

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

March 27, 2006

U-052-06

Mr.

26044 Old Fredericksburg Road Boerne, TX 78015

Subject: Sampling of Water Well OFR-1, Located at 26044 Old

Fredericksburg Road

Dear

Camp Stanley Storage Activity (CSSA) collected groundwater samples from your well (OFR-1) on 12/22/05. 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.

An abbreviated summary of analytical results compared to maximum contaminant levels (MCLs) allowed in drinking water by the U.S. EPA under the Safe Drinking Water Act is provided below:

| Date Sampled | VOC Compound | Result (ppb) | MCL (ppb) |
|--|--------------------------------------|--------------------|--------------|
| Well OFR-1 | l, Located at 26044 Old Fredericksbu | ırg Road | |
| 12/22/05 | Tetrachloroethene (PCE) | 0.29F | 5 |
| | Trichloroethene (TCE) | <0.05 (non-detect) | 5 |
| and the same of th | cis-1,2-Dichloroethene (DCE) | <0.07 (non-detect) | 7.0. |

*The "F" qualifier indicates the value is above the laboratory method detection limit, but below the laboratory reporting limit for the compound.

Based on the analytical data, low levels of the VOC PCE were identified in water samples from your well. These levels are below the applicable MCLs and do not affect usability of your well. Results from the laboratory analysis are provided as an attachment for the above sampling event.

In addition, a data qualifier, M, was placed on the analyte naphthalene for your well. The laboratory is required to follow certain quality assurance procedures, including a set of matrix spike and matrix spike duplicate analyses for every twenty wells sampled. The matrix spike analysis had naphthalene recovered below the acceptance criteria in another well from the same data package. The matrix spike duplicate was within the required quality assurance criteria. However, in accordance with the CSSA QAPP, the results for naphthalene were flagged "M" for all samples in this data package. All data for the December 2005 off-post groundwater monitoring event is considered usable. The "M" flag applied for naphthalene does not affect the usability of your well.

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 groundwater on- and offpost. As part of this effort, we may contact you in the future to schedule another sampling event for the well listed above.

Again, we would like to thank you for your 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 Glare Sanchez, Environmental Program Manager, at 698-5208.

Sincerely,

Jason D. Shirley

Installation Manager

Attachments

cc: Mr. Greq Lyssy, EPA Region 6

Mr. Sonny Rayos, TCEQ Central Office

Mr. Henry Karnei, TCEQ Region 13

Ms. Kyle Cunningham, San Antonio Metropolitan Health Dist.

Ms. Julie Burdey, Parsons

Ms. Kimberly Vaughn, Parsons

AFCEE ORGANIC ANALYSES DATA SHEET 2 RESULTS

Analytical Method: EPA 8260B

Preparatory Method:

AAB #: 060104AM-95232

Lab Name: APPL, Inc

5030B Contract #: F41624-03-D-8613, TO 08

Field Sample ID: OFR-1

Lab Sample ID: AX32994

Matrix: Water

% Solids: NA

Initial Calibration ID: M051231

Date Received: 23-Dec-05

Date Prepared: 05-Jan-06

Date Analyzed: 05-Jan-06

Concentration Units: ug/L

| Analyte | MDL | RL | Concentration | Dilution | Confirm | Qualifier |
|-------------------------|------|-----|---------------|----------|---------|-----------|
| 1,1-DCE | 0.12 | 1.2 | 0.12 | 1 | | U |
| Bromodichloromethane | 0.06 | 0.8 | 0.06 | 1 | | . " |
| Bromoform | 0.13 | 1.2 | 0.13 | 1 | | υ |
| Chloroform | 0.06 | 0.3 | 0.06 | 1 | | U |
| Cis-1,2-DCE | 0.07 | 1.2 | 0.07 | 1 | | U |
| Dibromochloromethane | 0.06 | 0.5 | 0.06 | 1 | | U |
| Dichlorodifluoromethane | 0.11 | 1.0 | 0.11 | 1 | | U |
| Methylene chloride | 0.51 | 2.0 | 0.51 | 1 | | U |
| Naphthalene | 0.07 | 0.4 | 0.07 | 1 | | M |
| TCE | 0.05 | 1.0 | 0.05 | 1 | | U |
| Tetrachloroethene | 0.06 | 1.4 | 0.29 | 1 | | F |
| Toluene | 0.06 | 1.1 | 0.06 | . 1 | | U |
| Trans-1,2-DCE | 0.08 | 0.6 | 0.08 | 1 | | U |
| Vinyl chloride | 0.08 | 1.1 | 0.08 | 1 | | U |

| Surrogate | Recovery | Control Limits | Qualifier |
|-------------------------|----------|----------------|-----------|
| 1,2-DCA-D4(S) | 102 | 69-139 | |
| 4-Bromofluorobenzene(S) | 96.7 | 75-125 | |
| Dibromofluoromethane(S) | 98.5 | 75-125 | |
| Toluene-D8(S) | 102 | 75-125 | |

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

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