyzed: 16-Aug-01
Concentration Units: ug/L

Analyte	MDL	RL	Concentration	Dilution	Confirm	Qualifier
Dibromochloromethane	0.09	0.5	0.09	1.0		U
Dibromomethane	0.10	2.4	0.10	1.0		U
Dichlorodifluoromethane	0.24	1.0	0.24	1.0		U
Ethylbenzene	0.06	0.6	0.06	1.0		U
Hexachlorobutadiene	0.19	1.1	0.19	1.0		U
Isopropylbenzene	0.08	0.5	0.08	1.0		U
m&p-Xylene	0.14	0.5	0.14	1.0		U
Methylene chloride	0.19	1.0	0.19	1.0		U
n-Butylbenzene	0.11	1.1	0.11	1.0		U
n-Propylbenzene	0.10	0.4	0.10	1.0		U
Naphthalene	0.08	0.8	0.08	1.0		U
o-Xylene	0.07	1.1	0.07	1.0		U
p-Isopropyltoluene	0.06	1.2	0.06	1.0		U
Sec-Butylbenzene	0.05	1.3	0.05	1.0		U
Styrene	0.07	0.4	0.07	1.0		U
TCE	0.14	1.0	0.34	1.0		
Tert-Butylbenzene	0.05	1.4	0.05	1.0		U
Tetrachloroethene	0.11	1.4	6.50	1.0		
Toluene	0.11	1.1	0.11	1.0		U
Trans-1,2-DCE	0.14	0.6	0.14	1.0		U
Trans-1,3-Dichloropropene	0.14	1.0	0.14	1.0		U
Trichlorofluoromethane	0.09	0.8	0.09	1.0		U
Vinyl chloride	0.27	1.1	0.27	1.0		U

Surrogate	Recovery	Control Limits	Qualifier
1,2-DCA-D4(S)	110	62-139	
4-Bromofluorobenzene(S)	99.2	75-125	
Dibromofluoromethane(S)	101	75-125	
Toluene-D8(S)	94.1	75-125	

Internal Std	Qualifier
1,4-Dichlorobenzene-D(IS)	
Chlorobenzene-D5(IS)	
Fluorobenzene(IS)	

Comments:

SAN ANTONIO TESTING LABORATORY, INC.

REPORT OF CHEMICAL ANALYSIS

Report No. 2001-08-223
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Camp Stanley Storage Facility
25800 Ralph Fair Rd
Boerne, TX 78015-4800
ATTN: Brian Murphy

Date Received:	Date Reported:	Sample Type:
08-28-2001	08-29-2001	Liquid

PROJECT NAME: Camp Stanley Storage Activity
PROJECT NUMBER: Off – Post Sample

SAMPLE ID #1: LS – 6 Post Gac

SURROGATE	% RECOVERY	LIMITS
Dibromofluoromethane	87	(86-115)
Toluene-d8	99	(81-121)
4-Bromofluorobenzene	104	(81-117)

VOLATILE TARGET COMPOUNDS	RESULTS (mg/L)	MDL (mg/L)	METHOD #	DATE ANALYZED
Dichlorodifluoromethane	<0.005	0.005	8260B	08-28-01
Chloromethane	<0.005	0.005	8260B	08-28-01
Vinyl Chloride	<0.002	0.002	8260B	08-28-01
Bromomethane	<0.005	0.005	8260B	08-28-01
Chloroethane	<0.005	0.005	8260B	08-28-01
Trichlorofluoromethane	<0.005	0.005	8260B	08-28-01
Acetone	<0.005	0.005	8260B	08-28-01
1,1-Dichloroethene	<0.005	0.005	8260B	08-28-01
Iodomethane	<0.005	0.005	8260B	08-28-01
Methylene Chloride	<0.005	0.005	8260B	08-28-01
Allyl Chloride	<0.005	0.005	8260B	08-28-01
Carbon Disulfide	<0.005	0.005	8260B	08-28-01
trans-1,2-Dichloroethene	<0.005	0.005	8260B	08-28-01
Methyl-t-butyl ether	<0.005	0.005	8260B	08-28-01
1,1-Dichloroethane	<0.005	0.005	8260B	08-28-01
Vinyl Acetate	<0.005	0.005	8260B	08-28-01
2-Butanone	<0.005	0.005	8260B	08-28-01
cis-1,2-Dichloroethene	<0.005	0.005	8260B	08-28-01
Bromochloromethane	<0.005	0.005	8260B	08-28-01
Chloroform	<0.005	0.005	8260B	08-28-01
2,2-Dichloropropane	<0.005	0.005	8260B	08-28-01
1,2-Dichloroethane	<0.005	0.005	8260B	08-28-01
1,1,1-Trichloroethane	<0.005	0.005	8260B	08-28-01
1,1-Dichloropropene	<0.005	0.005	8260B	08-28-01
Carbon tetrachloride	<0.005	0.005	8260B	08-28-01
Benzene	<0.005	0.005	8260B	08-28-01
Dibromoethane	<0.005	0.005	8260B	08-28-01
1,2-Dichloropropane	<0.005	0.005	8260B	08-28-01
Trichloroethene	<0.005	0.005	8260B	08-28-01
Bromodichloromethane	<0.005	0.005	8260B	08-28-01
2-Chloroethyl vinyl ether	<0.005	0.005	8260B	08-28-01
cis-1,3-Dichloropropene	<0.005	0.005	8260B	08-28-01

SAN ANTONIO
TESTING LABORATORY, INC.

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VOLATILE TARGET COMPOUNDS	RESULTS (mg/L)	MDL (mg/L)	METHOD #	DATE ANALYZED
4-Methyl-2-pentanone	<0.005	0.005	8260B	08-28-01
trans-1,3-Dichloropropene	<0.005	0.005	8260B	08-28-01
1,1,2-Trichloroethane	<0.005	0.005	8260B	08-28-01
Toluene	<0.005	0.005	8260B	08-28-01
1,3-dichloropropane	<0.005	0.005	8260B	08-28-01
2-Hexanone	<0.005	0.005	8260B	08-28-01
Dibromochloromethane	<0.005	0.005	8260B	08-28-01
1,2-Dibromoethane	<0.005	0.005	8260B	08-28-01
Tetrachloroethene	<0.005	0.005	8260B	08-28-01
1,1,1,2-Tetrachloroethane	<0.005	0.005	8260B	08-28-01
Chlorobenzene	<0.005	0.005	8260B	08-28-01
Ethylbenzene	<0.005	0.005	8260B	08-28-01
m-Xylene / p-Xylene	<0.010	0.010	8260B	08-28-01
Bromoform	<0.005	0.005	8260B	08-28-01
cis-1,4-Dichloro-2-Butane	<0.005	0.005	8260B	08-28-01
Styrene	<0.005	0.005	8260B	08-28-01
1,1,2,2-Tetrachloroethane	<0.005	0.005	8260B	08-28-01
o-Xylene	<0.005	0.005	8260B	08-28-01
1,2,3-Trichloropropane	<0.005	0.005	8260B	08-28-01
trans-1,4-Dichloro-2-Butane	<0.005	0.005	8260B	08-28-01
Pentachloroethane	<0.005	0.005	8260B	08-28-01
Isopropylbenzene	<0.005	0.005	8260B	08-28-01
Bromobenzene	<0.005	0.005	8260B	08-28-01
n-Propylbenzene	<0.005	0.005	8260B	08-28-01
2-Chlorotoluene	<0.005	0.005	8260B	08-28-01
4-Chlorotoluene	<0.005	0.005	8260B	08-28-01
1,3,5-Trimethylbenzene	<0.005	0.005	8260B	08-28-01
tert-Butylbenzene	<0.005	0.005	8260B	08-28-01
1,2,4-Trimethylbenzene	<0.005	0.005	8260B	08-28-01
sec- Butylbenzene	<0.005	0.005	8260B	08-28-01
1,4-Dichlorobenzene	<0.005	0.005	8260B	08-28-01
1,3-Dichlorobenzene	<0.005	0.005	8260B	08-28-01
p-Isopropyltoluene	<0.005	0.005	8260B	08-28-01
1,2-Dichlorobenzene	<0.005	0.005	8260B	08-28-01
n-Butylbenzene	<0.005	0.005	8260B	08-28-01
1,2-dibromo-3-chloropropane	<0.005	0.005	8260B	08-28-01
1,2,4-Trichlorobenzene	<0.005	0.005	8260B	08-28-01
Naphthalene	<0.005	0.005	8260B	08-28-01
Hexachlorobutadiene	<0.005	0.005	8260B	08-28-01
1,2,3-Trichlorobenzene	<0.005	0.005	8260B	08-28-01