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POST-CLOSURE SITE MONITORING REPORT
LANDFILL AND RESOURCE RECOVERY SUPERFUND SITE
NORTH SMITHFIELD, RHODE ISLAND

Prepared on Behalf of:

LANDFILL AND RESOURCE RECOVERY SITE GROUP

Prepared by:

O & M, Inc.
450 Montbrook Lane
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DECEMBER 2000



SDMS DocID 572990

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—
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Superfund Records Center
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December 11, 2000

Ms. Anna Krasko
US EPA
Mail Code: HBO
One Congress Street
Boston, MA 02114

**RE: Transmittal of December 2000 Post-Closure Site Monitoring Report
L & RR Superfund Site, North Smithfield, Rhode Island**

Dear Ms. Krasko:

Attached is one copy of the Post-Closure Site Monitoring Report for December 2000. This report includes the results from the July 2000 groundwater sampling event. The disk containing the groundwater quality data in Excel-compatible format will be forwarded under separate cover.

Please do not hesitate to call me at (908)735-9315 should you have any questions.

Sincerely,
de maximis, inc.

Melissa Williams
for John P. McBurney
Project Coordinator

DFG/mw

cc: D. Moreira - Waste Management, Inc. (w/o disk)
D. Peters - Metcalf & Eddy (w/o disk)
M. DiStefano, RIDEM (w/o disk)
T. Helgason - *de maximis, inc.* (w/o disk)

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1.0 INTRODUCTION

This Post-Closure Site Monitoring Report documents the Post-Closure Operation and Maintenance activities performed at the Landfill and Resource Recovery (L&RR) Superfund site from February 2000 through November 2000. During this period, *de maximis, inc.* coordinated all post-closure site monitoring and maintenance activities. Activities performed during the referenced period were conducted in accordance with the *Post-Closure Operation and Maintenance Plan; de maximis, inc.*; September 1996 ("Plan").

Groundwater and surface water sampling was performed April 4 through 7, 2000. The groundwater samples collected during the April groundwater sampling event were rejected due to the fact the field indicator parameters had not stabilized before the groundwater samples were collected. The groundwater sampling event was repeated from July 13 through 15, 2000. Sampling parameters for groundwater and surface water are presented in **Table 1** of this report.

Ambient air samples were collected June 8 and 9, 2000 from four locations at the landfill perimeter for TO-14 analysis. On August 14, 2000, a four hour composite sample was collected of the flare inlet gas using Method 18 sampling protocols and TO-14 analysis. Data from all the sampling programs conducted during the referenced period are presented in this report.

This report has been prepared in accordance with the requirements of the Consent Decree and Remedial Design/Remedial Action Statement of Work. This report addresses those requirements as indicated below:

| POST-CLOSURE SITE MONITORING REPORT REQUIREMENTS, SPECIFIED IN THE CONSENT DECREE AND RD/RA STATEMENT OF WORK | SECTION OF THIS REPORT WHERE THE REQUIREMENT IS ADDRESSED |
|--|--|
| a. map of the Site showing sample locations | Figure 1 Sample Location Plan |
| b. tabular representation of laboratory results by each media including comparison with any standard levels, with exceedances of maximum contaminant levels (MCLs) and other performance standards highlighted | 5.0 Groundwater Monitoring 6.0 Surface Water Monitoring |

| POST-CLOSURE SITE MONITORING REPORT REQUIREMENTS, SPECIFIED IN THE CONSENT DECREE AND RD/RA STATEMENT OF WORK (Cont'd) | SECTION OF THIS REPORT WHERE THE REQUIREMENT IS ADDRESSED (Cont'd) |
|--|--|
| c. laboratory results on a computer disc in a spreadsheet file such as Lotus 1-2-3 or D-Base | 5.0 Groundwater Monitoring |
| d. data validation packages | 4.0 Data Validation |
| e. results of statistical analysis of data | 8.0 Statistical Analysis |
| f. interpretation of any trends | 5.0 Groundwater Monitoring |
| g. inspection reports | 2.0 Landfill Inspection and Maintenance |
| h. description of maintenance activities completed | 2.0 Landfill Inspection and Maintenance |
| i. explanation of problems encountered in the field and measures taken to mitigate the problems | 7.0 Problems Encountered |
| j. activities planned for the next reporting period | 9.0 Activities Planned for Next Reporting Period |

2.0 LANDFILL INSPECTION AND MAINTENANCE

The landfill cover, drainage structures, gas system and groundwater monitoring wells were inspected monthly, with the results provided, on a monthly basis, to EPA and RIDEM. Overall, the components are in good condition. The following maintenance measures were performed:

- A timer was installed on the condensate pump to the injection system on February 7, 2000.
- The vegetative areas of the landfill cover were treated with slow-release fertilizer and grub control on May 25, 2000.
- Four small burrow holes, near gas wells 6 and 8, were filled in on March 3, 2000.
- Three small areas of sloughing were repaired on May 25, 2000
- The flexhose on gas well 7, which became hyperextended, was replaced on June 1, 2000.
- The flexhose on well 13 was removed and reconnected to avoid crimping on June 1, 2000.

- A louvre control switch on the flare was replaced the week of August 28, 2000.
- Monitoring wells MW-102A; MW-104A; and CW-6B were redeveloped on October 19 and 20, 2000.
- The flexhoses on wells 12, 14, and 18 were adjusted to avoid crimping and hyperextension on October 20, 2000.
- The landfill cover was mowed on November 9, 2000.

Copies of the landfill inspection reports are in **Appendix A**.

The annual settlement monument elevation survey was performed on October 31, 2000. Incremental settlement from the last settlement survey event (November, 1999), as well as total settlement are presented in the following table:

Settlement Monitoring Data

| Settlement Monument | Elevation (ft.) Following Installation* | Elevation (ft.) November, 1999 | Elevation (ft.) October, 2000 | Incremental Settlement (ft.) From November, 1999 to October, 2000 | Total Settlement (ft.) From Date of Installation |
|---------------------|---|--------------------------------|-------------------------------|---|--|
| SM-1 | 311.96 | 309.55 | 309.58 | +0.03 | 2.38 |
| SM-2 | 298.42 | 297.67 | 297.78 | +0.11 | 0.64 |
| SM-3 | 329.29 | 326.32 | 326.56 | +0.24 | 2.73 |
| SM-4 | 306.05 | 304.82 | 305.30 | +0.48 | 0.75 |
| SM-5 | 287.25 | 287.26 | 287.81 | +0.55 | +0.56 |
| SM-6 | 321.86 | 320.70 | 321.00 | +0.30 | 0.86 |
| SM-7 | 300.76 | 299.97 | 299.62 | 0.35 | 1.14 |
| SM-8 | 365.33 | 361.09 | 360.70 | 0.39 | 4.63 |
| SM-9 | 354.67 | 350.67 | 349.94 | 0.73 | 4.73 |
| SM-10 | 357.29 | 355.59 | 355.35 | 0.24 | 1.94 |
| SM-11 | 321.83 | 321.39 | 321.35 | 0.04 | 0.48 |
| SM-12 | 321.07 | 320.56 | 319.53 | 1.03 | 1.54 |
| SM-13** | 361.03 | 359.67 | No pipe | Not available | Not available |
| SM-14** | 360.20 | 358.54 | No pipe | Not available | Not available |

*All settlement monuments were installed in December, 1994 except for SM-13 and SM-14, which were installed on August 15, 1997.

**Settlement monuments SM-13 and SM-14 were inadvertently knocked over during a maintenance event and will be replaced and resurveyed.

3.0 LANDFILL GAS WELL AND PROBE MONITORING

The landfill gas extraction wells and methane migration probes were monitored periodically from February through November 2000. During this period, the results of the monitoring were provided on a monthly basis to EPA and RIDEM. Overall, the system continues to operate under negative pressure while controlling methane migration. Methane migration data appear below:

| Date | Methane Content (%) | | |
|----------|---------------------|------------|------------|
| | Probe GP-1 | Probe GP-4 | Probe GP-8 |
| 2/15/00 | 0.0 | 0.0 | 0.0 |
| 3/16/00 | 0.0 | 0.0 | 0.0 |
| 4/18/00 | 0.0 | 0.0 | 0.0 |
| 5/25/00 | 0.0 | 0.0 | 0.0 |
| 6/8/00 | 0.0 | 0.0 | 0.0 |
| 7/24/00 | 0.1 | 0.3 | 0.0 |
| 8/14/00 | 1.4 | 5.6 | 0.0 |
| 9/21/00 | 0.8 | 0.9 | 0.0 |
| 10/19/00 | 23.3 | 17.4 | 0.0 |
| 11/16/00 | 5.3 | 1.4 | 0.0 |

There were occasions where the methane level in a particular probe temporarily exceeded the action level of 1.25%. This was primarily due to a temporary flare shutdown or a gas well becoming temporarily inoperative due to a hyper-extended or crimped flex hose. In each case, the methane levels returned to lower levels in a few weeks following flare re-start or adjustment/replacement of the well flexhose. The methane levels in the probes will continue to be monitored on a monthly basis.

Gas well monitoring data is in **Appendix B**.

The condensate injection system, installed in December 1998, continued to treat about 1000 gallons of condensate every six to seven days.

On February 7, 2000, a timer was installed on the condensate pump to the injection system. That timer, set for a 10-hour injection cycle, effectively replaces the low level float switch in the condensate tank. An inspection of the flare was also made at that time, and the system was functioning properly. On August 4, 1999, the semi-annual flare inspection was performed by a representative of John Zink Co., the manufacturer of the flare. The results of that inspection indicate that the flare is in good condition and continues to

function properly. On August 28, 2000, a louvre control switch was replaced on the flare. Flare inspection reports and maintenance documentation is in **Appendix B**.

4.0 DATA VALIDATION

Analytical data for groundwater and surface water presented in this report were validated by Trillium, Inc. of Knoxville, Tennessee using the most current revision of the EPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Organics and Inorganic Analysis* (December 1996) as appropriate for each analysis. For analyses which do not have specific recommended guidelines for data review, results were evaluated based on the supporting data submitted with the data packages and on the professional judgement of the validator. Data qualifier codes by sample matrix applied by the Data Validator are presented in **Table 2**.

Analytical data for the Quality Assurance / Quality Control (QA/QC) samples were submitted as part of the analytical data package provided to Trillium, Inc. for data validation. The complete data validation reports are included in **Appendix C** (original field samples), **Appendix D** (split sample) and **Appendix E** (surface water samples).

5.0 GROUNDWATER MONITORING

Groundwater samples were collected on July 13 through 15, 2000, in accordance with the Plan. Prior to sample collection, groundwater elevations were measured at the monitoring wells. The water level monitoring data are presented in **Table 3**. **Figure 1** presents a groundwater sampling location plan.

Samples were collected from the seven monitoring wells in the groundwater monitoring network (MW-201, MW-202, MW-102A, MW-103A, MW-104A, CW-5B and CW-7A). The samples, along with the required QA/QC samples, were submitted to Severn Trent Laboratories (STL), formerly IEA Laboratories, in Monroe, Connecticut, for analysis. The specific parameters for groundwater samples and analytical methods are presented in **Table 1**.

5.1 Sample Collection

The monitoring wells were purged and sampled using the *EPA Region I Low Stress (low flow) Purging and Sampling Procedure for the Collection of Groundwater Samples from Monitoring Wells*, Revision 2. Purging and sampling of the monitoring wells was performed using dedicated, nitrogen-driven bladder pumps.

Field measurements (eH, pH, temperature, dissolved oxygen, turbidity and specific conductance) were collected using a YSI Model 6920 Closed Cell Water Quality Monitoring System as purging progressed. The field data collected during purging are presented in **Table 4**. The final reading for each location in **Table 4** is the reading collected immediately before sample collection.

Groundwater samples were shipped each day via Federal Express next day delivery to STL of Monroe, Connecticut, for analysis. A split sample from a single location (MW-104A) was provided to CompuChem Environmental Corporation (CompuChem) of Durham, North Carolina for the same suite of parameters as the original sample.

5.2 Quality Control Samples

All groundwater QA/QC samples for analysis, matrix spike and matrix spike duplicate analysis were collected from monitoring well CW-5B. A duplicate sample was collected from monitoring location MW-102A. A split sample was collected during the sampling event from monitoring location MW-104A for analysis of the same suite of parameters as the original sample. The split sample was shipped via overnight express mail to CompuChem. Trip blanks accompanied sample containers during shipment and were submitted for volatile organic analysis.

5.3 Groundwater Analytical Results

Analytical data packages provided by STL were submitted to Trillium, Inc. of Knoxville, Tennessee for validation. Data validation reports of the STL analyses for the July 2000 sampling event for groundwater are included in **Appendix C** of this report. Data qualifier code definitions are presented in **Table 2** and in the Attachments to each data validation report.

A summary of the validated groundwater analytical data for the July 2000, and previous sampling events under the Plan are presented in **Table 5**.

5.4 Split Sample Comparability Results

In accordance with the Plan, a split sample from monitoring location MW-104A was analyzed by the primary laboratory, STL, and by a secondary laboratory, CompuChem. The data from the primary and secondary laboratories were evaluated for comparability using the procedure outlined in the Plan. **Table 6** provides a comparison of the analytical results from both laboratories.

The data for all the analytes except for five are considered comparable using the criteria considered in the Plan. The five analytes for which the data are not considered comparable are: m, p xylene, Styrene, total lead, ammonia, and COD. The following presents potential causes for the non-comparable results.

5.4.1 m, p -xylene

The analytical data from Severn Trent Laboratories (STL) and CompuChem Laboratories both detected the compound, however, the Relative percent Difference (RPD) was 91.7 %. STL's higher concentration was qualified as "estimated" due to unacceptably high blank spike recoveries that may have result in a biased high result, therefore resulting in a high RPD.

5.4.2 Styrene

The result reported for styrene by STL was greater than twice the reporting limit of the non-detected result reported by Compuchem. The reported result for styrene by STL is very near the reporting limit. The detected estimated concentration of 2.1 $\mu\text{g/L}$ (from STL) is greater than two-times (2x) the nondetected result's reporting limit of 0.5 $\mu\text{g/L}$.

5.4.3 Total Lead

The reported concentrations of lead for both laboratories were estimated to be near the detection limits. The 91.6 % RPD between the reported results was outside the comparability criteria.

5.4.4 Ammonia and COD

The reported concentrations of ammonia and COD were outside the comparability criteria. The criteria for comparability is 30 %. Ammonia was calculated at 35.4 % RPD and COD was calculated at 60.4% RPD and therefore the results are not considered comparable.

6.0 SURFACE WATER MONITORING

Surface water samples were collected on April 7, 2000 from six surface water monitoring locations. Samples were not collected at locations LCH-2 or LCH-3, as those locations were dry.

Surface water samples were submitted to STL for analysis. Data from the surface water samples are presented in **Table 7**. Data validation reports for the surface water samples are presented in **Appendix E**.

7.0 PROBLEMS ENCOUNTERED

Groundwater samples collected during the April 2000 sampling event were rejected. After reviewing the field data, it was noted that at four wells, several field indicator parameters collected during the purging process had not stabilized as required by Low Flow sampling protocols. Therefore, a second groundwater sampling event was conducted during July 2000.

Except for a problem with the dissolved oxygen display at well MW-104A, no problems were encountered during the July sampling event. During the sampling of well MW-104A, the dissolved oxygen display on the YSI Model 6920 exhibited erratic behavior, the cause of which could not be readily determined at the time of sampling. Because all other parameters had stabilized, the sample was collected, recognizing the erratic dissolved oxygen reading.

8.0 STATISTICAL ANALYSIS

The DUMPSTAT statistical analysis, incorporating all of the groundwater quality data from October 1996 through October 1999, was submitted in February 2000. The next statistical analysis report will be submitted during the first quarter of the year 2001. That report will include the analysis of groundwater quality data from October 1996 through November 2000.

9.0 ACTIVITIES PLANNED FOR NEXT REPORTING PERIOD

The next reporting period will include the following activities from November 2000 through March 2001:

| <u>Activity</u> | <u>Schedule</u> |
|------------------------------------|-----------------|
| Landfill inspection | Quarterly |
| Gas Treatment System operation | Continuous |
| Extraction System monitoring | Monthly |
| Methane migration monitoring | Monthly |
| Semi-annual Groundwater Monitoring | November 2000 |

Tables

TABLE 1
POST-CLOSURE MONITORING/SAMPLING

| SAMPLING MEDIUM | PARAMETER AND METHOD | |
|---|---|---|
| GROUNDWATER | VOCs Arsenic (Total) Arsenic (Dissolved) Lead (Total) Lead (Dissolved) Chloride Iron Ammonia COD BOD | 8260B 7060A 7060A 7421 7421 352(.1, .2, .3) 6010B 350.2 410 (.1, .2) 405.1 |
| FIELD PARAMETERS for Surfacewater & Groundwater | eH pH DO Specific Conductance | |
| SURFACE WATER | VOCs Arsenic (Total) Arsenic (Dissolved) Chloride | 8260B 7060A 7060A 325(.1, .2, .3) |
| AMBIENT AIR/FLARE INLET | VOCs | TO-14 |
| FLARE INLET | VOCs | Method 18/TO-14 |

TABLE 2
DATA QUALIFIER DEFINITIONS BY MATRIX

GROUNDWATER AND SURFACE WATER

Volatile Organics in Groundwater and Surface Waters
STL Job Nos. 7000-1470A

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected. The associated numerical value is the sample quantitation limit. The sample quantitation limit accounts for the sample-specific dilution factors and percent solids corrections or sample sizes that deviate from those required by the method.
 - J - The associated numerical value is an estimated quantity.
 - R - The data are unusable (analyte may or may not be present). Resampling and reanalysis is necessary for verification. The R replaces the numerical value or sample quantitation limit.
 - UJ - The material was analyzed for, but was not detected. The sample quantitation limit is an estimated quantity.
-

TOTAL AND DISSOLVED METALS IN GROUNDWATER AND SURFACE WATERS

STL Case No. 1470A, Lot No. 9A08G369; Lot No 9A04G680

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.
 - J - The associated value is an estimated quantity.
 - R - The data are unusable (Note: Analyte may or may not be present).
 - UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.
-

AMMONIA, CHLORIDE, BOD AND COD IN GROUNDWATER AND SURFACE WATERS

STL Laboratory Case No. 1470A;; Lot No 9A04G680

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.
- J - The associated numerical value is an estimated quantity.
- R - The data are unusable (Note: Analyte may or may not be present).
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.

TABLE 2 Continued
DATA QUALIFIER DEFINITIONS BY MATRIX

GROUNDWATER
Compuchem Laboratories
Laboratory Case No. Q1458, SDG R1458
Volatile Organics in Groundwater

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected. The associated numerical value is the sample quantitation limit. The sample quantitation limit accounts for the sample-specific dilution factors and percent solids corrections or sample sizes that deviate from those required by the method.
 - J - The associated numerical value is an estimated quantity.
 - R - The data are unusable (compound may or may not be present). Resampling and reanalysis is necessary for verification. The R replaces the numerical value or the sample quantitation limit.
 - UJ - The material was analyzed for, but was not detected. The sample quantitation limit is an estimated quantity.
-

TOTAL AND DISSOLVED METALS IN GROUNDWATER
Compuchem Laboratories
Laboratory Case No. Q1458, SDG R1458 and S1458
Inorganic Analysis in Groundwater

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.
 - J - The associated numerical value is an estimated quantity.
 - R - The data are unusable (Note: Analyte may or may not be present).
 - UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.
-

CHLORIDE, BOD, COD AND AMMONIA IN GROUNDWATER
Compuchem Laboratories
Wet Chemistry In Groundwater
Laboratory Case No. r1458; Test America Project No. 200001

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.
- J - The associated numerical value is an estimated quantity.
- R - The data are unusable (Note: Analyte may or may not be present).
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.

TABLE 3: WATER LEVEL MONITORING DATA, L&RR SUPERFUND SITE, JULY 11, 2000

| Monitoring Location | Top of Casing Elevation (ftamsl) | Ground Surface Elevation (ftamsl) | Pump Intake Depth (fbtoc) | Total Depth ** (fbtoc) | Screened Interval (fbtoc) | Water Level (fbtoc) | Water Elevation (ftamsl) |
|---------------------|----------------------------------|-----------------------------------|---------------------------|------------------------|---------------------------|---------------------|--------------------------|
| MW -101 | 329.07 | 326.40 | 78 | 82.60 | 74.2 - 79.5 | 74.54 | 254.53 |
| MW - 102A | 258.03 | 255.90 | 69 | 78.12 | 62.7 - 73.3 | 11.45 | 246.58 |
| MW - 102B * | 253.74 | 251.10 | 35 | 42.05 | 28.9 - 39.4 | 7.41 | 246.33 |
| MW - 103A | 268.48 | 265.48 | 54 | 63.11 | 39.2 - 55.1 | 16.23 | 252.25 |
| MW - 103B * | 268.57 | 265.50 | 20 | 30.06 | 12.0 - 21.8 | 16.08 | 252.49 |
| MW - 104A | 263.54 | 260.54 | 52 | 55.63 | 43.5 - 54.0 | 16.46 | 247.08 |
| MW - 104B * | 263.77 | 260.77 | 22 | 25.72 | 14.5 - 24.0 | 12.51 | 251.26 |
| CW - 5A | 304.31 | 302.00 | 134 | 136.11 | 125.0 - 135.0 | 54.05 | 250.26 |
| CW - 5B * | 303.92 | 302.00 | 98 | 103.53 | 92.0 - 102.0 | 54.43 | 249.49 |
| CW - 5C | 303.98 | 302.40 | 62 | 69.41 | 48.5 - 68.5 | 53.45 | 250.53 |
| CW - 6A | 264.06 | 262.30 | 89 | 93.53 | 82.0 - 92.0 | 18.16 | 245.90 |
| CW - 6B * | 261.74 | 261.50 | 59 | 62.96 | 51.0 - 61.0 | 15.81 | 245.93 |
| CW - 6C | 263.98 | 261.30 | 25 | 34.61 | 13.0 - 33.0 | 15.18 | 248.80 |
| CW - 7A * | 255.59 | 254.90 | 46 | 57.96 | 37.0 - 47.0 | 7.81 | 247.78 |
| CW - 7B | 255.50 | 253.00 | None | 53.00 | 43.0 - 53.0 | 8.24 | 247.26 |
| CW - 7C | 255.05 | 254.50 | 26 | 36.55 | 7.0 - 27.0 | 8.63 | 246.42 |
| MW-201 * | 320.25 | 318.04 | 79 | 89.00 | 69.0 - 89.0 | 66.32 | 253.93 |
| MW-202 * | 253.26 | 251.43 | 28.5 | 36.00 | 21.0 - 36.0 | 11.21 | 242.05 |

Notes:

ftamsl - feet above mean sea level

fbtoc - feet below top of casing

* - Denotes unconsolidated aquifer monitoring locations.

** - As presented in well boring logs.

TABLE 4
WELL PURGING
FIELD WATER QUALITY MEASUREMENTS
July 2000

| L&RR SUPERFUND SITE | | | | | | | | | | DEPTH TO TOP & BOTTOM OF SCREEN - TOP 37'/BOTTOM 47' |
|-----------------------------|------------------------------|---|---|---------------------|--|------|----------|------------|--------------------|--|
| July 12, 2000 | | | | | | | | | | |
| WELL NUMBER - CW-7A | | | | | | | | | | PUMP INTAKE AT (FT. BELOW MP) - 46' |
| FIELD PERSONNEL - TK/SS | | | | | | | | | | PURGING DEVICE - Geotech/Bladder |
| SAMPLING ORGANIZATION - STL | | | | | | | | | | |
| CLOCK TIME | WATER DEPTH BELOW MP (FT) | PURGE RATE SPECIFY UNITS (ML/MIN) | CUM. VOLUME PURGED (SPECIFY UNITS) (ML) | TEMP (DEGREES C) | SPEC. COND. millimhos/cm AT 25 DEGREES C | pH | Eh mv | DO mg/l | TURBIDITY (NTU) | COMMENTS |
| 1325 | 7.81 | 400 | - | 16.22 | 311 | 6.45 | 207.5 | 2.55 | 34.73 | |
| 1330 | 7.81 | 400 | 2,000 | 16.11 | 349 | 6.3 | 111.1 | 2.1 | 31.78 | |
| 1335 | 7.81 | 400 | 4,000 | 15.58 | 333 | 6.13 | 21.6 | 1.1 | 12.42 | |
| 1340 | 7.81 | 400 | 6,000 | 15.99 | 332 | 6.06 | -25.1 | 0.7 | 8.52 | |
| 1345 | 7.81 | 400 | 8,000 | 15.77 | 334 | 6.08 | -47.1 | 0.8 | 7.55 | |
| 1350 | 7.81 | 400 | 10,000 | 15.30 | 334 | 6.06 | -68.1 | 0.4 | 7.96 | |
| 1355 | 7.81 | 400 | 12,000 | 15.10 | 338 | 6.07 | -76.0 | 0.4 | 7.64 | |
| 1400 | 7.81 | 400 | 14,000 | 15.24 | 342 | 6.07 | -78.1 | 0.4 | 1.85 | |
| 1405 | 7.81 | 400 | 16,000 | 15.33 | 343 | 6.07 | -79.3 | 0.4 | 2.27 | |
| 1410 | 7.81 | 400 | 18,000 | 15.46 | 342 | 6.09 | -84.7 | 0.4 | 2.12 | |
| 1415 | 7.81 | 400 | 20,000 | 15.52 | 342 | 6.09 | -88.6 | 0.4 | 2.07 | |

TABLE 4 (Continued)
WELL PURGING
FIELD WATER QUALITY MEASUREMENTS
July 2000

Page 1 of 1

| L&RR SUPERFUND SITE | | | | | | | | | | |
|--|------------------------------|---|---|---------------------|--|------|----------|------------|--------------------|----------|
| DEPTH TO TOP & BOTTOM OF SCREEN - TOP 62.9'/BOTTOM 73.3' | | | | | | | | | | |
| July 13, 2000 | | | | | | | | | | |
| PUMP INTAKE AT (FT. BELOW MP) - 69' | | | | | | | | | | |
| WELL NUMBER - MW-102A | | | | | | | | | | |
| PURGING DEVICE - Geotech/Bladder | | | | | | | | | | |
| FIELD PERSONNEL - TK/SS | | | | | | | | | | |
| SAMPLING ORGANIZATION - STL | | | | | | | | | | |
| CLOCK TIME | WATER DEPTH BELOW MP (FT) | PURGE RATE SPECIFY UNITS (ML/MIN) | CUM. VOLUME PURGED (SPECIFY UNITS) (ML) | TEMP (DEGREES C) | SPEC. COND. millimhos/cm AT 25 DEGREES C | pH | Eh mv | DO mg/l | TURBIDITY (NTU) | COMMENTS |
| 1330 | 11.45 | 400 | - | 14.76 | 476 | 5.92 | 13.4 | 3.45 | 11.57 | |
| 1335 | 11.75 | 400 | 2,000 | 14.13 | 441 | 5.83 | 29.3 | 2.76 | 9.44 | |
| 1340 | 11.63 | 400 | 4,000 | 14.28 | 434 | 5.85 | 28.2 | 3.34 | 10.9 | |
| 1345 | 11.61 | 400 | 6,000 | 13.78 | 434 | 5.73 | 38.0 | 1.39 | 10.64 | |
| 1350 | 11.61 | 400 | 8,000 | 13.72 | 433 | 5.86 | 28.1 | 1.08 | 12.78 | |
| 1355 | 11.61 | 400 | 10,000 | 13.28 | 433 | 5.84 | 26.4 | 0.69 | 10.12 | |
| 1400 | 11.61 | 400 | 12,000 | 13.55 | 433 | 5.76 | 35.9 | 0.92 | 20.33 | |
| 1405 | 11.61 | 400 | 14,000 | 13.66 | 435 | 5.84 | 26.6 | 0.97 | 13.84 | |
| 1410 | 11.61 | 400 | 16,000 | 13.80 | 434 | 5.87 | 27.3 | 0.73 | 11.35 | |
| 1415 | 11.61 | 400 | 18,000 | 13.78 | 436 | 5.85 | 20.7 | 0.36 | 17.63 | |
| 1420 | 11.61 | 400 | 20,000 | 13.55 | 438 | 5.87 | 22.0 | 0.45 | 17.33 | |
| 1425 | 11.61 | 400 | 22,000 | 13.72 | 439 | 5.93 | 18.7 | 0.82 | 17.37 | |

TABLE 4 (Continued)
WELL PURGING
FIELD WATER QUALITY MEASUREMENTS
July 2000

| | |
|--|--|
| L&RR SUPERFUND SITE July 12, 2000 WELL NUMBER - MW-201 FIELD PERSONNEL - TK/SS SAMPLING ORGANIZATION - STL | DEPTH TO TOP & BOTTOM OF SCREEN - TOP 69'/BOTTOM 89' PUMP INTAKE AT (FT. BELOW MP) - 79' PURGING DEVICE - Geotech/Bladder |
|--|--|

| CLOCK TIME | WATER DEPTH BELOW MP (FT) | PURGE RATE SPECIFY UNITS (ML/MIN) | CUM. VOLUME PURGED (SPECIFY UNITS) (ML) | TEMP (DEGREES C) | SPEC. COND. millimhos/cm At 25 Degrees C | pH | Eh mv | DO mg/l | TURBIDITY (NTU) | COMMENTS |
|------------|------------------------------|---|---|---------------------|--|------|----------|------------|--------------------|----------|
| 845 | 66.32 | 400 | - | 10.86 | 61 | 7.96 | 185.5 | 9.93 | 115.10 | |
| 850 | 66.40 | 400 | 2,000 | 11.02 | 62 | 7.55 | 187.9 | 8.04 | 146.00 | |
| 855 | 66.93 | 400 | 4,000 | 10.70 | 61 | 6.98 | 201.9 | 7.77 | 144.00 | |
| 900 | 66.35 | 400 | 6,000 | 10.89 | 62 | 6.72 | 208.6 | 8.53 | 103.00 | |
| 905 | 66.35 | 400 | 8,000 | 10.92 | 62 | 6.62 | 214.5 | 7.74 | 78.00 | |
| 910 | 66.35 | 400 | 10,000 | 10.91 | 60 | 6.54 | 219.7 | 7.72 | 47.02 | |
| 915 | 66.35 | 400 | 12,000 | 11.02 | 60 | 6.49 | 223.1 | 7.79 | 33.15 | |
| 920 | 66.35 | 400 | 14,000 | 11.04 | 60 | 6.43 | 226.1 | 7.83 | 31.61 | |
| 925 | 66.35 | 400 | 16,000 | 11.16 | 60 | 6.37 | 233.1 | 7.80 | 22.47 | |
| 930 | 66.35 | 400 | 18,000 | 11.26 | 60 | 6.32 | 233.6 | 7.49 | 21.70 | |
| 935 | 66.35 | 400 | 20,000 | 11.48 | 59 | 6.30 | 234.8 | 7.80 | 17.34 | |
| 940 | 66.35 | 400 | 22,000 | 11.59 | 59 | 6.28 | 235.9 | 7.70 | 16.97 | |
| 945 | 66.35 | 400 | 24,000 | 11.96 | 59 | 6.26 | 236.1 | 7.71 | 15.72 | |
| 950 | 66.35 | 400 | 26,000 | 12.04 | 59 | 6.23 | 237.7 | 7.35 | 14.88 | |
| 955 | 66.35 | 400 | 28,000 | 12.21 | 59 | 6.23 | 237.7 | 7.35 | 14.88 | |

TABLE 4 (Continued)
WELL PURGING
FIELD WATER QUALITY MEASUREMENTS
 July 2000

L&RR SUPERFUND SITE

DEPTH TO TOP & BOTTOM OF SCREEN - TOP 39.2'/BOTTOM 55.1

July 12, 2000

WELL NUMBER - MW-103A

PUMP INTAKE AT (FT. BELOW MP) - 54'

FIELD PERSONNEL - TK/SS

PURGING DEVICE - Geotech/Bladder

SAMPLING ORGANIZATION - STL

| CLOCK TIME | WATER DEPTH BELOW MP (FT) | PURGE RATE SPECIFY UNITS (ML/MIN) | CUM. VOLUME PURGED (SPECIFY UNITS) (ML) | TEMP (DEGREES C) | SPEC. COND. millimhos/cm AT 25 DEGREES C | pH | Eh mv | DO mg/l | TURBIDITY (NTU) | COMMENTS |
|------------|------------------------------|---|---|---------------------|--|------|----------|------------|--------------------|----------|
| 1215 | 16.23 | 300 | - | 11.38 | 174 | 6.21 | 280.5 | 6.75 | 0.0 | |
| 1220 | 20.05 | 300 | 1,500 | 12.27 | 174 | 6.46 | 266.1 | 5.81 | 0.0 | |
| 1225 | 20.05 | 300 | 3,000 | 12.67 | 175 | 6.64 | 265.1 | 5.10 | 0.0 | |
| 1230 | 20.05 | 300 | 4,500 | 12.59 | 175 | 6.77 | 250.0 | 5.30 | 0.0 | |
| 1235 | 20.05 | 300 | 6,000 | 12.55 | 175 | 6.81 | 248.4 | 4.60 | 0.0 | |
| 1240 | 20.05 | 300 | 7,500 | 12.80 | 175 | 6.89 | 243.8 | 4.62 | 0.0 | |
| 1245 | 20.05 | 300 | 9,000 | 13.01 | 175 | 6.95 | 239.9 | 4.41 | 0.0 | |
| 1250 | 20.05 | 300 | 10,500 | 13.12 | 176 | 6.98 | 239.3 | 4.42 | 0.0 | |

TABLE 4 (Continued)
WELL PURGING
FIELD WATER QUALITY MEASUREMENTS
July 2000

| L&RR SUPERFUND SITE | | | | | | | | | | |
|--|------------------------------|---|---|---------------------|--|------|----------|------------|--------------------|----------|
| July 13, 2000 | | | | | | | | | | |
| WELL NUMBER - CW-5B | | | | | | | | | | |
| FIELD PERSONNEL - TK/SS | | | | | | | | | | |
| SAMPLING ORGANIZATION - STL | | | | | | | | | | |
| DEPTH TO TOP & BOTTOM OF SCREEN - TOP92'/BOTTOM 102' | | | | | | | | | | |
| PUMP INTAKE AT (FT. BELOW MP) - 98' | | | | | | | | | | |
| PURGING DEVICE - Geotech/Bladder | | | | | | | | | | |
| CLOCK TIME | WATER DEPTH BELOW MP (FT) | PURGE RATE SPECIFY UNITS (ML/MIN) | CUM. VOLUME PURGED (SPECIFY UNITS) (ML) | TEMP (DEGREES C) | SPEC. COND. millimhos/cm AT 25 DEGREES C | pH | Eh mv | DO mg/l | TURBIDITY (NTU) | COMMENTS |
| 920 | 54.43 | 400 | -- | 13.94 | 180 | 5.78 | 42.0 | 6.95 | 0.01 | |
| 925 | 54.51 | 400 | 2,000 | 11.36 | 170 | 5.60 | 18.8 | 4.32 | 3.27 | |
| 930 | 54.51 | 400 | 4,000 | 11.58 | 169 | 5.66 | 18.7 | 3.35 | 0.00 | |
| 935 | 54.51 | 400 | 6,000 | 11.66 | 168 | 5.63 | 1.0 | 1.86 | 0.00 | |
| 940 | 54.51 | 400 | 8,000 | 11.60 | 165 | 5.69 | 22.5 | 2.24 | 0.00 | |
| 945 | 54.51 | 400 | 10,000 | 11.52 | 162 | 5.66 | 21.0 | 1.1 | 0.29 | |
| 950 | 54.51 | 400 | 12,000 | 11.44 | 161 | 5.69 | 28.0 | 1.34 | 0.00 | |
| 955 | 54.51 | 400 | 14,000 | 11.53 | 161 | 5.65 | 29.3 | 1.33 | 0.08 | |
| 1000 | 54.51 | 400 | 16,000 | 11.61 | 159 | 5.65 | 30.2 | 1.35 | 0.00 | |
| 1005 | 54.51 | 400 | 18,000 | 11.62 | 159 | 5.62 | 28.6 | 1.39 | 0.00 | |
| 1010 | 54.51 | 400 | 20,000 | 11.63 | 159 | 5.69 | 31.0 | 1.41 | 0.00 | |

TABLE 4 (Continued)
WELL PURGING
FIELD WATER QUALITY MEASUREMENTS
July 2000

| L&RR SUPERFUND SITE | | | | | | | | | | DEPTH TO TOP & BOTTOM OF SCREEN - TOP 44'/BOTTOM 54' |
|-----------------------------|------------------------------|---|---|---------------------|--|------|----------|------------|--------------------|--|
| July 13, 2000 | | | | | | | | | | |
| WELL NUMBER - MW-104A | | | | | | | | | | PUMP INTAKE AT (FT. BELOW MP) - 49' |
| FIELD PERSONNEL - TK/SS | | | | | | | | | | PURGING DEVICE - Geotech/Bladder |
| SAMPLING ORGANIZATION - STL | | | | | | | | | | |
| CLOCK TIME | WATER DEPTH BELOW MP (FT) | PURGE RATE SPECIFY UNITS (ML/MIN) | CUM. VOLUME PURGED (SPECIFY UNITS) (ML) | TEMP (DEGREES C) | SPEC. COND. millimhos/cm AT 25 DEGREES C | pH | Eh mv | DO mg/l | TURBIDITY (NTU) | COMMENTS |
| 1115 | 16.56 | 400 | -- | 17.52 | 2303 | 6.42 | -86.7 | 0.67 | 343 | |
| 1120 | 16.72 | 400 | 2,000 | 15.91 | 2256 | 6.32 | -78.1 | 0.44 | 754 | |
| 1125 | 16.72 | 400 | 4,000 | 15.88 | 2197 | 6.32 | -80.3 | 0.52 | 1000 | Cleaned Cell |
| 1130 | 16.72 | 400 | 6,000 | 15.70 | 2227 | 6.32 | -81.4 | 0.41 | -- | |
| 1135 | 16.72 | 400 | 8,000 | 15.62 | 2214 | 6.32 | -88.2 | 0.38 | -- | |
| 1140 | 16.72 | 400 | 10,000 | 15.56 | 2204 | 6.32 | -92.1 | 0.28 | 3260 | |
| 1145 | 16.72 | 400 | 12,000 | 15.52 | 2184 | 6.32 | -94.8 | 0.28 | 1370 | |
| 1150 | 16.72 | 400 | 14,000 | 15.31 | 2161 | 6.33 | -98.5 | 0.21 | 1665 | Cleaned Cell |
| 1155 | 16.72 | 400 | 16,000 | 15.07 | 2281 | 6.34 | -89.2 | 0.42 | 2360 | |
| 1200 | 16.72 | 400 | 18,000 | 15.83 | 2280 | 6.33 | -88.6 | 0.33 | 920 | |
| 1205 | 16.72 | 400 | 20,000 | 15.95 | 2267 | 6.33 | -93.1 | 0.20 | 835 | Cleaned Cell |
| 1210 | 16.72 | 400 | 22,000 | 15.82 | 2284 | 6.34 | -91.2 | 0.13 | 1425 | |
| 1215 | 16.72 | 400 | 24,000 | 15.97 | 2285 | 6.33 | -90.4 | 0.62 | 1515 | Cleaned Cell |
| 1220 | 16.72 | 400 | 26,000 | 15.52 | 2300 | 6.37 | -91.4 | 0.54 | 2015 | |
| 1225 | 16.72 | 400 | 28,000 | 15.68 | 2260 | 6.35 | -94.3 | 0.34 | 550 | |
| 1230 | 16.72 | 400 | 30,000 | 15.48 | 2226 | 6.34 | -100.2 | 0.31 | 995 | |
| 1235 | 16.72 | 400 | 32,000 | 15.32 | 2210 | 6.32 | -104.1 | 0.19 | 1020 | |
| 1240 | 16.72 | 400 | 34,000 | 15.59 | 2167 | 6.32 | -108.1 | 0.24 | 925 | |
| 1245 | 16.72 | 400 | 36,000 | 15.52 | 2150 | 6.33 | -109.8 | 0.26 | 1005 | |

TABLE 4 (Continued)
WELL PURGING
FIELD WATER QUALITY MEASUREMENTS
July 2000

| | |
|--|---|
| L&RR SUPERFUND SITE July 12, 2000 WELL NUMBER - MW-202 FIELD PERSONNEL - TK/SS SAMPLING ORGANIZATION - STL | DEPTH TO TOP & BOTTOM OF SCREEN - TOP 21.6'/BOTTOM 38.6' PUMP INTAKE AT (FT. BELOW MP) - 28.6 PURGING DEVICE - Geotech/Bladder |
|--|---|

| CLOCK TIME | WATER DEPTH BELOW MP (FT) | PURGE RATE SPECIFY UNITS (ML/MIN) | CUM. VOLUME PURGED (SPECIFY UNITS) (ML) | TEMP (DEGREES C) | SPEC. COND. millimhos/cm AT 25 DEGREES C | pH | Eh mv | DO mg/l | TURBIDITY (NTU) | COMMENTS |
|------------|------------------------------|---|---|---------------------|--|------|----------|------------|--------------------|----------|
| 1030 | 11.21 | 400 | - | 11.23 | 102 | 5.41 | 287.8 | 9.51 | 209 | |
| 1035 | 11.40 | 400 | 2,000 | 10.50 | 101 | 5.21 | 303.7 | 8.90 | 113 | |
| 1040 | 11.32 | 400 | 4,000 | 11.49 | 101 | 5.18 | 310.1 | 8.61 | 71 | |
| 1045 | 11.32 | 400 | 6,000 | 11.55 | 103 | 5.16 | 315.2 | 8.62 | 104 | |
| 1050 | 11.32 | 400 | 8,000 | 12.34 | 102 | 5.14 | 320.7 | 9.10 | 47 | |
| 1055 | 11.32 | 400 | 10,000 | 12.27 | 103 | 5.12 | 324.7 | 8.24 | 100 | |
| 1100 | 11.32 | 400 | 12,000 | 12.31 | 103 | 5.12 | 328.2 | 8.40 | 120 | |
| 1105 | 11.32 | 400 | 14,000 | 12.39 | 103 | 5.10 | 330.5 | 8.42 | 105 | |
| 1110 | 11.32 | 400 | 16,000 | 12.83 | 102 | 5.10 | 332.7 | 8.22 | 111 | |
| 1115 | 11.32 | 400 | 18,000 | 12.87 | 102 | 5.10 | 334.2 | 8.13 | 113 | |
| 1120 | 11.32 | 400 | 20,000 | 13.04 | 102 | 5.09 | 331.0 | 7.98 | 107 | |
| 1125 | 11.32 | 400 | 22,000 | 11.26 | 102 | 5.11 | 336.4 | 8.77 | 84 | |
| 1130 | 11.32 | 400 | 24,000 | 11.63 | 102 | 5.06 | 340.2 | 8.29 | 92 | |
| 1135 | 11.32 | 400 | 26,000 | 11.92 | 102 | 5.08 | 341.3 | 8.38 | 66 | |
| 1140 | 11.32 | 400 | 28,000 | 12.05 | 101 | 5.07 | 342.1 | 8.26 | 67 | |
| 1145 | 11.32 | 400 | 30,000 | 12.15 | 101 | 5.09 | 341.9 | 8.42 | 71 | |

TABLE 5: GROUNDWATER CHEMISTRY DATA, L&RR SUPERFUND SITE, JULY 2000

| MONITORING WELL | MCL (2) | MW-102A | | | | | | | | | | Duplicate July 2000 |
|------------------------------------|---------|--------------|---------------|-----------|--------------|---------------|-----------|--------------|------------|-----------|---------|---------------------|
| | | October 1996 | February 1997 | June 1997 | October 1997 | February 1998 | June 1998 | October 1998 | March 1999 | July 2000 | | |
| TCL VOLATILES (ug/L) | | | | | | | | | | | | |
| Acetone | | 5 U | 5 U J | 5 U | R | R | 2 UJ | 2 U | R | R | 100 U J | 69 U J |
| Benzene | 5 | 12 | 10 | 9 | 8 | 7 | 6 J | 7 | 9 | 7.6 | 10 | 10 |
| Bromobenzene | | 5 U | 5 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 0.5 U | 5 U |
| Chloroethane | | 100 | 58 | 56 | 53 | 47 | 32 | 34 | 25 J | 1 U | 28 | 27 |
| Chlorobenzene | 100 | 24 | 21 | 21 | 16 | 16 | 14 J | 12 | 15 | 11 | 11 | 11 |
| Chloromethane | | 5 U | 2 | 1 U | 1 U | 2 U | 1 UJ | 1 U | 1 UJ | 1 U | 5 U | 5 U |
| 2-Chlorotoluene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 5 U |
| Dichlorodifluoromethane | | 35 J | 22 | 32 | 2 U | 26 | 28 | 20 | 2 UJ | 46 J | 27 | 27 |
| Trichlorofluoromethane | | 5 U | 1 U | 1 U | 1 U | 1 J | 1 U | 1 U | 1 U | 2 | 5 U | 1.7 U |
| 1,1 - Dichloroethane | | 180 | 130 | 110 J | 110 | 120 J | 70 | 100 | 130 | 109 | 160 | 160 |
| 1,1 - Dichloroethene | 7 | 1 U | 1 U | 1 U | 1 U | 2 U | 1 U | 0.6 J | 1 U | 1 U | 5 U | 5 U |
| 1,2 - Dichloroethane | 5 | 18 | 16 | 13 | 13 | 12 | 11 | 10 | 12 | 1 U | 15 | 16 |
| 1,2 - Dichloropropane | 5 | 4 | 3 | 4 | 3 | 2 U | 3 | 3 | 4 | 4.1 | 6.4 | 6.6 |
| cis-1,2-Dichloroethene | 70 | 46 | 43 | 1 U | 100 | 110 J | 1 U | 130 | 180 | 170 | 230 | 240 |
| trans-1,2-Dichloroethene | 100 | 5 U | 2 | 2 | 1 U | 2 | 1 U | 2 | 3 | 2.4 | 3.6 J | 3.8 J |
| 1,2-dichlorobenzene | 600 | 5 U | 2 | 2 | 1 | 1 J | 1 U | 1 | 1 U | 1 U | 0.93 J | 1 J |
| 1,3-Dichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 5 U | 5 U |
| 1,4-Dichlorobenzene | 75 | 8 | 6 | 6 | 5 | 5 | 4 | 4 | 1 U | 4.2 | 5.5 | 5.7 |
| Ethylbenzene | 700 | 5 U | 1 | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 5 U |
| Isopropylbenzene | | 5 | 5 | 1 U | 4 | 4 | 3 | 3 | 1 U | 3.5 | 4.1 J | 4.2 J |
| p-Isopropyltoluene | | 5 U | 0.8 J | 0.7 J | 1 U | 2 U | 1 U | 1 U | 1 U | 5 U | 5 U | 5 U |
| Methylene Chloride | 5 | 5 U | 2 U J | 2 J | 4 J | 2 U | 2 U | 2 U J | 2 U | 1.5 J | 32 U J | 31 U J |
| Naphthalene | | 5 U | 2 J | 6 | 1 U | 1 J | 2 U J | 0.8 J | 2 U J | 2 U J | 5 U | 5 U |
| n-Propylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 5 U |
| Tetrachloroethene | 5 | 11 | 15 | 25 | 35 | 34 J | 34 | 46 | 70 | 53 | 55 J | 54 J |
| Trichloroethene | 5 | 15 | 18 | 25 | 31 | 31 | 33 | 38 | 52 | 50 | 77 | 79 |
| Toluene | 1000 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.95 J | 0.89 J |
| o-Xylene | (4) | 5 U | 1 J | 1 U | 1 U | 2 U | 1 U | 0.6 J | 1 U | 1 U | 5 U | 5 U |
| m,p Xylenes | (4) | 5 U | 3 | 2 | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 10 U | 10 U |
| 1, 2, 3 - Trichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U J | 1 U J | 5 U | 5 U |
| 1, 1, 1 - Trichloroethane | 200 | 5 U | 1 | 1 | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 5 U |
| 1,2,4-Trimethylbenzene | | 5 U | 2 | 1 | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 5 U |
| 1,3,5-Trimethylbenzene | | 5 U | 5 J | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 5 U |
| Styrene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 5 U |
| 2-Butanone | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | R | R |
| Vinyl Chloride | 2 | 20 | 20 | 24 | 26 | 30 | 18 | 21 | 40 | 26 | 46 | 45 |
| DISSOLVED TAL METALS (ug/L) | | | | | | | | | | | | |
| Arsenic | 50 | 18.6 J | 18.4 | 14.5 | 15.0 | 15.5 | 16.6 | 14.2 | 14.4 J | 18.6 | 17.6 J | 11.8 J |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.42 U | 1.42 U | 1.42 U | 0.464 U | 0.464 UJ | 1.30 | 3.0 U J | 0.60 U J |
| TOTAL TAL METALS (ug/L) | | | | | | | | | | | | |
| Arsenic | 50 | 18.6 J | 17.9 | 14.7 | 15.6 | 15.9 | 16.1 | 16.0 | 15.9 J | 18.0 | 18.5 | 14.2 |
| Iron | 300 (1) | 55,400 | 57,600 | 49,400 | 44,500 | 48,400 | 49,000 | 43,100 | 49,000 | 35,400 | 37,200 | 36,600 |
| Lead | 15 | 1.0 U | 1.41 U | 1.74 J | 1.42 U | 1.42 U | 1.42 U | 3.53 | R | 3.47 | 3.0 U J | 0.60 U J |
| INDICATOR PARAMETERS (mg/L) | | | | | | | | | | | | |
| Ammonia | | 0.824 | 0.87 | 0.60 | 0.68 | 0.62 | 0.58 J | 1.3 | 0.564 J | 0.496 | 0.409 | 0.403 |
| COD | | 26.9 | 12.2 | 14 | <10 | 10 U | 17 | 23 | 18 | <5 | 5 U | 5 U |
| BOD | | 4.2 | 9.1 | 10 | 7 J | 12 | 17 J | 9 | 13 | 10 | 14.6 | 11.6 |
| Chloride | 250(1) | 6.00 J | 13.00 | 2 | 8.0 | 8 J | 7 | <5 | 5 | 373 | 7.41 | 7.02 |

Notes

- (1) - Secondary standards are as noted.
 - (2) - Source: Drinking Water Regulations and Health Advisories, Office of Water, USEPA Document Number EPA 822-B-96-002, October 1996
 - (4) - Concentrations for o-xylene and m, p - xylenes should be added for total xylene concentration. Total Xylene MCL is 10,000 ug/l.
- F:\om353\qpr\table1099.wb2 tab gwdata

TABLE 5: GROUNDWATER CHEMISTRY DATA, L&RR SUPERFUND SITE, JULY 2000 (CONTINUED)

| MONITORING WELL | MCL (2) | MW-103A | | | | | | | | | |
|------------------------------------|---------|--------------|---------------|-----------|--------------|---------------|-----------|--------------|------------|--------------|-----------|
| | | October 1996 | February 1997 | June 1997 | October 1997 | February 1998 | June 1998 | October 1998 | March 1999 | October 1999 | July 2000 |
| TCL VOLATILES (ug/L) | | | | | | | | | | | |
| Acetone | | 5 U | 5 U J | 1 U | R | R | 2 U J | 1 U | R | R | R |
| Benzene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 2 U | 1 U | 1 U | 0.5 U |
| Bromobenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloroethane | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U J | 1 U | 0.5 U |
| Chlorobenzene | 100 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloromethane | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U J | 1 U | 1 U J | 1 U | 0.5 U |
| 2-Chlorotoluene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Dichlorodifluoromethane | | 5 U | 1 U | 1 U | 2 U | 2 U J | 2 U | 2 U | 2 U J | 2 U J | 0.5 U |
| Trichlorofluoromethane | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,1 - Dichloroethane | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.1 J |
| 1,1 - Dichloroethene | 7 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloroethane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloropropane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| cis-1,2-Dichloroethene | 70 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| trans-1,2-Dichloroethene | 100 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2-dichlorobenzene | 600 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3-Dichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,4-Dichlorobenzene | 75 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Ethylbenzene | 700 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Isopropylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| p-Isopropyltoluene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| Methylene Chloride | 5 | 5 U | 2 U J | 2 U | 1 U | 2 U J | 2 U | 2 U J | 2 U | 2 U | 1 U J |
| Naphthalene | | 5 U | 2 U J | 2 U | 1 U | 2 U J | 2 U J | 2 U | 2 U J | 2 U J | 0.5 U |
| n-Propylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Tetrachloroethene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Trichloroethene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Toluene | 1000 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| o-Xylene | (4) | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| m,p Xylenes | (4) | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 1 U |
| 1,2,3 - Trichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U J | 1 U J | 0.5 U |
| 1,1,1 - Trichloroethane | 200 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2,4-Trimethylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3,5-Trimethylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Styrene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 2-Butanone | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | R |
| Vinyl Chloride | 2 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| DISSOLVED TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 1.4 J | 3.31 U | 3.32 U | 3.32 U | 3.32 U | 3.32 U | R | 1.08 U | 2.87 U | 0.75 U |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.42 U | 1.42 U | 1.42 U | 0.464 U | 0.464 U J | 0.464 U | 0.60 U J |
| TOTAL TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 1.2 J | 3.31 U | 3.32 U | 3.32 U | 3.32 U | 3.32 U | R | 1.08 U | 2.31 U | 0.80 J |
| Iron | 300 (1) | 55.3 | 162 | 106 U | 29.8 J | 104 U | 24.4 U J | 116 | 126 | 68.2 J | 13.8 J |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.42 U | 1.42 U | 1.42 U | 0.690 J | 0.464 U J | 0.490 J | 0.60 U J |
| INDICATOR PARAMETERS (mg/L) | | | | | | | | | | | |
| Ammonia | | 0.04 U | 0.040 U | 0.040 U | 0.040 U | 0.040 U | 0.05 U J | 0.28 | 0.313 | 0.0400 U | 0.0400 U |
| COD | | 10.0 U J | <10.0 | 10 U | 10 | 10 U | <10 | <5 | <10 | <5 | 5 U |
| BOD | | 2 U | <2.0 | 2 U | 2 U J | <2 | <2 | <2 | <2 | 96 | 2.0 U J |
| Chloride | 250(1) | 3.00 | 13 | 7 | 5 | 5 | 3 | <5 | 3 | 3.65 | 3.58 |

Notes:

(1) - Secondary standards are as noted

(2) - Source, Drinking Water Regulations and Health Advisories, Office of Water, USEPA Document Number EPA 822-B-96-002, October 1996

(4) - Concentrations for o-xylene and m, p - xylenes should be added for total xylene concentration. Total Xylene MCL is 10,000 ug/L.

TABLE 5: GROUNDWATER CHEMISTRY DATA, L&RR SUPERFUND SITE, JULY 2000 (CONTINUED)

| MONITORING WELL | MCL (2) | MW-104A | | | | | | | | | | Split Sample July 2000 |
|------------------------------------|---------|--------------|---------------|-----------|--------------|---------------|-----------|--------------|------------|--------------|-----------|------------------------|
| | | October 1996 | February 1997 | June 1997 | October 1997 | February 1998 | June 1998 | October 1998 | March 1999 | October 1999 | July 2000 | |
| TCL VOLATILES (ug/L) | | | | | | | | | | | | |
| Acetone | | 5 U | 5 U J | 5 U | R | R | 12 J | 2 U | R | R | 160 U J | 16 U J |
| Benzene | 5 | 18 | 14 | 15 | 16 | 17 | 18 | 12 | 16 J | 12 | 24 | 25 J |
| Bromobenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 0.4 J |
| Chloroethane | | 110 | 58 | 62 | 45 | 48 | 44 | 24 | 40 J | 1 U | 49 | 44 J |
| Chlorobenzene | 100 | 7 | 5 | 7 | 8 | 9 | 10 | 8 | 9 J | 7.0 | 10 | 10 |
| Chloromethane | | 5 U | 1 U | 0.9 J | 1 U | 2 U | 1 U J | 1 U | 1 U J | 1 U | 5 U | 0.5 U |
| 2-Chlorotoluene | | 5 U | 1 U | 4 | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| Dichlorodifluoromethane | | 5 U | 1 U | 1 U | 2 U | 2 U | 2 U | 2 U | 2 U J | 11 J | 5 U | 0.2 J |
| Trichlorofluoromethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| 1,1 - Dichloroethane | | 45 | 36 | 38 | 32 | 36 | 41 | 27 | 42 | 27 | 40 | 41 J |
| 1,1 - Dichloroethene | 7 | 5 U | 1 U | 1 U | 1 | 2 U | 1 U | 1 J | 1 | 1 U | 5 U | 0.5 U |
| 1,2 - Dichloroethane | 5 | 13 | 11 | 10 | 9 | 9 | 12 | 8 | 11 | 1 U | 9 | 9 |
| 1,2 - Dichloropropane | 5 | 9 | 9 | 10 | 9 | 2 U | 10 | 7 | 10 | 1 U | 10 | 10 |
| cis-1,2-Dichloroethene | 70 | 5 | 1 U | 1 U | 1 U | 2 U | 1 U | 1 | 1 U | 1 U | 5 U | 0.9 |
| trans-1,2-Dichloroethene | 100 | 5 | 1 U | 2 | 1 U | 1 J | 1 | 0.9 J | 1 | 1.1 | 1.8 J | 2 |
| 1,2-dichlorobenzene | 600 | 5 U | 4 | 5 | 6 | 6 | 6 | 6 | 5 | 5.6 | 6.6 | 6 J |
| 1,3-Dichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 J | 1 U | 1 U | 0.88 J | 0.9 U J |
| 1,4-Dichlorobenzene | 75 | 18 | 15 | 17 | 20 | 22 | 26 | 24 | 1 U | 21 | 33 | 27 J |
| Ethylbenzene | 700 | 150 | 120 | 150 J | 140 | 130 J | 150 | 130 | 120 | 104 | 140 | 110 J |
| Isopropylbenzene | | 38 | 28 | 31 | 24 | 27 | 22 | 17 | 15 | 17 | 23 | 24 |
| p-Isopropyltoluene | | 10 J | 10 | 9 | 12 | 13 | 12 | 9 | 6 | 5 U | 5 U | 12 |
| Methylene Chloride | 5 | 5 U | 7 J | 7 J | 4 J | 4 J | 2 U | 1 J | 2 U | 2.2 J | 18 U J | 4 U J |
| Naphthalene | | 46 | 2 U J | 48 | 35 | 25 J | 53 | 41 | 25 J | 36 J | 59 | 58 J |
| n-Propylbenzene | | 5 U | 3 | 4 | 5 | 5 | 4 | 3 | 4 | 7.1 | 4.8 J | 5 |
| Tetrachloroethene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| Trichloroethene | 5 | 5 U | 0.7 J | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| Toluene | 1000 | 13 | 10 | 10 | 9 | 9 | 8 | 8 | 11 J | 9.1 | 11 | 12 |
| o-Xylene | (4) | 60 | 1 U | 62 J | 59 | 51 J | 62 | 43 | 46 | 44 | 63 | 60 J |
| m,p Xylenes | (4) | 140 | 120 J | 150 J | 150 | 130 J | 150 | 100 | 120 | 120 | 350 J | 130 J |
| 1,2,3 - Trichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U J | 0.5 J | 1 U J | 1 U J | 5 U | 0.5 U J |
| 1,1,1 - Trichloroethane | 200 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| 1,2,4-Trimethylbenzene | | 27 | 24 | 26 | 32 | 35 | 34 | 28 | 26 | 1 U | 42 | 40 J |
| 1,3,5-Trimethylbenzene | | 8 | 25 | 9 | 22 | 10 | 9 | 8 | 1 U | 24 J | 12 | 13 |
| Styrene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 2.1 J | 0.5 U |
| 2-Butanone | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 10 J |
| Vinyl Chloride | 2 | 5 U | 2 | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 5 U | 1 |
| DISSOLVED TAL METALS (ug/L) | | | | | | | | | | | | |
| Arsenic | 50 | 82.0 | 96.8 J | 79.5 | 106 J | 82.5 J | 79.7 J | 68.5 | 116 J | 98.9 J | 86.9 | 113.0 |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.42 U | 1.42 U | 1.42 U | 0.464 U | 0.464 U | 0.464 U | 3.0 U J | 2.3 J |
| TOTAL TAL METALS (ug/L) | | | | | | | | | | | | |
| Arsenic | 50 | 92.9 | 76.4 J | 85.0 | 92.9 J | 62.7 J | 71.2 J | 61.7 | 107 J | 72.3 J | 94.7 | 113 |
| Iron | 300 (1) | 115,000 | 98,400 | 117,000 | 86,400 | 104,000 | 104,000 | 126,000 | 136,000 | 102,000 | 89,100 | 99,200 |
| Lead | 15 | 8.6 | 6.92 | 16.5 | 5.79 | 10.6 | 10.0 | 42.3 | 16.1 J | 10.2 | 4.2 J | 11.3 J |
| INDICATOR PARAMETERS (mg/L) | | | | | | | | | | | | |
| Ammonia | | 22.8 | 20.5 | 19.0 | 18.8 | 18.3 | 20 J | 87 | 21.5 | 21.7 | 22.6 | 15.8 |
| COD | | 222 | 200 | 300 | 190 | 73 | 190 | 250 | 190 J | 151 | 220 | 118 |
| BOD | | 24 J | 42 J | 51 J | 30 J | 23 | 20 | 28 | 24 | 257 J | 22.3 | 27.0 J |
| Chloride | 250(1) | 344 | 320 | 380 | 370 | 360 | 250 | 330 | 340 | 374 | 419 | 386 |

Notes:

- (1) - Secondary standards are as noted.
- (2) - Source: Drinking Water Regulations and Health Advisories, Office of Water, USEPA Document Number EPA 822-B-98-002, October 1996
- (4) - Concentrations for o-xylene and m, p - xylenes should be added for total xylene concentration. Total Xylene MCL is 10,000 ug/L.

TABLE 5: GROUNDWATER CHEMISTRY DATA, L&RR SUPERFUND SITE, JULY 2000 (CONTINUED)

| MONITORING WELL | MCL (2) | MW-201 | | | | | | | | | |
|------------------------------------|---------|--------------|---------------|-----------|--------------|---------------|-----------|--------------|------------|--------------|-----------|
| | | October 1996 | February 1997 | June 1997 | October 1997 | February 1998 | June 1998 | October 1998 | March 1999 | October 1999 | July 2000 |
| TCL VOLATILES (ug/L) | | | | | | | | | | | |
| Acetone | | 5 U | 5 U J | 5 U | R | R | 2 U J | 2 U J | R | R | R |
| Benzene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Bromobenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloroethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U J | 1 U | 0.5 U |
| Chlorobenzene | 100 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloromethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U J | 1 U J | 1 U J | 1 U | 0.5 U |
| 2-Chlorotoluene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Dichlorodifluoromethane | | 5 U | 1 U | 1 U | 3 | 3 | 3 | 4 J | 2 U J | 2 U J | 4.6 |
| Trichlorofluoromethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,1 - Dichloroethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,1 - Dichloroethene | 7 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloroethane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloropropane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| cis-1,2-Dichloroethene | 70 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| trans-1,2-Dichloroethene | 100 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2-dichlorobenzene | 600 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3-Dichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,4-Dichlorobenzene | 75 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Ethylbenzene | 700 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Isopropylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| p-Isopropyltoluene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| Methylene Chloride | 5 | 5 U | 2 U J | 2 U | 1 U | 2 U | 2 U | 2 U J | 2 U | 2 U | 1 U J |
| Naphthalene | | 5 U | 16 J | 2 U | 1 U | 2 U | 2 U J | 1 U | 2 U J | 2 U J | 0.5 U |
| n-Propylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U | 1 U | 0.5 U |
| Tetrachloroethane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U | 1 U | 0.5 U |
| Trichloroethene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Toluene | 1000 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| o-Xylene | (4) | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| m,p Xylenes | (4) | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 1 U |
| 1,2,3 - Trichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U J | 1 U J | 0.5 U |
| 1,1,1 - Trichloroethane | 200 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2,4-Trimethylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3,5-Trimethylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Styrene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 2-Butanone | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | R |
| Vinyl Chloride | 2 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| DISSOLVED TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 1.2 U | 3.31 U | 3.32 U | 3.32 U | 3.32 U | 3.3 U | 1.70 U | 1.08 U J | 1.08 U | 0.75 U |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.42 U | 1.42 U | 1.4 U | 0.464 U | 0.464 U | 0.464 U | 0.60 U J |
| TOTAL TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 1.2 U | 3.31 U | 3.32 U | 3.32 U | 3.32 U | 3.3 U | 1.08 U | 1.08 U J | 1.08 U | 0.75 U |
| Iron | 300 (1) | 353 | 51.1 | 332 J | 100 | 324 | 221 | 261 | 395 | 1,190 | 1,530 |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.42 U | 1.42 U | 1.4 U | 0.464 U | 0.464 U J | 0.464 U | 2.5 J |
| INDICATOR PARAMETERS (mg/L) | | | | | | | | | | | |
| Ammonia | | 0.04 U | 0.040 U | 0.040 U | 0.040 U | 0.040 U | 0.5 U J | 0.13 | 0.0970 | 0.0400 U | 0.0400 U |
| COD | | 10.0 U | <10.0 | 10 U | <10 | 7 | <10 | <5 | <10 | <5 | 5 U |
| BOD | | 2.0 U | <2.0 | 4 U J | 2 U J | <2 | <2 | <2 | <2 | <2 | 2 U J |
| Chloride | 250(1) | 6.00 J | 2.0 | 4 | 2 | 7 | <2 | <5 | <2 | 3.00 U | 3.00 U |

Notes

- (1) - Secondary standards are as noted.
- (2) - Source: Drinking Water Regulations and Health Advisories, Office of Water, USEPA Document Number EPA 822-B-96-002, October 1996
- (4) - Concentrations for o-xylene and m, p - xylenes should be added for total xylene concentration. Total Xylene MCL is 10,000 ug/L.

TABLE 5: GROUNDWATER CHEMISTRY DATA, L&RR SUPERFUND SITE, JULY 2000 (CONTINUED)

| MONITORING WELL | MCL (2) | MW-202 | | | | | | | | | |
|------------------------------------|---------|--------------|---------------|-----------|--------------|---------------|-----------|--------------|------------|--------------|-----------|
| | | October 1996 | February 1997 | June 1997 | October 1997 | February 1998 | June 1998 | October 1998 | March 1999 | October 1999 | July 2000 |
| TCL VOLATILES (ug/L) | | | | | | | | | | | |
| Acetone | | 5 U | 5 U J | 5 U | R | R | 4 J | 2 U J | R | R | R |
| Benzene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Bromobenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloroethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U J | 1 U | 0.5 U |
| Chlorobenzene | 100 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloromethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U J | 1 U J | 1 U J | 1 U | 0.5 U |
| 2-Chlorotoluene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Dichlorodifluoromethane | | 5 U | 1 U | 1 U | 2 U | 2 U | 2 U J | 2 U J | 2 U J | 2 U | 0.5 U |
| Trichlorofluoromethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U | 1 U J | 0.5 U |
| 1,1 - Dichloroethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,1 - Dichloroethene | 7 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloroethane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloropropane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| cis-1,2-Dichloroethene | 70 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| trans-1,2-Dichloroethene | 100 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2-dichlorobenzene | 600 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U J | 0.5 U |
| 1,3-Dichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U J | 0.5 U |
| 1,4-Dichlorobenzene | 75 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U J | 0.5 U |
| Ethylbenzene | 700 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Isopropylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| p-Isopropyltoluene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 5 U J | 0.5 U |
| Methylene Chloride | 5 | 5 U | 2 U J | 2 U | 1 U | 2 U | 2 U | 2 U J | 2 U | 2 U | 1 U J |
| Naphthalene | | 5 U | 2 U J | 2 U | 1 U | 2 U | 2 U J | 1 U | 2 U J | 2 U J | 0.5 U |
| n-Propylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U | 1 U J | 0.5 U |
| Tetrachloroethene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U | 1 U | 0.5 U |
| Trichloroethene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Toluene | 1000 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| o-Xylene | (4) | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| M,P Xylenes | (4) | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 1 U |
| 1,2,3 - Trichlorobenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U J | 1 U | 0.5 U |
| 1,1,1 - Trichloroethane | 200 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2,4-Trimethylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3,5-Trimethylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U J | 0.5 U |
| Styrene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 2-Butanone | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | R |
| Vinyl Chloride | 2 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| DISSOLVED TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 1.2 U | 3.31 U | 3.32 U | 3.32 U | 3.32 U | 3.32 U | 1.80 U | 1.08 U | 1.20 U | 0.75 U |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.42 U | 1.42 U | 1.42 U | 0.464 U | 0.464 U J | 0.464 U | 0.82 J |
| TOTAL TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 2.1 J | 3.31 U | 3.32 U | 3.32 U | 3.32 U | 3.32 U | 1.08 J | 1.08 U | 1.52 U | 0.75 U |
| Iron | 300 (1) | 14,900 | 7,080 | 1,940 | 1,490 | 4,310 | 4,150 | 3,220 | 1,610 | 2,040 | 7,790 |
| Lead | 15 | 9.6 | 11.4 | 2.95 | 2.12 J | 2.97 | 2.3 J | 1.96 U | 0.910 J | 1.27 | 0.6 U J |
| INDICATOR PARAMETERS (mg/L) | | | | | | | | | | | |
| Ammonia | | 0.04 U | 0.040 U | 0.040 U | 0.040 U | 0.040 U | 0.05 U J | 0.30 | 0.292 | 0.765 | 0.0400 U |
| COD | | 10.0 U | <10.0 | 10 U | <10 | 10 U | <10 | <5 | <10 | <5 | 5 U |
| BOD | | 2.0 U | <2.0 | 2 U | 2 U J | <2 | <2 | <2 | <2 | <2 | 2.0 U |
| Chloride | 250(1) | 12.0 J | 8.0 | 12 | 9 | 11 | 9 | 14 | 10 | 12.3 | 14.9 |

Notes:

(1) - Secondary standards are as noted.

(2) - Source: Drinking Water Regulations and Health Advisories, Office of Water, USEPA Document Number EPA 822-B-96-002, October 1996

(4) - Concentrations for o-xylene and m, p - xylenes should be added for total xylene concentration. Total Xylene MCL is 10,000 ug/L.

TABLE 5: GROUNDWATER CHEMISTRY DATA, L&RR SUPERFUND SITE, JULY 2000 (CONTINUED)

| MONITORING WELL | MCL (2) | CW-5B | | | | | | | | | |
|------------------------------------|---------|--------------|---------------|-----------|--------------|---------------|-----------|--------------|------------|--------------|-----------|
| | | October 1996 | February 1997 | June 1997 | October 1997 | February 1998 | June 1998 | October 1998 | March 1999 | October 1999 | July 2000 |
| TCL VOLATILES (ug/L) | | | | | | | | | | | |
| Acetone | | 8 J | 5 U J | 5 U | R | R | 2 U J | 2 U | R | R | R |
| Benzene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 0.6 J | 1 U | 1 U | 0.55 |
| Bromobenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloroethane | | 5 U | 1 U | 1 U | 1 U | 2 U J | 3 | 3 | 1 U J | 1 U | 1.8 |
| Chlorobenzene | 100 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloromethane | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U J | 1 U | 0.5 U |
| 2-Chlorotoluene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U J | 1 U | 0.5 U |
| Dichlorodifluoromethane | | 80 J | 44 | 45 | 26 | 19 J | 22 | 13 | 1 J | 6.2 J | 8.7 |
| Trichlorofluoromethane | | 5 U | 1 U | 7 | 3 | 2 J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,1 - Dichloroethane | | 5 U | 18 | 16 | 22 | 19 J | 24 | 24 | 30 | 15 | 19 |
| 1,1 - Dichloroethene | 7 | 5 U | 0.6 J | 0.8 J | 1 U | 2 U J | 1 U | 0.5 J | 1 U | 1 U | 0.35 J |
| 1,2 - Dichloroethane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloropropane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| cis-1,2-Dichloroethene | 70 | 5 U | 9 | 1 U | 11 | 2 U J | 1 U | 12 | 14 | 6.6 | 11 |
| trans-1,2-Dichloroethene | 100 | 5 U | 1 U | 0.8 J | 1 U | 2 U J | 1 U | 0.6 J | 1 U | 1 U | 0.44 J |
| 1,2-dichlorobenzene | 600 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3-Dichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,4-Dichlorobenzene | 75 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Ethylbenzene | 700 | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 0.8 J | 0.9 J | 1 U | 0.12 J |
| Isopropylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.11 J |
| p-Isopropyltoluene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| Methylene Chloride | 5 | 5 U | 19 J | 15 J | 1 U | 6 J | 2 U | 2 U J | 7 U | 1.8 J | 3.8 U J |
| Naphthalene | | 5 U | 2 U J | 2 U | 1 U | 2 U J | 2 U J | 0.5 J | 2 U J | 2 U J | 0.5 U |
| n-Propylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Tetrachloroethene | 5 | 18 | 25 | 28 | 40 | 26 J | 36 | 35 | 47 | 29 | 24 J |
| Trichloroethene | 5 | 5 U | 11 | 11 | 14 | 10 J | 14 | 13 | 14 | 7.5 | 9.4 |
| Toluene | 1000 | 5 U | 1 | 1 J | 1 U | 2 U J | 1 U | 0.9 J | 1 U | 1 U | 0.5 U |
| o-Xylene | (4) | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.32 J |
| m,p Xylenes | (4) | 5 U | 0.6 J | 0.8 J | 1 | 2 U J | 2 | 2 | 2 | 1 U | 3.7 J |
| 1,2,3 - Trichlorobenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U J | 1 U J | 5 U |
| 1,1,1 - Trichloroethane | 200 | 6 | 5 | 5 | 4 | 3 J | 3 | 2 | 1 U | 1 U | 1 |
| 1,2,4-Trimethylbenzene | | 5 U | 1 U J | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3,5-Trimethylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Styrene | | 5 U | 1 U | 1 U | 1 U | 2 U J | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 2-Butanone | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | R |
| Vinyl Chloride | 2 | 5 U | 4 | 4 | 8 | 7 J | 6 | 6 | 12 | 4.4 | 6.8 |
| DISSOLVED TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 1.20 U | 3.31 U | 3.32 U | 3.32 U | 3.32 U | 3.32 U | 2.18 | 1.08 U | 1.95 U | 0.75 U |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.42 U | 1.42 U | 1.42 U | 0.464 U | 0.870 U | 0.464 U | 0.6 U J |
| TOTAL TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 1.20 U | 3.31 U | 3.32 U | 3.32 U | 3.32 U | 3.32 U | 4.94 | 1.08 U J | 1.61 U | 0.75 U |
| Iron | 300 (1) | 233 | 521 | 620 | 1,770 | 4,790 | 4070 | 4,340 | 3,560 | 2,330 | 3,010 |
| Lead | 15 | 1.0 U | 1.41 U | 2.91 | 1.42 U | 1.42 U | 1.42 U | 0.590 J | 0.464 U J | 0.464 U | 0.60 U J |
| INDICATOR PARAMETERS (mg/L) | | | | | | | | | | | |
| Ammonia | | 21.1 | 0.040 U | 0.040 U | 0.052 | 0.040 U | 0.05 U J | 0.13 | 0.0980 | 0.0400 U | 0.0400 U |
| COD | | 10.0 U | <10.0 | 10 U | <10 | 10 U | 13 | 10 | <10 | <5 | 5 U |
| BOD | | 2.0 U | <2.0 | 2 J | 2 J | 5 | 3 J | 7 | 7 | 12 | 4.6 |
| Chloride | 250(1) | 3.00 | 3.0 | 4 | 2 | 4 | 2 | <5 | <2 | 3.00 U | 3.00 U |

Notes

- (1) - Secondary standards are as noted.
- (2) - Source: Drinking Water Regulations and Health Advisories, Office of Water, USEPA Document Number EPA 822-B-96-002, October 1996
- (4) - Concentrations for o-xylene and m, p - xylenes should be added for total xylene concentration. Total Xylene MCL is 10,000 ug/L.

TABLE 5: GROUNDWATER CHEMISTRY DATA, L&RR SUPERFUND SITE, JULY 2000 (CONTINUED)

| MONITORING WELL | MCL (2) | CW-7A | | | | | | | | | |
|------------------------------------|---------|--------------|---------------|-----------|--------------|---------------|-----------|--------------|------------|--------------|-----------|
| | | October 1996 | February 1997 | June 1997 | October 1997 | February 1998 | June 1998 | October 1998 | March 1999 | October 1999 | July 2000 |
| TCL VOLATILES (ug/L) | | | | | | | | | | | |
| Acetone | | 5 U | 5 U J | 5 U | R | R | 2 U J | 2 U J | R | R | R |
| Benzene | 5 | 5 U | 1 | 1 | 2 | 1 J | 1 | 1 | 1 J | 1 U | 0.26 J |
| Bromobenzene | | 5 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Chloroethane | | 5 U | 0.7 J | 1 U | 1 | 2 U | 1 U | 1 J | 1 U J | 1 U | 0.15 J |
| Chlorobenzene | 100 | 5 U | 2 | 4 | 6 | 5 | 4 | 4 | 3 | 2 | 0.75 |
| Chloromethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U J | 1 U J | 1 U J | 1 U | 0.5 U |
| 2-Chlorotoluene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Dichlorodifluoromethane | | 5 U | 1 U | 1 U | 2 U | 2 U | 2 U | 2 U J | 2 U J | 2 U J | 0.5 U |
| Trichlorofluoromethane | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,1 - Dichloroethane | | 5 U | 0.8 J | 0.9 J | 1 | 2 U | 1 U | 0.6 J | 1 U | 1 U | 0.35 J |
| 1,1 - Dichloroethene | 7 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloroethane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2 - Dichloropropane | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| cis-1,2-Dichloroethene | 70 | 5 U | 1 U | 1 U | 0.9 J | 2 U | 1 U | 0.7 J | 1 U | 1 U | 0.12 J |
| trans-1,2-Dichloroethene | 100 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2-dichlorobenzene | 600 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3-Dichlorobenzene | | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,4-Dichlorobenzene | 75 | 5 U | 2 | 3 | 6 | 6 | 4 | 4 | 1 U | 2.0 | 0.83 |
| Ethylbenzene | 700 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Isopropylbenzene | | 5 U | 2 | 1 U | 3 | 3 | 2 | 2 | 1 U | 1 U | 0.24 J |
| p-Isopropyltoluene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 5 U | 0.5 U |
| Methylene Chloride | 5 | 5 U | 2 U J | 2 U | 1 U | 2 U | 2 U | 2 U J | 2 U | 2 U | 1 U J |
| Naphthalene | | 5 U | 2 U J | 2 J | 1 U | 2 | 2 U J | 1 U | 2 U J | 3.9 J | 0.5 U |
| n-Propylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U | 1 U | 0.5 U |
| Tetrachloroethene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U J | 1 U | 1 U | 0.5 U |
| Trichloroethene | 5 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| Toluene | 1000 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| o-Xylene | (4) | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| m,p Xylenes | (4) | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 1 U |
| 1,2,3 - Trichlorobenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U J | 1 U J | 0.5 U |
| 1,1,1 - Trichloroethane | 200 | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,2,4-Trimethylbenzene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 1,3,5-Trimethylbenzene | | 5 U | 0.6 J | 1 U | 1 U | 2 U | 1 U | 0.7 J | 1 U | 1 U | 0.5 U |
| Styrene | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| 2-Butanone | | 5 U | 1 U | 1 U | 1 U | 2 U | 1 U | 1 U | 1 U | 1 U | R |
| Vinyl Chloride | 2 | 5 U | 1 U | 1 U | 0.9 J | 2 U | 1 U | 1 U | 1 U | 1 U | 0.5 U |
| DISSOLVED TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 10.0 | 9.36 | 8.76 | 8.3 | 10.9 | 11.5 | R | 9.51 J | 9.37 | 9.1 |
| Lead | 15 | 1.0 U | 1.41 U | 1.42 U | 1.4 U | 1.42 U | 1.42 U | 0.464 U | 0.464 U | 0.464 U | 3.0 U J |
| TOTAL TAL METALS (ug/L) | | | | | | | | | | | |
| Arsenic | 50 | 10.0 | 9.47 | 8.39 | 9 | 10.3 | 11.9 | R | 10.3 J | 8.98 | 9.4 |
| Iron | 300 (1) | 25,400 | 31,500 | 35,900 | 36,100 | 29,200 | 31,700 | 30,800 | 24,900 | 18,000 | 15,200 |
| Lead | 15 | 1.0 U | 1.41 U | R | 1.4 U | 1.42 U | 1.42 U | 0.464 U | 0.464 U J | 0.464 U | 3.0 U J |
| INDICATOR PARAMETERS (mg/L) | | | | | | | | | | | |
| Ammonia | | 0.04 U | 0.55 | 0.62 | 0.0 | 0.77 | 0.6 J | 1.8 | 0.934 | 0.681 | 0.260 |
| COD | | 33.9 J | 28.3 | 42 | 30 | 13 | 24 | 30 | 14 | 5 | 5 U |
| BOD | | 4 | 6.9 | 8 | 6 J | 7 | 8 J | 6 | 7 | 9 | 2.6 J |
| Chloride | 250(1) | 26.0 | 32 | 17 | 33 | 4 J | 18 | 22 | 12 | 11.2 | 7.05 |

Notes:

(1) - Secondary standards are as noted

(2) - Source: Drinking Water Regulations and Health Advisories, Office of Water, USEPA Document Number EPA 822-B-96-002, October 1996

(4) - Concentrations for o-xylene and m, p - xylenes should be added for total xylene concentration. Total Xylene MCL is 10,000 ug/L.

TABLE 6: GROUNDWATER CHEMISTRY SPLIT SAMPLE COMPARISONS, L&RR SUPERFUND SITE, JULY 2000

| MONITORING WELL WELL | MW-104A Severn Trent Lab. | MW - 104A SPLIT Compuchem Environmental | RPD % Orig./Split | Orig./Split Comparable Yes/No (1) |
|------------------------------------|------------------------------|---|----------------------|---|
| TCL VOLATILES (ug/L) | | | | |
| Acetone | 160 U J | 16 U J | | Yes |
| Benzene | 24 | 25 J | 4.1 | Yes |
| Bromobenzene | 5 U | 0.4 J | | Yes |
| Chloroethane | 49 | 44 J | 10.8 | Yes |
| Chlorobenzene | 10 | 10 | 0.0 | Yes |
| Chloromethane | 5 U | 0.5 U | | Yes |
| 2-Chlorotoluene | 5 U | 0.5 U | | Yes |
| Dichlorodifluoromethane | 5 U | 0.2 J | | Yes |
| Trichlorofluoromethane | 5 U | 0.5 U | | Yes |
| 1,1 - Dichloroethane | 40 | 41 J | 2.5 | Yes |
| 1,1 - Dichloroethane | 5 U | 0.5 U | | Yes |
| 1,2 - Dichloroethane | 9 | 9 | 0.0 | Yes |
| 1,2 - Dichloropropane | 10 | 10 | 0.0 | Yes |
| cis-1,2-Dichloroethene | 5 U | 0.9 | | Yes |
| trans-1,2-Dichloroethene | 1.8 J | 2 | 10.5 | Yes |
| 1,2-dichlorobenzene | 6.6 | 6 J | 9.5 | Yes |
| 1,3-Dichlorobenzene | 0.88 J | 0.9 U J | | Yes |
| 1,4-Dichlorobenzene | 33 | 27 J | 20.0 | Yes |
| Ethylbenzene | 140 | 110 J | 24.0 | Yes |
| Isopropylbenzene | 23 | 24 | 4.3 | Yes |
| p-isopropyltoluene | 13 | 12 | 8.0 | Yes |
| Methylene Chloride | 5 U | 12 | | Yes |
| Naphthalene | 18 U J | 4 U J | | Yes |
| n-Propylbenzene | 59 | 58 J | 1.7 | Yes |
| Tetrachloroethene | 4.8 J | 5 | 4.1 | Yes |
| Trichloroethene | 5 U | 0.5 U | | Yes |
| Toluene | 5 U | 0.5 U | | Yes |
| o-Xylene | 11 | 12 | 8.7 | Yes |
| m,p Xylenes | 63 | 60 J | 4.9 | Yes |
| 1, 2, 3 - Trichlorobenzene | 350 J | 130 J | 91.7 | No |
| 1, 1, 1 - Trichloroethane | 5 U | 0.5 U J | | Yes |
| 1,2,4-Trimethylbenzene | 5 U | 0.5 U | | Yes |
| 1,3,5-Trimethylbenzene | 42 | 40 J | 4.9 | Yes |
| Styrene | 12 | 13 | 8.0 | Yes |
| 2-Butanone | 2.1 J | 0.5 U | | No |
| Vinyl Chloride | 5 U | 10 J | | Yes |
| | 5 U | 1 | | Yes |
| DISSOLVED TAL METALS (ug/L) | | | | |
| Arsenic | | | | |
| Lead | 86.9 | 113.0 | 26.1 | Yes |
| | 3.0 U J | 2.3 J | | Yes |
| TOTAL TAL METALS (ug/L) | | | | |
| Arsenic | | | | |
| Iron | 94.7 | 113 | 17.6 | Yes |
| Lead | 89,100 | 99,200 | 10.7 | Yes |
| | 4.2 J | 11.3 J | 91.6 | No |
| INDICATOR PARAMETERS (mg/L) | | | | |
| Ammonia | | | | |
| COD | 22.6 | 15.8 | 35.4 | No |
| BOD | 220 | 118 | 60.4 | No |
| Chloride | 22.3 | 27.0 J | 19.1 | Yes |
| | 419 | 386 | 8.2 | Yes |

Notes:

(1) - Comparability analysis performed according to the Sampling and Analysis Plan, L&RR Superfund Site, prepared by de maximus, inc., September 1996

TABLE 7: SURFACE WATER CHEMISTRY DATA, L&RR SUPERFUND SITE, APRIL 2000

| MONITORING LOCATION Sample Designation | Surface Water Criteria Fresh Water (ug/L) Criterion | | SW-5 | SW-8 | SW-10 | SW-16 | LCH-4 | Duplicate Sample of LCH-4 | LCH-5 |
|---|---|------------|---------|-------|--------|--------|--------|---------------------------------|--------|
| | Maximum | Continuous | | | | | | | |
| TCL VOLATILES (ug/L) | | | | | | | | | |
| 1,1-Dichloroethane | | | 0.5 U | 0.8 | 0.5 U | 0.5 U | 23 | 24 | 0.5 U |
| Chloroethane | | | 0.5 U | 1 | 0.8 | 0.8 | 17 | 19 | 0.5 U |
| p-Isopropyltoluene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.5 | 0.6 | 0.5 U |
| 1,2-Dichlorobenzene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.7 | 0.8 | 0.5 U |
| 1,4-Dichlorobenzene | | | 0.5 U | 3 | 0.5 U | 0.5 U | 2 | 2 | 0.5 U |
| Toluene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.7 | 0.8 | 0.5 U |
| m,p Xylene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 3 | 3 | 0.5 U |
| o-Xylene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 4 | 5 | 0.5 U |
| Isopropylbenzene | | | 0.5 U | 2 | 0.5 U | 0.5 U | 12 | 14 | 0.5 U |
| Ethylbenzene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 9 | 10 | 0.5 U |
| Vinyl Chloride | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 3 | 3 | 0.5 U |
| Acetone | | | R | R | R | R | 2 J | 2 J | R |
| Methylene Chloride | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.7 | 0.8 | 0.5 U |
| trans-1,2-Dichloroethene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.7 | 0.7 | 0.5 U |
| cis-1,2-Dichloroethene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 8 | 8 | 0.5 U |
| Benzene | | | 0.5 U | 0.7 | 0.5 U | 0.5 U | 3 | 3 | 0.5 U |
| 1,2-Dichloroethane | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.9 | 1 | 0.5 U |
| Trichloroethene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 1 | 1 | 0.5 U |
| 1,2-Dichloropropane | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 5 | 6 | 0.5 U |
| Chlorobenzene | | | 0.5 U | 2 | 0.5 U | 0.5 U | 3 | 3 | 0.5 U |
| 1,3,5-Trimethylbenzene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.8 | 1.0 | 0.5 U |
| 1,2,4-Trimethylbenzene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 2 | 2 | 0.5 U |
| Naphthalene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 7 | 7 | 0.5 U |
| n-Propylbenzene | | | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.5 UJ | 0.5 J | 0.5 U |
| TOTAL METALS (ug/L) | | | | | | | | | |
| Arsenic | 360 | 190 | 0.85 UJ | 23.2 | 0.85 U | 0.85 U | 6.6 J | 3.2 J | 4.4 |
| DISSOLVED METALS (ug/L) | | | | | | | | | |
| Arsenic | 360 | 190 | 1.1 J | 11.0 | 0.85 U | 0.85 U | 5.6 J | 8.8 J | 0.85 U |
| WET CHEMISTRY PARAMETERS | | | | | | | | | |
| Chlorides | | | 10.5 | 58.3 | 30.2 | 29.4 | 81 J | 49.6 J | 16.3 |
| FIELD PARAMETERS | | | | | | | | | |
| pH (Standard units) | | | 5.7 | 6.39 | 6.93 | 6.79 | 6.1 | 6.07 | 6.56 |
| Temperature (C) | | | 8.29 | 12.0 | 9.4 | 10.6 | 11 | 10.64 | 12.4 |
| Conductivity (mhos/cm) | | | 443 | 631 | 270 | 259 | 372 | 372 | 83.7 |

Notes:

(1) - Concentrations reference Federal Register Vol. 57, No. 246, December 22, 1992, Rules and Regulations 60911

TABLE 8: AMBIENT AIR SAMPLING DATA, L&RR SUPERFUND SITE, JUNE 8 AND JUNE 9, 2000

| Sampling Location | AS - 1 Upwind | AS - 1 DUPLICATE Upwind | AS - 2 Upwind | AS - 3 Downwind | AS - 4 Downwind |
|--|------------------|----------------------------|------------------|--------------------|--------------------|
| Start date | 6-8-00 | 6-8-00 | 6-8-00 | 6-8-00 | 6-8-00 |
| Start time (1) | 11:20 | 11:20 | 11:24 | 11:29 | 11:33 |
| Canister start pressure(Field) | -30.00 | -30.00 | -30.00 | -30.00 | -29.90 |
| End date | 6-9-00 | 6-9-00 | 6-9-00 | 6-9-00 | 6-9-00 |
| End time (2) | 9:20 | 9:20 | 9:24 | 9:29 | 9:33 |
| Canister end pressure (Field) | -7.50 | -7.25 | -7.00 | -7.00 | -6.90 |
| TO - 14 Compounds Detected (ppbv) * | | | | | |
| Chloromethane | 1.5 | 0.86 U | 0.80 U | 0.78 U | 0.84 U |
| Freon 12 | 1.0 | 0.86 U | 0.82 | 0.78 U | 0.91 |
| Methylene Chloride | 1.4 | 1.4 | 0.89 | 1.4 B | 0.90 |
| Toluene | 0.86 U | 4.9 | 0.80 U | 0.78 U | 0.84 U |
| Ethyl Benzene | 0.86 U | 3.3 | 0.80 U | 0.78 U | 0.84 U |
| m, p - xylene | 0.86 U | 23 | 0.80 U | 0.78 U | 0.84 U |
| o-Xylene | 0.86 U | 3.8 | 0.80 U | 0.78 U | 0.84 U |
| 1,3,5-Trimethylbenzene | 0.86 U | 2.1 | 0.80 U | 0.78 U | 0.84 U |
| 1,2,4-Trimethylbenzene | 0.86 U | 1.0 | 0.80 U | 0.78 U | 0.84 U |
| 1,4-Dichlorobenzene | 0.86 U | 0.99 | 0.80 U | 0.78 U | 0.84 U |
| 1,2-Dichlorobenzene | 0.86 U | 0.94 | 0.80 U | 0.78 U | 0.84 U |
| Chloroform | 0.86 U | 0.84 U | 0.80 U | 0.78 U | 0.84 U |
| Carbon Tetrachloride | 0.86 U | 0.84 U | 0.80 U | 0.78 U | 0.84 U |
| Acetone | 16 | 36 | 23 | 6.8 | 7.5 |
| Carbon Disulfide | 3.4 U | 4.1 | 3.2 U | 3.1 U | 3.4 U |
| 2-Butanone(MEK) | 3.4 U | 5.9 | 3.3 | 3.1 U | 3.4 U |

Notes:

* - Parts per billion by volume.

(1) - Start time weather information: 75-80 degrees F., Wind speed 0-4mph; Barometric pressure 30.13 inches.

(2) - End time weather information: 75-80 degrees F, Wind speed 0-9 mph; Barometric pressure 29.97 inches.

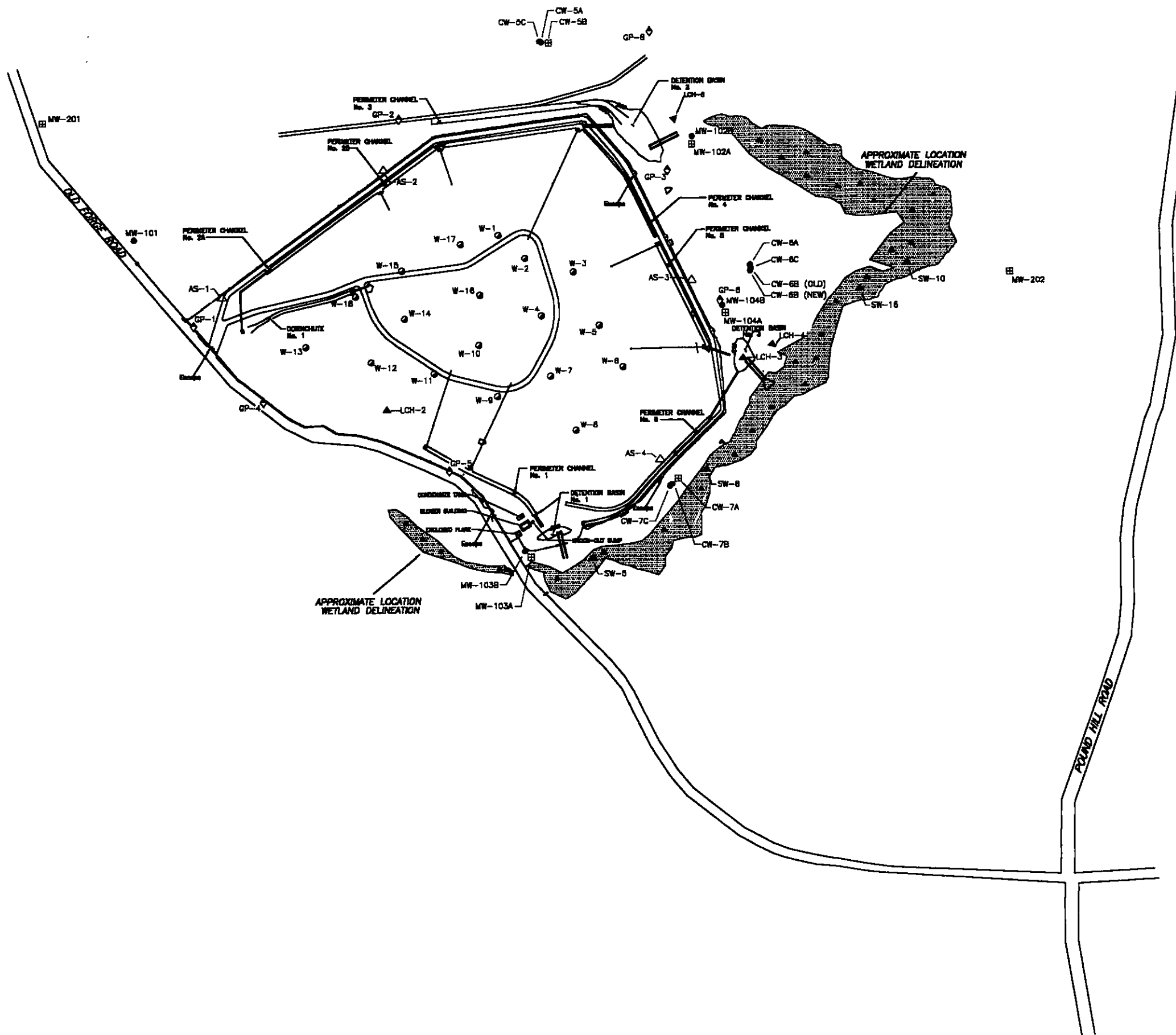
f:\om\353\qpro\tabs7200.wb2 TAB: AIR

TABLE 9: FLARE INLET SAMPLING DATA, L&RR SUPERFUND SITE, AUGUST 15 2000

| Sampling Location | Flare Inlet |
|--|-----------------|
| Start date | August 15, 2000 |
| End date | August 15, 2000 |
| TO - 14 Compounds Detected (ppbv) * | |
| Freon 12 | 220 |
| Freon 114 | 156 |
| Methylene Chloride | 1337 |
| Vynyl Chloride | 1680 |
| Ethyl Chloride | 205 |
| Dichloromethane | 682 |
| 1,1-Dichloroethane | 504 |
| cis, 1,2-Dichloroethylene | 34 |
| Benzene | 267 |
| Trichloroethylene | 349 |
| Toluene | 56300 |
| Trans-1,3-Dichloropropene | 22 |
| Tetrachloroethylene | 488 |
| Chlorobenzene | 47 |
| Ethyl Benzene | 3730 |
| 1,3-Xylene+1,4-Xylene | 7030 |
| 1,2-Xylene | 1610 |
| 4-Ethyltoluene | 251 |
| Non-target Compounds Detected | |
| Trans-1,2-Dichloroethylene | 1080 |

Notes:

Figure



LEGEND:

- W-1 GAS WELL
- ◇ GP-6 GAS PROBE
- MW-104B GROUNDWATER MONITORING WELL
- POST-CLOSURE GROUNDWATER MONITORING WELL
- ▲ SW-5 SURFACE WATER SAMPLING LOCATION
- △ AS-1 AMBIENT AIR MONITORING LOCATION

NOTE:

- 1) TOPOGRAPHIC DATA, LOCATIONS, AND PROPERTY BOUNDARIES SURVEYED BY DANSON SURVEYING AND ENGINEERING COMPANY, NOVEMBER 1994, SEPTEMBER 1995 AND SEPTEMBER 1998.
- 2) ftamal - Feet above mean sea level.



SAMPLE LOCATION PLAN

L&RR SUPERFUND SITE
NORTH SMITHFIELD, RHODE ISLAND

PREPARED FOR

THE L&RR GROUP

O & M, Inc. Environmental Operations and Maintenance Management

| | | |
|-----------------|----------|----------------------|
| DATE: 1-25-00 | FIGURE 1 | DRAWING NAME |
| SCALE: AS SHOWN | | SAMPLE LOCATION PLAN |

Appendix A

Inspection Reports

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith

Date: 2/15/2000 Time On Site: 11:00 AM

Weather: Clear

Signature: Clayton Smith

| CORRECTIVE MEASURES | | | | | | |
|--|---|--|-------------------------------|-------------------------|------------------------------|---------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition (Except where noted) | Surface Feature (see note) | Sloughing (see note) | Document and Observe area | 2/15/00 |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition | | | | |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Leaving Tilting/Leaving | Good Condition | | | | |
| COMMENTS: A small area of sloughing, approximately 10' X 10' and one foot deep was observed near gas well #8. The area maintains positive drainage and was not ponding water. The area will be observed for any further sloughing. | | | | | | |

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith

Date: 3/15/2000 Time On Site: 8:00 AM

Weather: Clear

Signature: Clayton Smith

| CORRECTIVE MEASURES | | | | | | |
|--|---|-------------------------------------|----------------------------|---------------------------|---------------------|---------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition (Except where noted) | Surface Feature (see note) | Animal Burrows (see note) | Filled burrow holes | 3/15/00 |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition | | | | |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Heaving Tilting /Heaving | Good Condition | | | | |
| COMMENTS: Four small burrow holes were noted on the landfill surface. Corrective action was to fill each hole and apply grass seed to the area. The small area of sloughing, noted during the Febuary Inspection (3/15/2000), near gas well #8 was repaired (filled with topsoil and seeded) on 3/3/2000. Routine maintenace to Blower Number 2 was completed this period. | | | | | | |

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith
 Date: 4/20/2000 Time On Site: 8:00 AM
 Weather: Clear
 Signature: *Clayton Smith*

| CORRECTIVE MEASURES | | | | | | |
|---|---|--|-------------------------------|-------------------------|-----------------------------|---------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition (Except where noted) | Surface Feature (see note) | Sloughing (see note) | Scheduled corrective action | 4/20/00 |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition | | | | |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Heaving Tilting /Heaving | Good Condition | | | | |
| <p>COMMENTS: Three small areas of sloughing were noted on the landfill surface. The areas are approximately 20 LF apart from each other and located on a fairly level area approximately 50 LF northeast of gas probe #4. The first area noted is approximately 12 feet in diameter and 1.5 feet deep, the second area is approximately 8 feet in diameter and .5 feet deep and the third area is approximately 5 feet in diameter and .5 foot deep. Corrective action will be to fill each area with topsoil and apply grass seed to any disturbed areas.</p> | | | | | | |

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith

Date: 5/25/2000 Time On Site: 8:00 AM

Weather: Clear

Signature: Clayton Smith

| CORRECTIVE MEASURES | | | | | | |
|--|---|-------------------------------------|--------------------------------|-------------------------------------|---|----------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition | | | | |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition (Except where noted) | LFG Extraction Well (see note) | Well Head & Flex Coupler (see note) | Closed Well #13 and Scheduled corrective action | 5/25/000 |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Heaving Tilting /Heaving | Good Condition | | | | |
| COMMENTS: The three small areas of sloughing noted on the 4/20/00 Inspection Log were repaired. Extraction Well #13 recorded high percent oxygen content and was closed on 5/25/00. The flex coupler on extraction well #7 appeared to be overstressed. Extension of the flex coupler on extraction well #7 was scheduled for 5/1/00. The extent of any necessary corrective action repair to extracton Well #13 will be determined on 5/1/00 and the well will remain closed until determination of further corrective action can be made. | | | | | | |

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith

Date: 6/30/2000 Time On Site: 8:00 AM

Weather: Clear

Signature: Clayton Smith

| CORRECTIVE MEASURES | | | | | | |
|--|---|----------------|------------------|------------------------|--------|------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition | | | | |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition | | | | |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Heaving Tilting /Heaving | Good Condition | | | | |
| COMMENTS: The hyper extended flex coupler and corrective action repair inspection noted on the 5/25/00-inspection log concerning gas extraction wells #7 and # 13 were completed on 5/1/00. | | | | | | |

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith
 Date: 7/24/00 Time On Site: 8:30 AM
 Weather: Clear
 Signature: *Clayton Smith*

| CORRECTIVE MEASURES | | | | | | |
|--|---|----------------|------------------|------------------------|--------|------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition | | | | |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition | | | | |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Heaving Tilting /Heaving | Good Condition | | | | |
| COMMENTS: | | | | | | |
| None | | | | | | |

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith
 Date: 8/14/00 Time On Site: 8:30 AM
 Weather: Cloudy/Lt. Rain
 Signature: *Clayton Smith*

| CORRECTIVE MEASURES | | | | | | |
|--|---|--|------------------|------------------------------|--------------------------------|------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition | | | | |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition (Except where noted) | Flare | Intermittent flare shutdowns | Scheduled John Zink Technician | |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Heaving Tilting/Heaving | Good Condition | | | | |
| COMMENTS: The flare has been experiencing intermittent shutdowns. A John Zink Co. technician is scheduled to complete the semi-annual flare inspection and to further investigate the intermittent flare shutdown problem the week of August 28. | | | | | | |

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith
 Date: 9/12/00 Time On Site: 8:30 AM
 Weather: Clear
 Signature: C. Smith

| CORRECTIVE MEASURES | | | | | | |
|---|---|----------------|------------------|------------------------|--------|------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition | | | | |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition | | | | |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Heaving Tilting/Heaving | Good Condition | | | | |
| <p>COMMENTS: As noted in the August 14, 2000 inspection report, the flare was experiencing intermittent shutdowns and a John Zink Co. technician was scheduled to investigate the problem. The John Zink technician determined that the cause for the intermittent shutdowns was a faulty louvre control. The louvre control automatically regulates the air intake to the flare which in turn controls the flare temperature. At times, the faulty louver did not allow enough air to the flare causing a low temperature flare shutdown. The louvre control was replaced during the week of August 28, 2000. The flare has been in continuous operation following replacement of the louver control.</p> | | | | | | |

**TABLE 2-1
L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: Clayton Smith

Date: 10/19/00 Time On Site: 8:30 AM

Weather: Clear

Signature: Clayton Smith

| CORRECTIVE MEASURES | | | | | | |
|--|---|--|----------------------|------------------------|----------------------|------------------------------|
| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
| 1. Security System a. Gate b. Fence c. Locks d. Signs | Inoperative Holes Inoperative Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | Animal Burrows, Other Holes, Cracks Washouts and Sloughing Brushes/Tree Growth, Bare Spots Washouts and Discoloration | Good Condition | | | | |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | Ponding Water, Filling and Sediment Filling with Sediment, Blocked by Debris Filling with Sediment Filling with Sediment, Riprap Lining Disturbed Blocked, Damaged, Riprap Outlets Disturbed Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | Broken, No Lock Cracked, Missing Cracked, Missing Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | Physical Damage to Casing, Wellhead, Sampling Port Physical Damage to Casing Recording Paper and Pens Empty | Good Condition (Except where noted) | LFG Extraction Wells | Strained Flex Coupler | Schedule Maintenance | Target Maint. Date: 10/20/00 |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | Tilting/Heaving Tilting/Heaving | Good Condition | | | | |
| COMMENTS: As noted on the 10/19/00 Field Well Monitoring Data Sheet (Data Sheet), methane concentration in monitoring probes #1 and #4 were above criteria established in the PC/O&M Plan. Also noted on the Data Sheet the flex couplers on gas wells 12, 14 and 18 were strained. Immediate corrective action was to schedule maintenance of the flex couplers noted. | | | | | | |

**L&RR SUPERFUND SITE
INSPECTION LOG**

Inspector Name: John Hunt
 Date: 11/16/00 Time On Site: 0800
 Weather: SUNNY
 Signature: [Signature]

CORRECTIVE MEASURES

| Feature | Trouble Signs | Status | Problem Location | Description of Problem | Action | Date |
|--|---|----------------|------------------|------------------------|--------|------|
| 1. Security System a. Gate b. Fence c. Locks d. Signs | a. Inoperative b. Holes c. Inoperative d. Missing, Unreadable | Good Condition | | | | |
| 2. Cover Integrity a. Surface Features b. Slopes c. Vegetation d. Breakouts | a. Animal Burrows, Other Holes, Cracks b. Washouts and Sloughing c. Brushes/Tree Growth, Bare Spots d. Washouts and Discoloration | Good Condition | | | | |
| 3. Stormwater Management System a. Diversion Swales b. Catch Basins c. Stilling Wells d. Perimeter Channels e. Culverts f. Detention Basins | a. Ponding Water, Filling and Sediment b. Filling with Sediment, Blocked by Debris c. Filling with Sediment d. Filling with Sediment, Riprap Lining Disturbed e. Blocked, Damaged, Riprap Outlets Disturbed f. Filling with Sediment, Riprap Outlets Disturbed | Good Condition | | | | |
| 4. Groundwater Monitoring Wells a. Locking Cap b. Protective Casing c. Concrete Collar d. Local Erosion | a. Broken, No Lock b. Cracked, Missing c. Cracked, Missing d. Ponding, Water Channels | Good Condition | | | | |
| 5. Landfill Gas Monitoring and Collection System a. LFG Extraction Wells b. LFG Migration Probes c. Control Panel | a. Physical Damage to Casing, Wellhead, Sampling Port b. Physical Damage to Casing c. Recording Paper and Pens Empty | Good Condition | | | | |
| 6. Permanent Monuments a. Bench Marks b. Settlement Monuments | a. Tilting/Heaving b. Tilting/Heaving | Good Condition | | | | |

COMMENTS:



Appendix B

Gas Monitoring Data

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 2/15/00

By: C. Smith

Weather Conditions: Clear Temp: 40° Barometric Pressure: 29.9

System Conditions: Good Control Temp (TIC201): 1603 Flame Temp (TISH101): 1589 Flow (cfm): 426

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|-------|---------------------|---------------------|--------------------|-----------|--------------------------------|-------------|
| 1 | 11:00 | 52.7 | 43.1 | 0.2 | 78° | -12.1 | |
| 2 | | 53.7 | 41.0 | 0.1 | 101 | -5.1 | |
| 3 | | 48.6 | 41.5 | 0.0 | 88 | -7.8 | |
| 4 | | 50.5 | 38.2 | 0.1 | 90 | -5.8 | |
| 5 | | 35.6 | 40.2 | 0.0 | 38 | -5.0 | |
| 6 | | 3.0 | 11.5 | 11.2 | 45 | -0.3 | Well Closed |
| 7 | | 54.2 | 42.1 | 0.0 | 40 | -7.1 | |
| 8 | | 0.5 | 16.7 | 19.5 | 40 | -1.2 | Well Closed |
| 9 | | 40.1 | 37.1 | 0.0 | 80 | -7.1 | |
| 10 | | 50.1 | 38.2 | 0.1 | 82 | -6.0 | |
| 11 | | 43.2 | 38.5 | 0.0 | 70 | -7.2 | |
| 12 | | 53.3 | 36.2 | 0.0 | 88 | -5.8 | |
| 13 | | 50.5 | 43.2 | 0.0 | 80 | -1.3 | |
| 14 | | 52.5 | 35.1 | 0.0 | 80 | -5.7 | |
| 15 | | 50.5 | 40.6 | 0.1 | 80 | -6.5 | |
| 16 | | 0.5 | 0.3 | 16.7 | 35 | -0.4 | Well Closed |
| 17 | | 55.4 | 42.0 | 0.1 | 90 | -7.2 | |
| 18 | 12:30 | 57.8 | 40.2 | 0.0 | 76 | -6.8 | |
| Probe | | | | | | | |
| 1 | | 0.0 | 3.4 | 17.8 | | | |
| 2 | | | | | | | |
| 3 | | | | | | | |
| 4 | | 0.0 | 5.4 | 14.9 | | | |
| 5 | | | | | | | |
| 6 | | | | | | | |
| 7 | | | | | | | |
| 8 | | 0.0 | 1.1 | 15.4 | | | |
| Flare | | 54.5 | 42.7 | 0.0 | | | |
| Blower | | 50.8 | 41.8 | 0.2 | | | |

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 3/16/00

By: Clayton Smith

Weather Conditions: Clear Temp: 52° Barometric Pressure: 29.9

System Conditions: Good Control Temp (TIC201): 1608 Flame Temp (TISH101): 1580 Flow (cfm): 435

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|-------------|
| 1 | 1030 | 53.5 | 42.9 | 0.0 | 84° | -10.4 | |
| 2 | | 50.6 | 41.2 | 0.3 | 106 | -8.0 | |
| 3 | | 36.6 | 33.8 | 1.2 | 92 | -1.1 | |
| 4 | | 46.7 | 36.5 | 0.4 | 90 | -8.0 | |
| 5 | | 38.2 | 28.4 | 5.4 | 42 | -0.5 | |
| 6 | | 0.7 | 8.5 | 15.5 | 50 | 0.0 | Well closed |
| 7 | | 41.9 | 32.3 | 0.0 | 98 | -1.3 | |
| 8 | | 0.0 | 0.0 | 19.0 | 50 | -2.3 | Well closed |
| 9 | | 42.6 | 34.6 | 0.0 | 92 | -4.2 | |
| 10 | | 48.7 | 38.9 | 0.0 | 88 | -1.6 | |
| 11 | | 41.3 | 34.5 | 0.0 | 90 | -2.3 | |
| 12 | | 32.2 | 30.9 | 0.0 | 98 | -0.3 | |
| 13 | | 45.7 | 38.3 | 0.0 | 62 | -0.6 | |
| 14 | | 55.0 | 40.8 | 0.1 | 84 | -0.2 | |
| 15 | | 51.1 | 41.9 | 0.0 | 82 | -3.5 | |
| 16 | | 5.2 | 3.6 | 19.2 | 86 | -1.4 | Well closed |
| 17 | | 52.8 | 41.5 | 0.0 | 92 | -3.6 | |
| 18 | | 54.2 | 41.0 | 0.0 | 80 | -2.4 | |
| Probe | | | | | | | |
| ① | | 0.0 | 3.6 | 18.6 | | | |
| ② | | | | | | | |
| ③ | | | | | | | |
| ④ | | 0.0 | 1.4 | 18.7 | | | |
| ⑤ | | | | | | | |
| ⑥ | | | | | | | |
| ⑦ | | | | | | | |
| ⑧ | | 0.0 | 2.7 | 9.6 | | | |
| Flare | ▼ | 46.0 | 36.9 | 0.0 | | | |
| Blower | 1330 | 46.8 | 38.7 | 0.2 | | | |

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 4/18/2000

By: C. Smith

Weather Conditions: Cloudy Temp: 50° Barometric Pressure: 30.1

System Conditions: Good Control Temp (TIC201): 160.1 Flame Temp (TISH101): 1537 Flow (cfm): 420

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|-------------|
| 1 | 0830 | 51.7 | 38.8 | 0.0 | 82 | -7.2 | |
| 2 | | 53.6 | 38.9 | 0.1 | 108 | -4.2 | |
| 3 | | 50.1 | 39.1 | 0.1 | 90 | -2.1 | |
| 4 | | 53.4 | 36.1 | 6.3 | 90 | -3.1 | |
| 5 | | 55.7 | 41.4 | 0.2 | 40 | -0.8 | |
| 6 | | 3.0 | 7.2 | 15.6 | 50 | -0.1 | Well closed |
| 7 | | 55.5 | 42.0 | 0.0 | 100 | -0.5 | |
| 8 | | 53.9 | 40.4 | 12.0 | 40 | -0.2 | Well closed |
| 9 | | 47.6 | 35.4 | 0.2 | 85 | -1.0 | |
| 10 | | 41.8 | 31.8 | 0.0 | 85 | -1.4 | |
| 11 | | 55.2 | 32.4 | 0.4 | 78 | -6.7 | |
| 12 | | 53.8 | 41.3 | 0.0 | 90 | -1.6 | |
| 13 | | 55.1 | 41.1 | 0.0 | 50 | -2.1 | |
| 14 | | 49.7 | 38.7 | 0.0 | 80 | -2.4 | |
| 15 | | 50.7 | 38.1 | 0.0 | 84 | -2.1 | |
| 16 | | 55.3 | 38.3 | 15.8 | 40 | -0.9 | Well closed |
| 17 | | 49.7 | 36.7 | 0.0 | 92 | -2.7 | |
| 18 | | 51.8 | 38.9 | 0.0 | 80 | -2.8 | |
| Probe | | | | | | | |
| ① | | 0.0 | 6.1 | | | | |
| 2 | | | | | | | |
| 3 | | | | | | | |
| ④ | | 0.0 | 4.3 | | | | |
| 5 | | | | | | | |
| 6 | | | | | | | |
| 7 | | | | | | | |
| ⑧ | | 0.0 | 2.0 | 15.8 | | | |
| Flare | | 49.9 | 40.8 | 0.1 | | | |
| Blower | 1045 | 47.7 | 38.7 | 1.1 | | | |

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 5/25/00

By: C. Smith

Weather Conditions: Cloudy Temp: 64°F Barometric Pressure: _____

System Conditions: Good Control Temp (TIC201): 1602 Flame Temp (TISH101): 1684 Flow (cfm): 450

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|-----------------------------|
| 1 | 0800 | 45.5 | 39.6 | 0.0 | 92° | -16.1 | |
| 2 | | 36.6 | 34.9 | 0.0 | 110 | -12.6 | |
| 3 | | 21.4 | 26.2 | 0.0 | 90 | -2.1 | |
| 4 | | 33.0 | 31.2 | 0.0 | 90 | -12.8 | |
| 5 | | 45.7 | 40.1 | 0.4 | 52 | -1.2 | |
| 6 | | 0.4 | 9.1 | 15.1 | 60 | -0.6 | Well Closed |
| 7 | | 44.0 | 36.2 | 1.7 | 100 | -4.8 | Flex Coupler Rigid |
| 8 | | 0.1 | 0.0 | 18.6 | 60 | -3.6 | Well Closed |
| 9 | | 22.1 | 26.5 | 0.1 | 92 | -8.5 | |
| 10 | | 38.5 | 36.9 | 0.0 | 89 | -7.7 | |
| 11 | | 32.1 | 29.8 | 0.0 | 80 | -7.3 | |
| 12 | | 29.4 | 32.1 | 0.0 | 98 | -6.7 | |
| 13 | | 0.3 | 0.2 | 18.8 | 75 | -2.6 | Closed Well this Inspection |
| 14 | | 43.6 | 39.5 | 0.0 | 90 | -6.3 | |
| 15 | | 38.0 | 34.4 | 0.0 | 80 | -7.1 | |
| 16 | | 38.6 | 36.5 | 15.4 | 62 | -6.5 | Well Closed |
| 17 | | 29.2 | 35.0 | 0.3 | 98 | -6.7 | |
| 18 | | 42.6 | 37.5 | 0.0 | 80 | -5.3 | |
| Frce | | | | | | | |
| 1 | | 0.0 | 6.4 | 10.3 | | | |
| 2 | | | | | | | |
| 3 | | | | | | | |
| 4 | | 0.0 | 3.0 | 15.5 | | | |
| 5 | | | | | | | |
| 6 | | | | | | | |
| 7 | | | | | | | |
| 8 | | 0.0 | 7.8 | 14.6 | | | |
| Flare | | 38.5 | 34.2 | 0.2 | | | |
| Blower | 1130 | 37.6 | 32.8 | 0.8 | | | |

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 6/8/00

By: C. Smith

Weather Conditions: Clear/Calm Temp: _____ Barometric Pressure: 30.1

System Conditions: good Control Temp (TIC201): 1595 Flame Temp (TISH101): 1655 Flow (cfm): 425

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|-------------|
| 1 | 0800 | 47.5 | 39.9 | 0.8 | 92 | -14.2 | |
| 2 | | 41.2 | 39.8 | 0.0 | 110 | -11.2 | |
| 3 | | 30.8 | 29.9 | 0.3 | 100 | -9.2 | |
| 4 | | 32.8 | 29.0 | 1.3 | 90 | -11.1 | |
| 5 | | 40.7 | 31.8 | 4.0 | 60 | -5.5 | |
| 6 | | 0.7 | 14.7 | 14.5 | 60 | -0.6 | Well closed |
| 7 | | 49.1 | 40.6 | 0.6 | 100 | -10.2 | |
| 8 | | 0.3 | 0.1 | 19.9 | 60 | -3.9 | Well closed |
| 9 | | 22.7 | 26.7 | 0.2 | 90 | -7.8 | |
| 10 | | 38.7 | 34.4 | 0.1 | 90 | -7.0 | |
| 11 | | 33.2 | 30.1 | 0.1 | 85 | -6.8 | |
| 12 | | 34.8 | 32.7 | 0.0 | 100 | -6.3 | |
| 13 | | 38.2 | 18.8 | 7.9 | 90 | -2.2 | Well closed |
| 14 | | 45.1 | 37.4 | 0.2 | 92 | -5.3 | |
| 15 | | 40.5 | 32.9 | 0.8 | 90 | -6.0 | |
| 16 | | 0.5 | 0.1 | 20.0 | 60 | -6.1 | Well closed |
| 17 | | 43.2 | 34.4 | 0.8 | 100 | -6.7 | |
| 18 | | 45.1 | 38.7 | 0.1 | 85 | -4.9 | |
| Probe | | | | | | | |
| 1 | | 0.0 | 4.8 | 12.8 | | | |
| 2 | | | | | | | |
| 3 | | | | | | | |
| 4 | | 0.0 | 3.6 | 16.1 | | | |
| 5 | | | | | | | |
| 6 | | | | | | | |
| 7 | | | | | | | |
| 8 | | 0.0 | 7.8 | 15.2 | | | |
| Flare | | 44.5 | 34.2 | 0.2 | | | |
| Blower | 1130 | 43.2 | 33.3 | 0.3 | | | |

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 7/24/00

By: C. Smith

Weather Conditions: Clear Temp: 65° Barometric Pressure: 30.1

System Conditions: Good Control Temp (TIC201): 1618 Flame Temp (TISH101): 1716 Flow (cfm): 410

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|-------------|
| 1 | 0845 | 46.3 | 40.7 | 1.5 | 98 | -11.4 | |
| 2 | | 43.7 | 37.2 | 1.0 | 110 | -9.0 | |
| 3 | | 39.6 | 35.4 | 0.2 | 98 | -1.6 | |
| 4 | | 37.6 | 30.2 | 1.4 | 90 | -9.4 | |
| 5 | | 33.3 | 26.0 | 8.5 | 60 | -4.4 | |
| 6 | | 2.2 | 13.2 | 13.1 | 60 | -0.1 | Well Closed |
| 7 | | 50.2 | 40.8 | 0.3 | 100 | -8.3 | |
| 8 | | 0.2 | 0.0 | 20.6 | 60 | -3.3 | Well Closed |
| 9 | | 29.2 | 28.4 | 0.5 | 94 | -6.9 | |
| 10 | | 44.9 | 35.6 | 0.2 | 92 | -4.3 | |
| 11 | | 36.1 | 33.3 | 0.4 | 90 | -6.1 | |
| 12 | | 35.7 | 37.2 | 0.1 | 98 | -5.1 | |
| 13 | | 22.5 | 12.7 | 13.1 | 89 | -2.0 | Well Closed |
| 14 | | 47.2 | 40.6 | 0.1 | 89 | -3.4 | |
| 15 | | 43.6 | 34.1 | 0.4 | 90 | -8.6 | |
| 16 | | 0.5 | 0.0 | 20.3 | 75 | -5.7 | Well Closed |
| 17 | | 46.9 | 37.7 | 0.5 | 100 | -6.0 | |
| 18 | | 51.4 | 41.2 | 0.1 | 82 | -4.8 | |
| Probe | | | | | | | |
| ① | | 0.1 | 4.1 | 15.8 | | | |
| 2 | | | | | | | |
| 3 | | | | | | | |
| ④ | | 0.3 | 9.3 | 9.1 | | | |
| 5 | | | | | | | |
| 6 | | | | | | | |
| 7 | | | | | | | |
| ⑧ | | 0.0 | 8.4 | 15.1 | | | |
| Flare | | 46.4 | 37.8 | 0.1 | | | |
| Blower | 1330 | 44.2 | 36.7 | 0.3 | | | |

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 8/14/00

By: C. Smith

Weather Conditions: Cloudy / Lt. Rain Temp: 67° Barometric Pressure: 29.95

System Conditions: Fair Control Temp (TIC201): 1502 Flame Temp (TISH101): 1615 Flow (cfm): 440

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|-------------|
| 1 | 0900 | 55.2 | 44.0 | 0.7 | 98 | -4.5 | |
| 2 | | 52.7 | 40.7 | 0.2 | 110 | -1.8 | |
| 3 | | 46.3 | 40.2 | 0.3 | 98 | -2.1 | |
| 4 | | 42.6 | 39.8 | 0.8 | 92 | -3.3 | |
| 5 | | 36.6 | 42.9 | 0.4 | 65 | -0.1 | |
| 6 | | 5.0 | 43.8 | 16.1 | 62 | -4.7 | Well closed |
| 7 | | 52.3 | 42.4 | 0.3 | 100 | -0.4 | |
| 8 | | 1.3 | 0.2 | 20.1 | 60 | -0.2 | well closed |
| 9 | | 28.6 | 27.4 | 0.0 | 95 | -3.4 | |
| 10 | | 41.6 | 37.6 | 0.1 | 92 | -0.2 | |
| 11 | | 42.6 | 38.5 | 0.2 | 90 | -3.4 | |
| 12 | | 47.3 | 35.6 | 0.0 | 98 | -3.1 | |
| 13 | | 20.5 | 14.7 | 12.8 | 90 | -1.8 | well closed |
| 14 | | 41.9 | 40.6 | 0.0 | 90 | -2.6 | |
| 15 | | 48.6 | 42.8 | 0.1 | 90 | -2.4 | |
| 16 | | 0.1 | 2.8 | 18.6 | 75 | -3.5 | well closed |
| 17 | | 48.1 | 41.6 | 0.2 | 100 | -3.7 | |
| 18 | | 42.8 | 43.7 | 0.1 | 82 | -2.4 | |
| Probe | | | | | | | |
| 1 | | 1.4 | 17.8 | 6.9 | | | |
| 2 | | | | | | | |
| 3 | | | | | | | |
| 4 | | 5.6 | 15.4 | 4.8 | | | |
| 5 | | | | | | | |
| 6 | | | | | | | |
| 7 | | | | | | | |
| 8 | | 0.0 | 6.8 | 15.4 | | | |
| Flare | | 52.1 | 43.2 | 0.6 | | | |
| Blower | 1050 | 47.2 | 38.7 | 0.3 | | | |

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 9/21/00

By: C. Smith

Weather Conditions: cloudy Temp: 75° Barometric Pressure: 29.67

System Conditions: Good Control Temp (TIC201): 1587 Flame Temp (TISH101): 1542 Flow (cfm): 435

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|-------------|
| 1 | 1030 | 53.8 | 41.2 | 0.2 | 100 | -6.8 | |
| 2 | | 54.2 | 42.1 | 0.4 | 110 | -6.2 | |
| 3 | | 55.1 | 43.0 | 0.2 | 108 | -4.3 | |
| 4 | | 54.2 | 40.3 | 0.4 | 92 | -5.4 | |
| 5 | | 52.3 | 42.4 | 0.1 | 60 | -1.3 | |
| 6 | | 4.7 | 14.9 | 14.1 | 70 | -0.8 | Well closed |
| 7 | | 55.6 | 41.3 | 0.0 | 100 | -3.4 | |
| 8 | | 41.1 | 32.8 | 6.1 | 64 | -1.3 | Well closed |
| 9 | | 46.6 | 38.9 | 0.0 | 92 | -3.9 | |
| 10 | | 52.4 | 37.6 | 0.1 | 86 | -1.5 | |
| 11 | | 48.7 | 39.6 | 0.0 | 82 | -3.7 | |
| 12 | | 47.6 | 38.6 | 10.2 | 98 | -1.2 | |
| 13 | | 30.4 | 17.6 | 0.2 | 84 | -0.9 | |
| 14 | | 52.4 | 42.6 | 0.1 | 92 | -2.8 | |
| 15 | | 50.1 | 44.5 | 0.4 | 90 | -3.2 | |
| 16 | | 48.6 | 33.0 | 18.5 | 72 | -0.7 | Well closed |
| 17 | | 50.1 | 40.6 | 0.5 | 98 | -2.5 | |
| 18 | | 48.0 | 42.6 | 0.4 | 88 | -2.9 | |
| Probe | | | | | | | |
| 1 | | 0.8 | 6.4 | 15.6 | | | |
| 2 | | | | | | | |
| 3 | | | | | | | |
| 4 | | 0.9 | 1.8 | 16.7 | | | |
| 5 | | | | | | | |
| 6 | | | | | | | |
| 7 | | | | | | | |
| 8 | | 0.0 | 6.4 | 16.2 | | | |
| Flare | ↓ | 52.2 | 39.7 | 1.2 | | | |
| Blower | 1245 | 49.1 | 37.6 | 1.3 | | | |

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 10/19/00

By: C Smith

Weather Conditions: Clean Temp: _____ Barometric Pressure: 29.89

System Conditions: Good Control Temp (max): 1604 Flame Temp (max): 1627 Flow (cfm): 421

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|-----------------------|
| 1 | 0955 | 55.4 | 43.2 | 0.5 | 108 | -9.8 | |
| 2 | | 47.6 | 38.3 | 0.5 | 110 | -9.40 | |
| 3 | | 46.6 | 36.4 | 0.5 | 108 | -7.2 | |
| 4 | | 53.2 | 40.6 | 0.5 | 94 | -7.5 | |
| 5 | | 51.3 | 39.6 | 2.3 | 50 | -3.1 | Well Closed |
| 6 | | 4.4 | 17.2 | 13.1 | 58 | -0.6 | |
| 7 | | 55.1 | 42.1 | 0.8 | 96 | -5.60 | |
| 8 | | 8.0 | 5.9 | 17.3 | 50 | -1.60 | Well Closed |
| 9 | | 33.6 | 32.0 | 0.5 | 80 | -4.20 | |
| 10 | | 52.1 | 40.6 | 0.4 | 82 | -4.50 | |
| 11 | | 39.7 | 34.5 | 0.4 | 80 | -4.60 | |
| 12 | | 40.4 | 37.0 | 1.0 | 92 | -4.5 | Flex Coupler Strained |
| 13 | | 39.8 | 26.4 | 6.7 | 62 | -0.7 | |
| 14 | | 52.9 | 41.6 | 0.4 | 82 | -2.6 | Flex Coupler Strained |
| 15 | | 50.8 | 39.20 | 0.5 | 90 | -6.8 | |
| 16 | | 49.9 | 33.6 | 5.0 | 52 | -3.5 | Well Closed |
| 17 | | 50.3 | 38.4 | 1.0 | 92 | -4.3 | |
| 18 | | 55.6 | 41.5 | 0.8 | 80 | -3.0 | Flex Coupler Strained |
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(1) 23.3 24.7 0.5

(4) 17.4 21.3 4.6

(8) 0.0 1.5 16.8

Flare 48.0 34.3 1.4

Boiler 45.4 36.2 2.0

LANDFILL & RESOURCE RECOVERY SUPERFUND SITE FIELD GAS WELL MONITORING DATA SHEET

Date: 11/16/00

By: J Hunt

Weather Conditions: Sunny

Temp: 3°

Barometric Pressure: _____

System Conditions: Good Control Temp (TIC201): 1597

Flame Temp (TISH101): 1644

Flow (cfm): 379
1600

| Well No. | Time | CH ₄ (%) | CO ₂ (%) | O ₂ (%) | Temp (°F) | Pressure (in H ₂ O) | Comments |
|----------|------|---------------------|---------------------|--------------------|-----------|--------------------------------|----------|
| 1 | 0900 | 55.5 | 41.1 | 0.4 | 102° | -9.00 | |
| 2 | | 45.5 | 36.5 | 0.7 | 110° | -8.70 | |
| 3 | | 37.5 | 34.6 | 0.4 | 104° | -6.90 | |
| 4 | | 48.2 | 39.4 | 0.4 | 92° | -7.0 | |
| 5 | | 45.6 | 34.7 | 4.4 | 40° | -2.50 | |
| 6 | | 3.60 | 12.7 | 13.7 | 52° | +1.0 | |
| 7 | | 55.8 | 41.3 | 1.0 | 94° | -5.50 | |
| 8 | | 0 | 0.1 | 19.4 | 38°f | -1.2 | |
| 9 | | 29.1 | 30.5 | 0.8 | 80° | -4.30 | |
| 10 | | 51.0 | 39.1 | 0.4 | 80° | -3.80 -4.0 | |
| 11 | | 36.4 | 32.4 | 0.4 | 68° | -3.70 | |
| 12 | | 39.5 | 35.3 | 0.8 | 88° | -3.10 | |
| 13 | | 18.4 | 15.1 | 13.3 | 50° | -0.80 | |
| 14 | | 52.8 | 39.5 | 0.6 | 80° | -3.0 | |
| 15 | | 48.7 | 36.9 | 0.4 | 88° | -5.70 | |
| 16 | | 43.4 | 27.8 | 7.0 | 38° | -3.30 | |
| 17 | | 48.4 | 38.8 | 1.0 | 88° | -4.0 | |
| 18 | | 57.1 | 41.4 | 0.4 | 76° | -2.80 | |
| Probe | | | | | | | |
| 1 | | 5.3 | 19.5 | 2.70 | - | +2.30 | |
| 2 | | | | | | | |
| 3 | | | | | | +2.30 | |
| 4 | | 1.4 | 11.0 | 10.6 | - | +2.70 | |
| 5 | | | | | | | |
| 6 | | | | | | | |
| 7 | | | | | | | |
| 8 | | 0.0 | 1.5 | 17.0 | - | +1.30 | |
| are | | 44.1 | 35.0 | 1.1 | - | +6.40 | |
| Blower | 1050 | 41.4 | 33.3 | 2.4 | - | -15.40 | |

JOHN ZINK

A KOCH INDUSTRIES COMPANY
 1820 East Apache - Tulsa, Oklahoma 74121-1220
 918/234-1800 FAX 918/234-1968

Field Service Engineer:
Tim Levanduski

| | | | |
|-----------------|---------------|-----------------------------|--------|
| Original SO # | | New SO # | 910749 |
| Customer | de maximis | P.O. # | |
| Invoice Address | | Site Address | |
| Invoice City | Waltham | Site City | |
| State, Zip | MA | State, Zip | |
| Requested by | Clayton Smith | Site Contact & Phone Number | |

Are T&C's in place? Yes No If No, have the Customer sign KJZ T&C's prior to the start of ANY work. If Yes, what type of T&C's?
 Technical Asst Agmt PM Contract @ (rate) Blanket Contract @ (rate) Warranty Other

Type of Work: (Check at least one in each section)

| | | | |
|--|----------------------------------|-------------------------------------|--|
| <input type="checkbox"/> Internal | <input type="checkbox"/> TOSG | <input type="checkbox"/> Inspection | <input checked="" type="checkbox"/> PM |
| <input checked="" type="checkbox"/> External | <input type="checkbox"/> Burners | <input type="checkbox"/> Startup | <input type="checkbox"/> Sales Call |
| | <input type="checkbox"/> Flares | <input type="checkbox"/> Callout | <input type="checkbox"/> Non-JZ Equip |
| | <input type="checkbox"/> Vapor | <input type="checkbox"/> Training | <input type="checkbox"/> |

| Date | Time Interval | Chargeable Hours | | | Warranty Hours | Non Rev | Description of Work Performed |
|-----------------|---------------|------------------|------|----|----------------|---------|---|
| | | Reg. | Wknd | OT | | | |
| 8/9 | 7:30 | 10 | | | | | BI-ANNUAL INSPECTION OF Flare SYSTEM. Cleaned Flame arrester to Flare Checked Blowers, greased all Bearings. Cleaned Condensate gun and all strainers. Checked Full operation of Flare Unit running properly. |
| Sub-Total Hours | | 10 | | | | | |

Sub-Total Hours: 10
 Total Hours Worked: 10

RATE INFORMATION:
 Regular Days @ \$ _____ Weekend Days @ \$ _____
 Reg OT Hours @ \$ _____ Wknd OT Hours @ \$ _____

Required Parts and Sales Price _____
 Comments or Recommendations _____

Work Verified By: _____
 Work Submitted By: John Zink Representative Date: 8/9/00
 Phone (Fax) _____
 Add'l Info. Attached

Trip Expenses:

| | |
|---------------|--|
| Air Fare: | |
| Mileage: | |
| Tolls/Taxi: | |
| Car Rental: | |
| Hotel: | |
| Phone/Tips: | |
| Fuel/Parking: | |
| Meals: | |
| Misc.: | |
| Sub-Total: | |

(_____ Miles @ _____ /mile)

Meals:

| Date | Amount | Date | Amount |
|------|--------|------|--------|
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |

Leave Open
 Close

Failure Codes: _____
 Repair Codes: _____

Refer to Code List for Valid Codes

Labor Total: _____
 Expenses Total: _____
 Parts Total: _____
 Grand Total: _____

| | | | |
|-----------------|---------------|-----------------------------|----------------|
| Original SO # | | New SO # | 930289 |
| Customer | de maximis | P.O. # | |
| Invoice Address | | Site Address | LR+R Land fill |
| Invoice City | Waltham | Site City | N. Smithfield |
| State, Zip | MA | State, Zip | RI |
| Requested by | Clayton Smith | Site Contact & Phone Number | |

Field Service Engineer:
Tim Levanduski

Are T&C's in place? Yes No If No, have the Customer sign KJZ T&C's prior to the start of ANY work. If Yes, what type of T&C's?
 Technical Asst Agmt PM Contract @ (rate) Blanket Contract @ (rate) Warranty Other

Type of Work:

(Check at least one in each section)

- | | | | |
|--|--|---|---------------------------------------|
| <input type="checkbox"/> Internal | <input type="checkbox"/> TOSG | <input type="checkbox"/> Inspection | <input type="checkbox"/> PM |
| <input checked="" type="checkbox"/> External | <input type="checkbox"/> Burners | <input type="checkbox"/> Startup | <input type="checkbox"/> Sales Call |
| | <input checked="" type="checkbox"/> Flares | <input checked="" type="checkbox"/> Callout | <input type="checkbox"/> Non-JZ Equip |
| | <input type="checkbox"/> Vapor | <input type="checkbox"/> Training | <input type="checkbox"/> |

| Date | Time Interval | Chargeable Hours | | | Warranty Hours | Non Rev | Description of Work Performed |
|------|---------------|------------------|------|----|----------------|---------|--|
| | | Reg. | Wknd | OT | | | |
| 8/28 | 8:30 | 8 | | | | | Flare shutting down on Low Flare Temp. |
| | | | | | | | checked all Flame scanner operation -OK |
| | | | | | | | Found signal board inside Damper motor Control Defective |
| | | | | | | | The motor would go Full open with loss of signal on Board. |
| | | | | | | | Replaced Damper motor with spare at site. |
| | | | | | | | Unit operating properly. |

Sub-Total Hours 8
 Total Hours Worked: 8

RATE INFORMATION:
 Regular Days @ \$ _____
 Weekend Days @ \$ _____
 Reg OT Hours @ \$ _____
 Wknd OT Hours @ \$ _____

Required Parts and Sales Price _____
 Comments or Recommendations _____

Work Verified By: Not Available 8/28/00
 Customer Representative Date
 Work Submitted By: [Signature] 8/28/00
 John Zink Representative Date

Trip Expenses:

| | |
|---------------|--|
| Air Fare: | |
| Mileage: | |
| Tolls/Taxi: | |
| Car Rental: | |
| Hotel: | |
| Phone/Tips: | |
| Fuel/Parking: | |
| Meals: | |
| Misc.: | |
| Sub-Total: | |

130 Miles @ _____ mile)

Leave Open
 Close

Meals:

| Date | Amount | Date | Amount |
|------|--------|------|--------|
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |

Add'l Info. Attached

Refer to Code List for Valid Codes

Failure Codes: _____
 Repair Codes: _____

Labor Total: _____
 Expenses Total: _____
 Parts Total: _____
 Grand Total: _____



Appendix C

Groundwater Data Validation Reports Severn Trent Laboratories



TRILLIUM INC.
Consultants in Environmental Chemistry

356 FARRAGUT CROSSING DR.
KNOXVILLE, TN 37922

(423) 966-8880
FAX (423) 966-8885
cerikson@trilliuminc.com

September 11, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review Report - STL Report No. 7000-1470A - VOAs in Ground Waters

Dear Dan:

I have completed my evaluation of the volatile organics analysis data prepared by the Severn Trent Laboratory (STL) in Monroe, Connecticut, for eight ground water (GW) samples and two trip blanks (TB) from the L&RR Site, which were reported in a single data package under Report No. 7000-1470A. The data package was received by Trillium for review on August 28, 2000, with additional requested information received on September 1-2, 2000 and again on September 6, 2000. The following samples were reported:

| | |
|--------------|--------------|
| MW-201 | MW-202 |
| MW-103A | CW-7A |
| TBL+RR071200 | CW-5B |
| MW-104A | MW-102A |
| DUP-L+RR | TBL+RR071300 |

Sample identifications for both trip blanks were shortened by the laboratory throughout the data package by removing "L+RR" from the identifications found on the chain of custody records. The chain of custody identifications are listed above and are used throughout this evaluation report for all samples.

Analyses were performed according to EPA Method 8260B (SW-846, Third Edition, Update III); my evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced method, and the *Region I, EPA-New England Data Validation Functional Guidelines for Evaluating Environmental Analyses* (12/96). Qualifiers consistent with those defined in the Region I document were applied as necessary and appropriate.

HOME OFFICE:

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OFFICES IN:

LOUISIANA • MARYLAND • NEW JERSEY • NORTH CAROLINA • PENNSYLVANIA • TENNESSEE • TEXAS

Mr. Dan Garrigan

Evaluation of STL Report No. 7000-1470A (Volatiles in GW)

September 11, 2000

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No volatiles analysis raw data were provided in the data package. As specified by the QAPP, my review was limited to a Level III evaluation of the summary forms as presented.

Based on the evaluation, the following qualifiers were applied:

- Results for 2-butanone in all samples in this data set and for acetone in MW-201, MW-202, MW-103A, CW-7A, and CW-5B were rejected (R) as unreliable based on very low relative response factors in the associated calibration standards. Results for 2-butanone also warranted qualification as estimated due to unacceptably low recoveries in the associated blank spike, but the "R" qualifiers take precedence.
- Results for acetone in TBL+RR071200, TBL+RR071300, MW-104A, MW-102A, and DUP-L+RR were qualified as estimated (J) based on low relative response factors in the associated calibration standards.
- Results for tetrachloroethene in CW-5B, MW-102A, and DUP-L+RR were qualified as estimated (J) based on high percent relative standard deviations in the associated initial calibration.
- Results for methylene chloride in all samples were qualified as less than the reported values or less than the sample-specific reporting limit (RL), whichever was greater, based on associated method blank contamination and as estimated based on high percent relative standard deviations in the associated initial calibration and unacceptably high blank spike recoveries (UJ).
- Results for acetone in MW-102A, MW-104A, and DUP-L+RR were qualified as less than the reported values or less than the sample-specific RL, whichever was greater, based on associated method and trip blank contamination (U).
- Results for trichlorofluoromethane in MW-102A and DUP-L+RR were qualified as less than the reported values or less than the sample-specific RL, whichever was greater, based on associated method blank contamination (U).
- Results for m/p-xylenes in CW-5B and MW-104A were qualified as estimated (J) based on the unacceptably high blank spike recoveries; these results may be biased high.
- Results for 4-methyl-2-pentanone in all samples were qualified as estimated (UJ) based on unacceptably low recoveries in the associated blank spike.

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Evaluation of STL Report No. 7000-1470A (Volatiles in GW)

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- Reporting limits for vinyl chloride and methylene chloride were corrected to 0.5 µg/L and 1 µg/L, respectively (from 0.2 µg/L and 0.3 µg/L, respectively), to reflect the lowest concentration calibration standards run in association with these analyses.

All qualifiers are reflected on the Organic Analysis Data Sheets (Form I) included as Attachment A to this report. Region I qualifier definitions are also provided in this attachment.

At the discretion of the data user, the following points may warrant consideration by laboratory and field personnel prior to future sampling efforts:

- Documentation of sample preservation in the field on all chain of custody records.
- Documentation of complete cooler temperatures and sample pHs on receipt at the laboratory directly on the chain of custody records or, as a minimum, documentation of cooler temperatures directly on the applicable chain of custody records and inclusion of copies of run logs or other laboratory documentation containing sample pH measurements at the time of analysis.
- Use of linear (or higher order) regressions for calibration of target analytes for which the %RSD from initial calibration exceeds 15%, in accordance with the requirements of the referenced method.
- Use of reporting limits for non-detected analytes that reflect, at best, the lowest concentration initial calibration standard run for each target analyte.
- Documentation of applicable method detection limits (MDLs) in the data package.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation, Sample Integrity: Copies of the two applicable chain of custody (COC) records were included in the data package, documenting sample collection dates of 7/12/00 and 7/13/00. An acceptable temperature (4°C) on laboratory receipt was documented on the COC for the samples collected on 7/12/00. On the COC for samples collected 7/13/00, "7°" was documented by the laboratory. It was assumed that this was also a centigrade measurement. Although slightly above the generally accepted temperature range of 4°C±2°C, the exceedance is very minor and no action was taken on this basis.

Acidification of the samples for volatiles analysis with hydrochloric acid was recorded by the sampler on only one of the COCs and documentation of sample pHs on laboratory receipt or at the time of

Mr. Dan Garrigan

Evaluation of STL Report No. 7000-1470A (Volatiles in GW)

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analysis was not included in the data package. At the request of the evaluator, the laboratory provided their sample receiving preservative records for all samples. However, these records contained pH measurements for only those sample containers preserved with sulfuric acid or nitric acid; they offered no information regarding verification of preservation for the sample containers intended for volatile organics analysis. The laboratory was contacted again by the evaluator, on 9/5/00. On 9/6/00, copies of the laboratory's "GC/MS Volatiles Injection Logs" were received from STL/Connecticut via facsimile. Sample pHs were documented on these records and were verified to be <2 in all cases. No further action was taken with regard to this issue.

All volatiles analyses were performed on 7/21/00 and 7/24/00. Both dates are within the required 14-day holding time for chemically preserved water samples.

GC/MS Instrument Performance Check: Summary forms (Form V) were included for four bromofluorobenzene (BFB) instrument performance checks on instrument "HP5971L," reflecting each 12-hour period during which samples, associated standards, and associated quality control samples were analyzed. Reported relative abundances for all four performance checks were acceptable.

Calibration: Summary results (Form VI) for an initial calibration (IC) performed on instrument HP5971L on 7/20/00 were reported in support of the sample analyses. Results for five standards were reported, at concentrations of 0.5, 2, 10, 20, and 30 $\mu\text{g/L}$ for most target analytes, establishing a calibration range of 0.5 $\mu\text{g/L}$ to 30 $\mu\text{g/L}$ for most target analytes. There was no indication in the data package that any target analytes were included in the standards at higher concentrations, however the laboratory was asked for clarification of the individual analyte concentrations and responded (memorandum dated 9/1/00) that the ketones (acetone, 2-butanone, 4-methyl-2-pentanone, and 2-hexanone) were present at five times the concentration of most target analytes (i.e., 2.5 $\mu\text{g/L}$ to 150 $\mu\text{g/L}$ and that methylene chloride was calibrated at concentrations twice those of most target analytes (i.e., 1 $\mu\text{g/L}$ to 60 $\mu\text{g/L}$).

For the relevant target analytes, reported average relative response factors (RRFs) were acceptable (greater than 0.05) in the IC except for acetone (0.019) and 2-butanone (0.037). Results for acetone in MW-201, MW-202, MW-103A, CW-7A, and CW-5B and for 2-butanone in all 10 samples were rejected (R) as unreliable on this basis. Acetone was detected in TBL+RR071200, TBL+RR071300, MW-104A, MW-102A, and DUP-L+RR, therefore these results were qualified as estimated (J) based on the low responses in the associated calibration standards.

Percent relative standard deviations (%RSDs) were below the method-specified maximum acceptance criterion of 15% for all target analytes except bromomethane (21.8%), acetone (21.7%), methylene chloride (34.6%), 2-butanone (18.5%), bromochloromethane (26.1%), tetrachloroethene (25.4%),

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Evaluation of STL Report No. 7000-1470A (Volatiles in GW)

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1,2-dibromo-3-chloropropane (25.6%), and 1,2-dichloroethane- d_4 (24.1%). No evidence that linear (or higher order) regressions were performed was provided in the data package. No positive results were reported for bromomethane, bromochloromethane, or 1,2-dibromo-3-chloropropane in any of the samples and the %RSDs for these compounds did not exceed the less restrictive validation criterion of 30%, therefore no action was taken with respect to the results for these analytes. Results for 2-butanone in all samples were previously rejected and results for acetone in all samples were rejected or estimated based on unacceptably low average RRFs; no further action was warranted with respect to these analytes. 1,2-Dichloroethane- d_4 is a system monitoring compound and recoveries were not adversely affected in any of the samples, therefore no action was warranted based on the high %RSD for this compound. Results for methylene chloride in all samples and for tetrachloroethene in CW-5B, MW-102A, and DUP-L+RR were qualified as estimated (J) based on the high %RSDs; tetrachloroethene was not detected in any other samples and the %RSD did not exceed the less restrictive validation criterion of 30%, therefore no further action was warranted.

Form VII-equivalents were provided for three continuing calibration (CC) standards, run on 7/21/00 at 08:56, on 7/21/00 at 22:31, and on 7/24/00 at 07:56. Reported RRFs were acceptable (greater than 0.05) except for acetone (0.016, 0.016, and 0.018) and 2-butanone (0.034, 0.035, and 0.032) in all three CCs. All sample results for acetone and 2-butanone were previously qualified based on similar responses in the ICs and no additional action was taken based on the CC results.

Reported percent differences (%Ds) from the IC were less than the method-specified acceptance criterion (20%) in each of the CCs with the following exceptions:

| | |
|----------------|-------------------------------------|
| 7/21/00-08:56: | 1,2-dibromo-3-chloropropane - 22.4% |
| 7/21/00-22:31: | 2,2-dichloropropane - 21.0% |
| 7/24/00-07:56: | 1,2-dibromo-3-chloropropane - 20.8% |

All of the %Ds listed above are below the maximum acceptance limit specified by the validation guidelines (25%). Therefore, no sample results were qualified based on these responses.

Blanks: Results for three method blanks (MBs) associated with the site sample analyses were provided in the data package. Methylene chloride was detected in all three MBs, at concentrations ranging from 1.1 $\mu\text{g/L}$ to 1.4 $\mu\text{g/L}$. In addition, acetone (6.1 $\mu\text{g/L}$) and trichlorofluoromethane (0.14 $\mu\text{g/L}$) were reported in the MB run on 7/24/00. Results for methylene chloride in all samples, for acetone in MW-102A, MW-104A, and DUP-L+RR, and for trichlorofluoromethane in MW-102A and DUP-L+RR were qualified as less than the reported values or less than the sample-specific reporting limit (RL), whichever was greater, based on the associated MB contamination (U). In all

Mr. Dan Garrigan

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cases, the qualified sample results were less than ten times (methylene chloride and acetone) or five times (trichlorofluoromethane) the concentration found in the associated MB. All comparisons were made based on sample results prior to adjustment for the applicable dilution factor.

Two trip blanks (TBs) were submitted with this data set. After qualifications based on MB contamination, chloroform (0.72 µg/L), toluene (0.1 µg/L), and acetone (12 µg/L) were reported in TBL+RR071200 and chloroform (0.72 µg/L) and acetone (4.7 µg/L) were reported in TBL+RR071300. Results for acetone in MW-102A, MW-104A, and DUP-L+RR warranted qualification as less than the reported values or less than the sample-specific RL, whichever was greater based on the TB results; however, these results were similarly qualified based on associated MB contamination and no additional action was taken. Chloroform was not detected in any of the site samples and toluene was not detected in any of the samples associated with TBL+RR071200, therefore no further action was warranted on this basis.

Surrogates: Reported recoveries (Form II) were acceptable for all site samples and quality control (QC) analyses.

Matrix Spike/Matrix Spike Duplicate (MS/MSD): Percent recovery (%R) and relative percent difference (RPD) results for MS/MSD analyses performed on sample CW-5B were reported on Form III. Percent recoveries (100-106%) for the five reported spiked target analytes (1,1-dichloroethene, trichloroethene, benzene, toluene, and chlorobenzene at 10 µg/L each) were within the specified acceptance limits in both spiked analyses. In addition, relative percent differences (RPDs) were below the laboratory-specified maximum acceptance limits for all five reported spiked analytes.

Based on the Form I results provided for the MS and the MSD analyses, it is apparent that most, if not all, of the target analytes were actually included in the spiking solution. However, true values were provided only for the five compounds listed above and no further evaluation of these data could be made.

Blank Spikes: Recoveries for 36 analytes in a blank spike analysis performed in association with the site sample analyses were also reported in the data package. The 36 analytes were included in the spiking solution at 10 µg/L or 50 µg/L. Percent recoveries were within the laboratory-specified acceptance limits (QC 60-140%) in all cases except carbon disulfide (0%), methylene chloride (200%), 2-butanone (40%), 4-methyl-2-pentanone (42%), 2-hexanone (42%), and m/p-xylenes (205%). Carbon disulfide and 2-hexanone are not project-specific target analytes, therefore no action was warranted with respect to these two analytes. Results for methylene chloride in all samples and for m/p-xylenes in CW-5B and MW-104A were qualified as estimated (J) based on the unacceptably high recoveries; these results may be biased high. Results for 2-butanone and 4-methyl-2-pentanone

Mr. Dan Garrigan

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in all samples were qualified as estimated (UJ) based on the unacceptably low recoveries in the blank spike.

Field Duplicate: Sample DUP-L+RR was identified as a field duplicate of MW-102A based on information provided by the client. Positive paired results for benzene (0 RPD), chlorobenzene (0 RPD), chloroethane (3.6 RPD), 1,2-dichlorobenzene (7.3 RPD), 1,4-dichlorobenzene (3.6 RPD), dichlorodifluoromethane (0 RPD), 1,1-dichloroethane (0 RPD), 1,2-dichloroethane (6.5 RPD), cis-1,2-dichloroethene (4.3 RPD), trans-1,2-dichloroethene (5.4 RPD), 1,2-dichloropropane (3.1 RPD), isopropylbenzene (2.4 RPD), tetrachloroethene (1.8 RPD), toluene (6.5 RPD), trichloroethene (2.6 RPD), and vinyl chloride (2.2 RPD) showed excellent reproducibility.

Internal Standard Responses: Internal standard areas and retention times (RTs) were reported on Form VIII for all site samples reported in this data set and the related quality control analyses; all areas and RTs were within acceptable limits as reported on the forms.

Sample Results: Results for all GW samples and both TBs were reported on Form Is, all of which indicated a sample analysis volume of 25 mL. Samples MW-102A, MW-104A, and DUPL+RR were run at 1:10 dilutions, and only the diluted analysis results were reported. Reporting limits for target analytes that were not detected were appropriately adjusted to reflect the dilution factor in these three sample analyses.

Target analytes that were not detected in the samples were reported by the laboratory as less than 0.5 µg/L (0.5 U) except for acetone (12 U), 2-butanone (6 U), 4-methyl-2-pentanone (5 U), 1,2-dibromo-3-chloropropane (2 U), vinyl chloride (0.2 U), and methylene chloride (0.3 U).

No MDL study results were provided in the data package. With the exceptions of vinyl chloride and methylene chloride, all RLs reported by the laboratory were equivalent to or above the lowest concentration IC standard associated with these analyses and no positive sample results less than the RLs were reported. Therefore, the reported results for all target analytes except vinyl chloride and methylene chloride are fully supported by the available data and are acceptable as reported.

Vinyl chloride was present at 0.5 µg/L in the lowest concentration calibration standard and methylene chloride was present in this standard at 1.0 µg/L. In the absence of MDL data for these two analytes, the RLs reported by the laboratory for these two analytes are not supportable. Therefore, RLs for vinyl chloride and methylene chloride were corrected to 0.5 µg/L and 1 µg/L, respectively, to reflect the lowest concentration calibration standards run in association with these analyses.

Documentation: Data package documentation was consistent with the specifications of the QAPP; specifically, a Level III data package (summary forms only - no supporting raw data) was provided for review.

Mr. Dan Garrigan

Evaluation of STL Report No. 7000-1470A (Volatiles in GW)

September 11, 2000

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COC records provided in the data package included all reported samples and were properly completed except that sample preservation was recorded on only one of the COCs and improper corrections were noted on both documents. In the future, all corrections should be made by drawing a single line through the incorrect information, inserting the correct information, and initialing and dating the change. "Write-overs" are not legally defensible.

Documentation of sample pHs on laboratory receipt (or at the time of analysis) was not provided in the data package as received for review. This information was verbally requested by the validator on 8/30/00; sample receiving preservation logs were received from the laboratory via facsimile on 9/2/00. These records were inserted into the data package by the evaluator as pages 55a and 55b, but did not address the samples intended for volatile organics analyses. The laboratory was contacted by the evaluator again on 9/5/00, and copies of the laboratory's GC/MS injection logs were provided via facsimile on 9/6/00. These logs are included in Attachment B.

Sample identifications for both trip blanks were shortened by the laboratory throughout the data package by removing "L+RR" from the identifications found on the COC records. The chain of custody identifications are used throughout this evaluation report for all samples, and the Form Is for the trip blanks were corrected by the evaluator to reflect the COC identifications.

On Form II, the identification of SMC2 at the bottom of the page was corrected to dibromofluoromethane (from 1,2-dichloroethane-d4) by the evaluator. A copy of the corrected form is included in Attachment B to this report.

On Form III, a result of 0.03 µg/L was reported for toluene in the unspiked analysis of CW-5B. On the Form I for this sample, toluene is reported as not detected (0.5 U). Therefore, the Form III result is incorrect. However, it is so low that it does not affect the calculated recovery, and no action was taken by the evaluator.

Form VII was not included in the data package for two of the three CCs associated with the sample analyses. The missing forms were requested from the laboratory on 8/30/00, and were received via facsimile on 9/1/00. They were inserted into the data package by the evaluator as pages 74a-74f.

Please let me know if you have any questions regarding this data package review report.

Sincerely,



Carol A. Erikson

Quality Assessment Manager

CAE/ekd

95406\Apr00\GWVoa

ATTACHMENT A

**ORGANIC ANALYSIS DATA SHEETS (Form I)
STL/Connecticut Case No. 1470A**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected. The associated numerical value is the sample quantitation limit. The sample quantitation limit accounts for sample-specific dilution factors and percent solids corrections or sample sizes that deviate from those required by the method.**
- J - The associated numerical value is an estimated quantity.**
- R - The data are unusable (analyte may or may not be present). Resampling and reanalysis is necessary for verification. The R replaces the numerical value or sample quantitation limit.**
- UJ - The analyte was analyzed for, but was not detected. The sample quantitation limit is an estimated quantity.**

VOLATILE ORGANICS ANALYSIS DATA SHEET

MW-201

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____ SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-01

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2009

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____(uL)

Soil Aliquot Volume: _____(uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|-----------------------------|--|---|
| 71-43-2 | Benzene | .5 | U |
| 108-86-1 | Bromobenzene | .5 | U |
| 74-97-5 | Bromochloromethane | .5 | U |
| 75-27-4 | Bromodichloromethane | .5 | U |
| 75-25-2 | Bromoform | .5 | U |
| 74-83-9 | Bromomethane | .5 | U |
| 104-51-8 | n-Butylbenzene | .5 | U |
| 135-98-8 | Sec-Butylbenzene | .5 | U |
| 98-06-6 | Tert-Butylbenzene | .5 | U |
| 56-23-5 | Carbon Tetrachloride | .5 | U |
| 108-90-7 | Chlorobenzene | .5 | U |
| 124-48-1 | Dibromochloromethane | .5 | U |
| 75-00-3 | Chloroethane | .5 | U |
| 67-66-3 | Chloroform | .5 | U |
| 74-87-3 | Chloromethane | .5 | U |
| 95-49-8 | 2-Chlorotoluene | .5 | U |
| 106-43-4 | 4-Chlorotoluene | .5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2 | U |
| 106-93-4 | 1,2-Dibromoethane | .5 | U |
| 74-95-3 | Dibromomethane | .5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | .5 | U |
| 541-73-1 | 1,3-Dichlorobenzene | .5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | .5 | U |
| 75-71-8 | Dichlorodifluoromethane | 4.6 | U |
| 75-34-3 | 1,1-Dichloroethane | .5 | U |
| 107-06-2 | 1,2-Dichloroethane | .5 | U |
| 75-35-4 | 1,1-Dichloroethene | .5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | .5 | U |
| 156-60-5 | trans-1,2-Dichloroethene | .5 | U |
| 78-87-5 | 1,2-Dichloropropane | .5 | U |
| 142-28-9 | 1,3-Dichloropropane | .5 | U |
| 544-20-7 | 2,2-Dichloropropane | .5 | U |
| 563-58-6 | 1,1-Dichloropropene | .5 | U |
| 100-41-4 | Ethylbenzene | .5 | U |
| 87-68-3 | Hexachlorobutadiene | .5 | U |

0005

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

MW-201

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water) WATER

Lab Sample ID: 001470A-01

Sample wt/vol: 25 (g/mL) ML

Lab File ID: >L2009

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|---------------------------|--|---|
| 98-82-8 | Isopropylbenzene | .5 | U |
| 99-87-6 | 4-Isopropyltoluene | .5 | U |
| 75-09-2 | Methylene Chloride | 1.6 | U |
| 91-20-3 | Naphthalene | .5 | U |
| 103-65-1 | N-Propylbenzene | .5 | U |
| 100-42-5 | Styrene | .5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | .5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | .5 | U |
| 127-18-4 | Tetrachloroethene | .5 | U |
| 108-88-3 | Toluene | .5 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | .5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | .5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | .5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | .5 | U |
| 79-01-6 | Trichloroethene | .5 | U |
| 75-69-4 | Trichlorofluoromethane | .23 | J |
| 96-18-4 | 1,2,3-Trichloropropane | .5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | .5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | .5 | U |
| 75-01-4 | Vinyl Chloride | 0.5 | U |
| NO CAS | mp-Xylene | 1 | U |
| 95-47-6 | o-xylene | .5 | U |
| 67-64-1 | Acetone | 12 | U |
| 78-93-3 | 2-Butanone | 6 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 5 | U |

Case 9/8/00

CASEYKSON 9/8/00

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

0004

CLIENT ID

MW-202

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-02

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2010

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____(uL)

Soil Aliquot Volume: _____(uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|-----------------------------|--|---|
| 71-43-2 | Benzene | .5 | U |
| 108-86-1 | Bromobenzene | .5 | U |
| 74-97-5 | Bromochloromethane | .5 | U |
| 75-27-4 | Bromodichloromethane | .5 | U |
| 75-25-2 | Bromoform | .5 | U |
| 74-83-9 | Bromomethane | .5 | U |
| 104-51-8 | n-Butylbenzene | .5 | U |
| 135-98-8 | Sec-Butylbenzene | .5 | U |
| 98-06-6 | Tert-Butylbenzene | .5 | U |
| 56-23-5 | Carbon Tetrachloride | .5 | U |
| 108-90-7 | Chlorobenzene | .5 | U |
| 124-48-1 | Dibromochloromethane | .5 | U |
| 75-00-3 | Chloroethane | .5 | U |
| 67-66-3 | Chloroform | .5 | U |
| 74-87-3 | Chloromethane | .5 | U |
| 95-49-8 | 2-Chlorotoluene | .5 | U |
| 106-43-4 | 4-Chlorotoluene | .5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | .2 | U |
| 106-93-4 | 1,2-Dibromoethane | .5 | U |
| 74-95-3 | Dibromomethane | .5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | .5 | U |
| 541-73-1 | 1,3-Dichlorobenzene | .5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | .5 | U |
| 75-71-8 | Dichlorodifluoromethane | .5 | U |
| 75-34-3 | 1,1-Dichloroethane | .5 | U |
| 107-06-2 | 1,2-Dichloroethane | .5 | U |
| 75-35-4 | 1,1-Dichloroethene | .5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | .5 | U |
| 156-60-5 | trans-1,2-Dichloroethene | .5 | U |
| 78-87-5 | 1,2-Dichloropropane | .5 | U |
| 142-28-9 | 1,3-Dichloropropane | .5 | U |
| 544-20-7 | 2,2-Dichloropropane | .5 | U |
| 563-58-6 | 1,1-Dichloropropene | .5 | U |
| 100-41-4 | Ethylbenzene | .5 | U |
| 87-68-3 | Hexachlorobutadiene | .5 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

0005

CLIENT ID

MW-202

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____ SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-02

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2010

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION | Q |
|----------|---------------------------|---------------|---|
| 98-82-8 | Isopropylbenzene | .5 | U |
| 99-87-6 | 4-Isopropyltoluene | .5 | U |
| 75-09-2 | Methylene Chloride | 1 .5 | U |
| 91-20-3 | Naphthalene | .5 | U |
| 103-65-1 | N-Propylbenzene | .5 | U |
| 100-42-5 | Styrene | .5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | .5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | .5 | U |
| 127-18-4 | Tetrachloroethene | .5 | U |
| 108-88-3 | Toluene | .5 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | .5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | .5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | .5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | .5 | U |
| 79-01-6 | Trichloroethene | .5 | U |
| 75-69-4 | Trichlorofluoromethane | .5 | U |
| 96-18-4 | 1,2,3-Trichloropropane | .5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | .5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | .5 | U |
| 75-01-4 | Vinyl Chloride | 0.5 .2 | U |
| NO CAS | mp-Xylene | 1 | U |
| 95-47-6 | o-xylene | .5 | U |
| 67-64-1 | Acetone | R 12 | U |
| 78-93-3 | 2-Butanone | R -6 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 5 | U |

CAE
9/9/00

CAE
9/8/00

CAE
9/8/00

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

0006

CLIENT ID

MW-103A

Lab Name: STL/CT

Contract: _____

Lab Code: IACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-03

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2011

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg)UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg)UG/L | Q |
|----------|-----------------------------|---|---|
| 71-43-2 | Benzene | .5 | U |
| 108-86-1 | Bromobenzene | .5 | U |
| 74-97-5 | Bromochloromethane | .5 | U |
| 75-27-4 | Bromodichloromethane | .5 | U |
| 75-25-2 | Bromoform | .5 | U |
| 74-83-9 | Bromomethane | .5 | U |
| 104-51-8 | n-Butylbenzene | .5 | U |
| 135-98-8 | Sec-Butylbenzene | .5 | U |
| 98-06-6 | Tert-Butylbenzene | .5 | U |
| 56-23-5 | Carbon Tetrachloride | .5 | U |
| 108-90-7 | Chlorobenzene | .5 | U |
| 124-48-1 | Dibromochloromethane | .5 | U |
| 75-00-3 | Chloroethane | .5 | U |
| 67-66-3 | Chloroform | .5 | U |
| 74-87-3 | Chloromethane | .5 | U |
| 95-49-8 | 2-Chlorotoluene | .5 | U |
| 106-43-4 | 4-Chlorotoluene | .5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2 | U |
| 106-93-4 | 1,2-Dibromoethane | .5 | U |
| 74-95-3 | Dibromomethane | .5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | .5 | U |
| 541-73-1 | 1,3-Dichlorobenzene | .5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | .5 | U |
| 75-71-8 | Dichlorodifluoromethane | .5 | U |
| 75-34-3 | 1,1-Dichloroethane | .1 | J |
| 107-06-2 | 1,2-Dichloroethane | .5 | U |
| 75-35-4 | 1,1-Dichloroethene | .5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | .5 | U |
| 156-60-5 | trans-1,2-Dichloroethene | .5 | U |
| 78-87-5 | 1,2-Dichloropropane | .5 | U |
| 142-28-9 | 1,3-Dichloropropane | .5 | U |
| 544-20-7 | 2,2-Dichloropropane | .5 | U |
| 563-58-6 | 1,1-Dichloropropene | .5 | U |
| 100-41-4 | Ethylbenzene | .5 | U |
| 87-68-3 | Hexachlorobutadiene | .5 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

MW-103A

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-03

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2011

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg)UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg)UG/L | Q |
|----------|---------------------------|---|------|
| 98-82-8 | Isopropylbenzene | .5 | U |
| 99-87-6 | 4-Isopropyltoluene | .5 | U |
| 75-09-2 | Methylene Chloride | 1 .38 | JB U |
| 91-20-3 | Naphthalene | .5 | U |
| 103-65-1 | N-Propylbenzene | .5 | U |
| 100-42-5 | Styrene | .5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | .5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | .5 | U |
| 127-18-4 | Tetrachloroethene | .5 | U |
| 108-88-3 | Toluene | .5 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | .5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | .5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | .5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | .5 | U |
| 79-01-6 | Trichloroethene | .5 | U |
| 75-69-4 | Trichlorofluoromethane | .5 | U |
| 96-18-4 | 1,2,3-Trichloropropane | .5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | .5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | .5 | U |
| 75-01-4 | Vinyl Chloride | 0.5 .2 | U |
| NO CAS | mp-Xylene | 1 | U |
| 95-47-6 | o-xylene | .5 | U |
| 67-64-1 | Acetone | R .12 | U |
| 78-93-3 | 2-Butanone | R .6 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 5 | U |

CAE
9/8/00CAE
9/8/00CAE
9/8/00

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

CW-7A

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-04

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2012

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|-----------------------------|--|---|
| 71-43-2 | Benzene | .26 | J |
| 108-86-1 | Bromobenzene | .5 | U |
| 74-97-5 | Bromochloromethane | .5 | U |
| 75-27-4 | Bromodichloromethane | .5 | U |
| 75-25-2 | Bromoform | .5 | U |
| 74-83-9 | Bromomethane | .5 | U |
| 104-51-8 | n-Butylbenzene | .5 | U |
| 135-98-8 | Sec-Butylbenzene | .5 | U |
| 98-06-6 | Tert-Butylbenzene | .5 | U |
| 56-23-5 | Carbon Tetrachloride | .5 | U |
| 108-90-7 | Chlorobenzene | .75 | |
| 124-48-1 | Dibromochloromethane | .5 | U |
| 75-00-3 | Chloroethane | .15 | J |
| 67-66-3 | Chloroform | .5 | U |
| 74-87-3 | Chloromethane | .5 | U |
| 95-49-8 | 2-Chlorotoluene | .5 | U |
| 106-43-4 | 4-Chlorotoluene | .5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | .2 | U |
| 106-93-4 | 1,2-Dibromoethane | .5 | U |
| 74-95-3 | Dibromomethane | .5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | .5 | U |
| 541-73-1 | 1,3-Dichlorobenzene | .5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | .83 | |
| 75-71-8 | Dichlorodifluoromethane | .5 | U |
| 75-34-3 | 1,1-Dichloroethane | .35 | J |
| 107-06-2 | 1,2-Dichloroethane | .5 | U |
| 75-35-4 | 1,1-Dichloroethene | .5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | .12 | J |
| 156-60-5 | trans-1,2-Dichloroethene | .5 | U |
| 78-87-5 | 1,2-Dichloropropane | .5 | U |
| 142-28-9 | 1,3-Dichloropropane | .5 | U |
| 544-20-7 | 2,2-Dichloropropane | .5 | U |
| 563-58-6 | 1,1-Dichloropropene | .5 | U |
| 100-41-4 | Ethylbenzene | .5 | U |
| 87-68-3 | Hexachlorobutadiene | .5 | U |

VOLATILE ORGANICS ANALYSIS DATA SHEET

CW-7A

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____ SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-04

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2012

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|---------------------------|--|-----------------|
| 98-82-8 | Isopropylbenzene | .24 | J |
| 99-87-6 | 4-Isopropyltoluene | .5 | U |
| 75-09-2 | Methylene Chloride | 1 -44 | JB U |
| 91-20-3 | Naphthalene | .5 | U |
| 103-65-1 | N-Propylbenzene | .5 | U |
| 100-42-5 | Styrene | .5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | .5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | .5 | U |
| 127-18-4 | Tetrachloroethene | .5 | U |
| 108-88-3 | Toluene | .5 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | .5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | .5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | .5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | .5 | U |
| 79-01-6 | Trichloroethene | .5 | U |
| 75-69-4 | Trichlorofluoromethane | .5 | U |
| 96-18-4 | 1,2,3-Trichloropropane | .5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | .5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | .5 | U |
| 75-01-4 | Vinyl Chloride | 0.5 -2 | U |
| NO CAS | mp-Xylene | 1 | U |
| 95-47-6 | o-xylene | .5 | U |
| 67-64-1 | Acetone | R 12 | U |
| 78-93-3 | 2-Butanone | R 6 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 5 | UJ |

CAE
9/8/00CAE
9/8/00CAE
9/8/00

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

0010

CLIENT ID

TBLR071200
TB071200

Lab Name: STL/CT

Contract: _____

06/18/00

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-05

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2008

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____(uL)

Soil Aliquot Volume: _____(uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|-----------------------------|--|---|
| 71-43-2 | Benzene | .5 | U |
| 108-86-1 | Bromobenzene | .5 | U |
| 74-97-5 | Bromochloromethane | .5 | U |
| 75-27-4 | Bromodichloromethane | .5 | U |
| 75-25-2 | Bromoform | .5 | U |
| 74-83-9 | Bromomethane | .5 | U |
| 104-51-8 | n-Butylbenzene | .5 | U |
| 135-98-8 | Sec-Butylbenzene | .5 | U |
| 98-06-6 | Tert-Butylbenzene | .5 | U |
| 56-23-5 | Carbon Tetrachloride | .5 | U |
| 108-90-7 | Chlorobenzene | .5 | U |
| 124-48-1 | Dibromochloromethane | .5 | U |
| 75-00-3 | Chloroethane | .5 | U |
| 67-66-3 | Chloroform | .72 | |
| 74-87-3 | Chloromethane | .5 | U |
| 95-49-8 | 2-Chlorotoluene | .5 | U |
| 106-43-4 | 4-Chlorotoluene | .5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2 | U |
| 106-93-4 | 1,2-Dibromoethane | .5 | U |
| 74-95-3 | Dibromomethane | .5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | .5 | U |
| 541-73-1 | 1,3-Dichlorobenzene | .5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | .5 | U |
| 75-71-8 | Dichlorodifluoromethane | .5 | U |
| 75-34-3 | 1,1-Dichloroethane | .5 | U |
| 107-06-2 | 1,2-Dichloroethane | .5 | U |
| 75-35-4 | 1,1-Dichloroethene | .5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | .5 | U |
| 156-60-5 | trans-1,2-Dichloroethene | .5 | U |
| 78-87-5 | 1,2-Dichloropropane | .5 | U |
| 142-28-9 | 1,3-Dichloropropane | .5 | U |
| 544-20-7 | 2,2-Dichloropropane | .5 | U |
| 563-58-6 | 1,1-Dichloropropene | .5 | U |
| 100-41-4 | Ethylbenzene | .5 | U |
| 87-68-3 | Hexachlorobutadiene | .5 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

Lab Name: STL/CT

Contract: _____

TBL+R2071200
TB071200

CAE 9/8/00

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-05

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2008

Level: (low/med) LOW

Date Received: 07/13/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____(uL)

Soil Aliquot Volume: _____(uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg)UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg)UG/L | Q |
|----------|---------------------------|---|-----|
| 98-82-8 | Isopropylbenzene | .5 | U |
| 99-87-6 | 4-Isopropyltoluene | .5 | U |
| 75-09-2 | Methylene Chloride | 3.4 | BU |
| 91-20-3 | Naphthalene | .5 | U |
| 103-65-1 | N-Propylbenzene | .5 | U |
| 100-42-5 | Styrene | .5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | .5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | .5 | U |
| 127-18-4 | Tetrachloroethene | .5 | U |
| 108-88-3 | Toluene | .1 | J |
| 87-61-6 | 1,2,3-Trichlorobenzene | .5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | .5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | .5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | .5 | U |
| 79-01-6 | Trichloroethene | .5 | U |
| 75-69-4 | Trichlorofluoromethane | .5 | U |
| 96-18-4 | 1,2,3-Trichloropropane | .5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | .5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | .5 | U |
| 75-01-4 | Vinyl Chloride | 0.5 | U |
| NO CAS | mp-Xylene | 1 | U |
| 95-47-6 | o-xylene | .5 | U |
| 67-64-1 | Acetone | 12 | J |
| 78-93-3 | 2-Butanone | R 6 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 5 | BUJ |

CAE 9/8/00

CAE 9/8/00

CAE 9/8/00

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

| |
|-------|
| CW-5B |
|-------|

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-06

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2006

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.

COMPOUND

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|-----------------------------|--|---|
| 71-43-2 | Benzene | .55 | |
| 108-86-1 | Bromobenzene | .5 | U |
| 74-97-5 | Bromochloromethane | .5 | U |
| 75-27-4 | Bromodichloromethane | .5 | U |
| 75-25-2 | Bromoform | .5 | U |
| 74-83-9 | Bromomethane | .5 | U |
| 104-51-8 | n-Butylbenzene | .5 | U |
| 135-98-8 | Sec-Butylbenzene | .5 | U |
| 98-06-6 | Tert-Butylbenzene | .5 | U |
| 56-23-5 | Carbon Tetrachloride | .5 | U |
| 108-90-7 | Chlorobenzene | .5 | U |
| 124-48-1 | Dibromochloromethane | .5 | U |
| 75-00-3 | Chloroethane | 1.8 | |
| 67-66-3 | Chloroform | .5 | U |
| 74-87-3 | Chloromethane | .5 | U |
| 95-49-8 | 2-Chlorotoluene | .5 | U |
| 106-43-4 | 4-Chlorotoluene | .5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2 | U |
| 106-93-4 | 1,2-Dibromoethane | .5 | U |
| 74-95-3 | Dibromomethane | .5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | .5 | U |
| 541-73-1 | 1,3-Dichlorobenzene | .5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | .5 | U |
| 75-71-8 | Dichlorodifluoromethane | 8.7 | |
| 75-34-3 | 1,1-Dichloroethane | 19 | |
| 107-06-2 | 1,2-Dichloroethane | .5 | U |
| 75-35-4 | 1,1-Dichloroethene | .35 | J |
| 156-59-2 | cis-1,2-Dichloroethene | 11 | |
| 156-60-5 | trans-1,2-Dichloroethene | .44 | J |
| 78-87-5 | 1,2-Dichloropropane | .5 | U |
| 142-28-9 | 1,3-Dichloropropane | .5 | U |
| 544-20-7 | 2,2-Dichloropropane | .5 | U |
| 563-58-6 | 1,1-Dichloropropene | .5 | U |
| 100-41-4 | Ethylbenzene | .12 | J |
| 87-68-3 | Hexachlorobutadiene | .5 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

CW-5B

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-06

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2006

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____(uL)

Soil Aliquot Volume: _____(uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|---------------------------|--|-----------------|
| 98-82-8 | Isopropylbenzene | .11 | J |
| 99-87-6 | 4-Isopropyltoluene | .15 | J |
| 75-09-2 | Methylene Chloride | 3.8 | U UJ |
| 91-20-3 | Naphthalene | .5 | U |
| 103-65-1 | N-Propylbenzene | .5 | U |
| 100-42-5 | Styrene | .5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | .5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | .5 | U |
| 127-18-4 | Tetrachloroethene | 24 | J |
| 108-88-3 | Toluene | .5 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | .5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | .5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 1 | U |
| 79-00-5 | 1,1,2-Trichloroethane | .5 | U |
| 79-01-6 | Trichloroethene | 9.4 | U |
| 75-69-4 | Trichlorofluoromethane | .5 | U |
| 96-18-4 | 1,2,3-Trichloropropane | .5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | .5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | .5 | U |
| 75-01-4 | Vinyl Chloride | 6.8 | U |
| NO CAS | mp-Xylene | 3.7 | J |
| 95-47-6 | o-xylene | .32 | J |
| 67-64-1 | Acetone | R 12 | U |
| 78-93-3 | 2-Butanone | R 6 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 5 | U UJ |

9/8/00
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VOLATILE ORGANICS ANALYSIS DATA SHEET

Lab Name: STL/CT

Contract: _____

MW-104A

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water) WATER

Lab Sample ID: 001470A-07

Sample wt/vol: 25 (g/mL) ML

Lab File ID: >L2032

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/24/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 10.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO.

COMPOUND

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|-----------------------------|--|---|
| 71-43-2 | Benzene | 24 | |
| 108-86-1 | Bromobenzene | 5 | U |
| 74-97-5 | Bromochloromethane | 5 | U |
| 75-27-4 | Bromodichloromethane | 5 | U |
| 75-25-2 | Bromoform | 5 | U |
| 74-83-9 | Bromomethane | 5 | U |
| 104-51-8 | n-Butylbenzene | 5 | U |
| 135-98-8 | Sec-Butylbenzene | 5 | U |
| 98-06-6 | Tert-Butylbenzene | 5 | U |
| 56-23-5 | Carbon Tetrachloride | 5 | U |
| 108-90-7 | Chlorobenzene | 10 | |
| 124-48-1 | Dibromochloromethane | 5 | U |
| 75-00-3 | Chloroethane | 49 | |
| 67-66-3 | Chloroform | 5 | U |
| 74-87-3 | Chloromethane | 5 | U |
| 95-49-8 | 2-Chlorotoluene | 5 | U |
| 106-43-4 | 4-Chlorotoluene | 5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 20 | U |
| 106-93-4 | 1,2-Dibromoethane | 5 | U |
| 74-95-3 | Dibromomethane | 5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 6.6 | |
| 541-73-1 | 1,3-Dichlorobenzene | .88 | J |
| 106-46-7 | 1,4-Dichlorobenzene | 33 | |
| 75-71-8 | Dichlorodifluoromethane | 5 | U |
| 75-34-3 | 1,1-Dichloroethane | 40 | |
| 107-06-2 | 1,2-Dichloroethane | 9 | |
| 75-35-4 | 1,1-Dichloroethene | 5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | 5 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.8 | J |
| 78-87-5 | 1,2-Dichloropropane | 10 | |
| 142-28-9 | 1,3-Dichloropropane | 5 | U |
| 544-20-7 | 2,2-Dichloropropane | 5 | U |
| 563-58-6 | 1,1-Dichloropropene | 5 | U |
| 100-41-4 | Ethylbenzene | 140 | |
| 87-68-3 | Hexachlorobutadiene | 5 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

MW-104A

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-07

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2032

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/24/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 10.0

Soil Extract Volume: _____(uL)

Soil Aliquot Volume: _____(uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|---------------------------|--|----|
| 98-82-8 | Isopropylbenzene | 23 | |
| 99-87-6 | 4-Isopropyltoluene | 13 | |
| 75-09-2 | Methylene Chloride | 18 | UJ |
| 91-20-3 | Naphthalene | 59 | |
| 103-65-1 | N-Propylbenzene | 4.8 | J |
| 100-42-5 | Styrene | 2.1 | J |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | U |
| 127-18-4 | Tetrachloroethene | 5 | U |
| 108-88-3 | Toluene | 11 | |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | U |
| 79-01-6 | Trichloroethene | 5 | U |
| 75-69-4 | Trichlorofluoromethane | 5 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 42 | |
| 108-67-8 | 1,3,5-Trimethylbenzene | 12 | |
| 75-01-4 | Vinyl Chloride | 5.2 | U |
| NO CAS | mp-Xylene | 350 | J |
| 95-47-6 | o-xylene | 63 | |
| 67-64-1 | Acetone | 160 | UJ |
| 78-93-3 | 2-Butanone | R 60 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 50 | UJ |

9/8/00
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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

0016

CLIENT ID

MW-102A

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-08

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2030

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/24/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 10.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|-----------------------------|--|---|
| 71-43-2 | Benzene | 10 | |
| 108-86-1 | Bromobenzene | 5 | U |
| 74-97-5 | Bromochloromethane | 5 | U |
| 75-27-4 | Bromodichloromethane | 5 | U |
| 75-25-2 | Bromoform | 5 | U |
| 74-83-9 | Bromomethane | 5 | U |
| 104-51-8 | n-Butylbenzene | 5 | U |
| 135-98-8 | Sec-Butylbenzene | 5 | U |
| 98-06-6 | Tert-Butylbenzene | 5 | U |
| 56-23-5 | Carbon Tetrachloride | 5 | U |
| 108-90-7 | Chlorobenzene | 11 | |
| 124-48-1 | Dibromochloromethane | 5 | U |
| 75-00-3 | Chloroethane | 28 | |
| 67-66-3 | Chloroform | 5 | U |
| 74-87-3 | Chloromethane | 5 | U |
| 95-49-8 | 2-Chlorotoluene | 5 | U |
| 106-43-4 | 4-Chlorotoluene | 5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 20 | U |
| 106-93-4 | 1,2-Dibromoethane | 5 | U |
| 74-95-3 | Dibromomethane | 5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | .93 | J |
| 541-73-1 | 1,3-Dichlorobenzene | 5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 5.5 | |
| 75-71-8 | Dichlorodifluoromethane | 27 | |
| 75-34-3 | 1,1-Dichloroethane | 160 | |
| 107-06-2 | 1,2-Dichloroethane | 15 | |
| 75-35-4 | 1,1-Dichloroethene | 5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | 230 | |
| 156-60-5 | trans-1,2-Dichloroethene | 3.6 | J |
| 78-87-5 | 1,2-Dichloropropane | 6.4 | |
| 142-28-9 | 1,3-Dichloropropane | 5 | U |
| 544-20-7 | 2,2-Dichloropropane | 5 | U |
| 563-58-6 | 1,1-Dichloropropene | 5 | U |
| 100-41-4 | Ethylbenzene | 5 | U |
| 87-68-3 | Hexachlorobutadiene | 5 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

0017

CLIENT ID

MW-102A

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-08

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2030

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/24/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 10.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|---------------------------|--|-----------------|
| 98-82-8 | Isopropylbenzene | 4.1 | J |
| 99-87-6 | 4-Isopropyltoluene | 5 | U |
| 75-09-2 | Methylene Chloride | 32 | U UJ |
| 91-20-3 | Naphthalene | 5 | U |
| 103-65-1 | N-Propylbenzene | 5 | U |
| 100-42-5 | Styrene | 5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | U |
| 127-18-4 | Tetrachloroethene | 55 | J |
| 108-88-3 | Toluene | .95 | J |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | U |
| 79-01-6 | Trichloroethene | 77 | |
| 75-69-4 | Trichlorofluoromethane | 1.7 | U U |
| 96-18-4 | 1,2,3-Trichloropropane | 5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 5 | U |
| 75-01-4 | Vinyl Chloride | 46 | |
| NO CAS | mp-Xylene | 10 | U |
| 95-47-6 | o-xylene | 5 | U |
| 67-64-1 | Acetone | 100 | U UJ |
| 78-93-3 | 2-Butanone | R 60 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 50 | U UJ |

9/8/00
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VOLATILE ORGANICS ANALYSIS DATA SHEET

DUP-L+RR

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-09

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2031

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/24/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 10.0

Soil Extract Volume: _____(uL)

Soil Aliquot Volume: _____(uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg)UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg)UG/L | Q |
|----------|-----------------------------|---|---|
| 71-43-2 | Benzene | 10 | |
| 108-86-1 | Bromobenzene | 5 | U |
| 74-97-5 | Bromochloromethane | 5 | U |
| 75-27-4 | Bromodichloromethane | 5 | U |
| 75-25-2 | Bromoform | 5 | U |
| 74-83-9 | Bromomethane | 5 | U |
| 104-51-8 | n-Butylbenzene | 5 | U |
| 135-98-8 | Sec-Butylbenzene | 5 | U |
| 98-06-6 | Tert-Butylbenzene | 5 | U |
| 56-23-5 | Carbon Tetrachloride | 5 | U |
| 108-90-7 | Chlorobenzene | 11 | |
| 124-48-1 | Dibromochloromethane | 5 | U |
| 75-00-3 | Chloroethane | 27 | |
| 67-66-3 | Chloroform | 5 | U |
| 74-87-3 | Chloromethane | 5 | U |
| 95-49-8 | 2-Chlorotoluene | 5 | U |
| 106-43-4 | 4-Chlorotoluene | 5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 20 | U |
| 106-93-4 | 1,2-Dibromoethane | 5 | U |
| 74-95-3 | Dibromomethane | 5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 1 | J |
| 541-73-1 | 1,3-Dichlorobenzene | 5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 5.7 | |
| 75-71-8 | Dichlorodifluoromethane | 27 | |
| 75-34-3 | 1,1-Dichloroethane | 160 | |
| 107-06-2 | 1,2-Dichloroethane | 16 | |
| 75-35-4 | 1,1-Dichloroethene | 5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | 240 | |
| 156-60-5 | trans-1,2-Dichloroethene | 3.8 | J |
| 78-87-5 | 1,2-Dichloropropane | 6.6 | |
| 142-28-9 | 1,3-Dichloropropane | 5 | U |
| 544-20-7 | 2,2-Dichloropropane | 5 | U |
| 563-58-6 | 1,1-Dichloropropene | 5 | U |
| 100-41-4 | Ethylbenzene | 5 | U |
| 87-68-3 | Hexachlorobutadiene | 5 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

DUP-L+RR

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-09

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2031

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/24/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 10.0

Soil Extract Volume: _____(uL)

Soil Aliquot Volume: _____(uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|---------------------------|--|-----------------|
| 98-82-8 | Isopropylbenzene | 4.2 | J |
| 99-87-6 | 4-Isopropyltoluene | 5 | U |
| 75-09-2 | Methylene Chloride | 31 | U UJ |
| 91-20-3 | Naphthalene | 5 | U |
| 103-65-1 | N-Propylbenzene | 5 | U |
| 100-42-5 | Styrene | 5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5 | U |
| 127-18-4 | Tetrachloroethene | 54 | J |
| 108-88-3 | Toluene | .89 | J |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5 | U |
| 79-01-6 | Trichloroethene | 79 | U |
| 75-69-4 | Trichlorofluoromethane | 1.7 | U U |
| 96-18-4 | 1,2,3-Trichloropropane | 5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 5 | U |
| 75-01-4 | Vinyl Chloride | 45 | U |
| NO CAS | mp-Xylene | 10 | U |
| 95-47-6 | o-xylene | 5 | U |
| 67-64-1 | Acetone | 69 | U UJ |
| 78-93-3 | 2-Butanone | R 60 | U UJ |
| 108-10-1 | 4-Methyl-2-Pentanone | 50 | U UJ |

0018/b
 0020X/0020

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

TBL+R2071300
TB071300

06/18/00

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-10

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2007

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|-----------------------------|--|---|
| 71-43-2 | Benzene | .5 | U |
| 108-86-1 | Bromobenzene | .5 | U |
| 74-97-5 | Bromochloromethane | .5 | U |
| 75-27-4 | Bromodichloromethane | .5 | U |
| 75-25-2 | Bromoform | .5 | U |
| 74-83-9 | Bromomethane | .5 | U |
| 104-51-8 | n-Butylbenzene | .5 | U |
| 135-98-8 | Sec-Butylbenzene | .5 | U |
| 98-06-6 | Tert-Butylbenzene | .5 | U |
| 56-23-5 | Carbon Tetrachloride | .5 | U |
| 108-90-7 | Chlorobenzene | .5 | U |
| 124-48-1 | Dibromochloromethane | .5 | U |
| 75-00-3 | Chloroethane | .5 | U |
| 67-66-3 | Chloroform | .72 | U |
| 74-87-3 | Chloromethane | .5 | U |
| 95-49-8 | 2-Chlorotoluene | .5 | U |
| 106-43-4 | 4-Chlorotoluene | .5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 2 | U |
| 106-93-4 | 1,2-Dibromoethane | .5 | U |
| 74-95-3 | Dibromomethane | .5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | .5 | U |
| 541-73-1 | 1,3-Dichlorobenzene | .5 | U |
| 106-46-7 | 1,4-Dichlorobenzene | .5 | U |
| 75-71-8 | Dichlorodifluoromethane | .5 | U |
| 75-34-3 | 1,1-Dichloroethane | .5 | U |
| 107-06-2 | 1,2-Dichloroethane | .5 | U |
| 75-35-4 | 1,1-Dichloroethene | .5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | .5 | U |
| 156-60-5 | trans-1,2-Dichloroethene | .5 | U |
| 78-87-5 | 1,2-Dichloropropane | .5 | U |
| 142-28-9 | 1,3-Dichloropropane | .5 | U |
| 544-20-7 | 2,2-Dichloropropane | .5 | U |
| 563-58-6 | 1,1-Dichloropropene | .5 | U |
| 100-41-4 | Ethylbenzene | .5 | U |
| 87-68-3 | Hexachlorobutadiene | .5 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT ID

TBL+RRO71300
TB071300
CAEA/9/00

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix: (soil/water)WATER

Lab Sample ID: 001470A-10

Sample wt/vol: 25 (g/mL)ML

Lab File ID: >L2007

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/21/00

GC Column: 007-624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| CAS NO. | COMPOUND | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L | Q |
|----------|---------------------------|--|---|
| 98-82-8 | Isopropylbenzene | .5 | U |
| 99-87-6 | 4-Isopropyltoluene | .5 | U |
| 75-09-2 | Methylene Chloride | 3 | U |
| 91-20-3 | Naphthalene | .5 | U |
| 103-65-1 | N-Propylbenzene | .5 | U |
| 100-42-5 | Styrene | .5 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | .5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | .5 | U |
| 127-18-4 | Tetrachloroethene | .5 | U |
| 108-88-3 | Toluene | .5 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | .5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | .5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | .5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | .5 | U |
| 79-01-6 | Trichloroethene | .5 | U |
| 75-69-4 | Trichlorofluoromethane | .5 | U |
| 96-18-4 | 1,2,3-Trichloropropane | .5 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | .5 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | .5 | U |
| 75-01-4 | Vinyl Chloride | 0.5 → 2 | U |
| NO CAS | mp-Xylene | 1 | U |
| 95-47-6 | o-xylene | .5 | U |
| 67-64-1 | Acetone | 4.7 | J |
| 78-93-3 | 2-Butanone | R 6 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 5 | U |

CAE
9/8/00

CAE
9/8/00

CAE
9/8/00

ATTACHMENT B

**CORRECTED FORM II
GC/MS Volatiles Injection Logs
STL/Connecticut Case No. 1470A**

2A
 WATER VOLATILE SYSTEM MONITORING COMPOUND RECOVERY ⁰⁰⁵⁰

Lab Name: STL/CT

Contract: _____

Lab Code: IEACT

Case No.: 1470A

SAS No.: _____

SDG No.: A1470

| | EPA SAMPLE NO. | SMC1 (DCA) # | SMC2 (DBFM) # | SMC3 (TOL) # | SMC4 (BFB) # | TOT OUT |
|----|-------------------|-----------------|------------------|-----------------|-----------------|------------|
| 01 | VBLKLN | 100 | 107 | 106 | 106 | 0 |
| 02 | CW-5BFMS | 93 | 102 | 99 | 95 | 0 |
| 03 | CW-5BFMSD | 88 | 91 | 85 | 85 | 0 |
| 04 | VBLKLO | 115 | 123 | 121 | 122 | 0 |
| 05 | MW-102A | 113 | 115 | 107 | 108 | 0 |
| 06 | DUP-L+RR | 114 | 115 | 107 | 108 | 0 |
| 07 | MW-104A | 105 | 109 | 104 | 100 | 0 |
| 08 | VBLKLM | 89 | 90 | 88 | 91 | 0 |
| 09 | CW-5B | 109 | 110 | 105 | 110 | 0 |
| 10 | TB071300 | 111 | 108 | 105 | 108 | 0 |
| 11 | TB071200 | 110 | 110 | 106 | 110 | 0 |
| 12 | MW-201 | 112 | 110 | 106 | 108 | 0 |
| 13 | MW-202 | 111 | 110 | 107 | 111 | 0 |
| 14 | MW-103A | 110 | 109 | 106 | 109 | 0 |
| 15 | CW-7A | 110 | 110 | 104 | 107 | 0 |
| 16 | | | | | | |
| 17 | | | | | | |
| 18 | | | | | | |
| 19 | | | | | | |
| 20 | | | | | | |
| 21 | | | | | | |
| 22 | | | | | | |
| 23 | | | | | | |
| 24 | | | | | | |
| 25 | | | | | | |
| 26 | | | | | | |
| 27 | | | | | | |
| 28 | | | | | | |
| 29 | | | | | | |
| 30 | | | | | | |

QC LIMITS

SMC1 (DCA) = 1,2-Dichloroethane-d4 (83-143)
 SMC2 (DBFM) = ~~1,2-Dichloroethane-d4~~ (75-125)
 SMC3 (TOL) = Toluene-d8 (75-125)
 SMC4 (BFB) = Bromofluorobenzene (75-125)

*Dibromofluoro-
methane
CAE 8/30/00*

Column to be used to flag recovery values

* Values outside of contract required QC limits

GC/MS VOLATILES INJECTION LOG

Instr. HP5971L

| | | |
|---------------------|--------------------------------|--------------------|
| Standards Codes: | Routine Maintenance Performed: | Date: 7/21/00 |
| P: 042400:51B:MSZ.8 | P: 071800:82A:MSZ | QC Batch: L0835 |
| P: 070500:71B | P: 82B | IDfile: L821008.05 |
| P: 71C | P: 83C | Calib file: |
| P: 071900:83C | | Method file: |

Calibrated for the following (please check): Protocol = SW846 CLP NYSDEC 524.2 Other
 List = Talk PP 524.2 Other

| Date file | CLP Data file | Client/Job# | Sample ID | Inj. Time | ALS # | DF | QF | Analysis | Comments | pH/WT |
|-----------|---------------|-------------|-----------|-----------|-------|----|----|----------|----------|-------|
| LB201 | | | 55M82B | | 15 | 1 | 1 | ✓ | | |
| LB202 | | | 7 | | 16 | 1 | 1 | | | |
| LB203 | | | | | 17 | 1 | 1 | | | |
| LB204 | | | 914 | | 18 | 1 | 1 | | | |
| L20154 | | | NSTD01LN | | 19 | 1 | 1 | | ok | |
| L20165 | | | NSTD01LN | | 20 | 1 | 1 | | | |
| L20176 | | | PCE 1pt | | 21 | 1 | 1 | | | |
| L20187 | | | VBKLN | | 22 | 1 | 1 | | | |
| L20198 | | | VBKLN | | 23 | 1 | 1 | | ok | |
| L2019 | | RES | 010020 | | 24 | 1 | 1 | | | |
| L2020 | | 00N70A | A-10FMS | | 25 | 1 | 1 | | EW-5B | 1.6 |
| L2021 | | | A-10FMSD | | 26 | 1 | 1 | | ↓ | 1.6 |
| L2022 | | | dup blk | | 27 | 1 | 1 | | blk. C/O | - |
| L2023 | | | A-8 | | 28 | 1 | 1 | | MW-102A | 1.9 |
| L2024 | | | A-9 | | 29 | 1 | 1 | | dup-LTR | 1.6 |
| L2025 | | | A-10 | | 30 | 1 | 1 | | IBL-102A | 1.9 |
| L2026 | | | | | 31 | | | | | |
| L2027 | | | | | 32 | | | | | |

✓ OK
9/8/00

Witnessed by: Ram Date: 7/27/00



TRILLIUM INC.
Consultants in Environmental Chemistry

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August 31, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review Report - STL/Illinois Lot No. 9A08G369, STL/Connecticut Case No. 1470A - Total and Dissolved Metals in Ground Water

Dear Dan:

I have completed my evaluation of the inorganics analysis data prepared by the Severn Trent Laboratory (STL) in University Park, Illinois (under subcontract to STL/Connecticut), for eight ground water (GW) samples from the L&RR Site. These data were reported in a single data package under STL/Illinois Lot No. 9A08G369 (STL/Connecticut Case No. 1470A) which was received by Trillium for review on August 28, 2000. The following GW samples were reported:

| | |
|---------|----------|
| MW-201 | MW-202 |
| MW-103A | CW-7A |
| CW-5B | MW-104A |
| MW-102A | DUP-L+RR |

Analyses were performed according to EPA Methods 7060A (total and dissolved arsenic) and 7421A (total and dissolved lead); both methods are found in the Third Edition of SW-846, Update III. My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced methods, and the USEPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses* (6/88; Rev. 2/89). Qualifiers consistent with those defined in the Region I document were applied as necessary and appropriate.

Although the QAPP specifies that Level III data packages are required for all analyses except the 5-year sampling events, full raw data deliverables (i.e., GFAA instrument print-outs) were provided for all of the metals analyses. ***Unless otherwise noted, these raw data were not reviewed by Trillium as part of this evaluation.*** As specified in Section 9.2 of the QAPP (page 23), a Level III evaluation, limited to an overview of the information provided on the summary forms, was performed.

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Mr. Dan Garrigan

Evaluation of STL/Illinois Lot No. 9A08G369, STL/Connecticut Case No. 1470A (Metals in GW)

August 31, 2000

Page 2

Results from analysis of the dissolved metals sample fractions were distinguished by the laboratory from analysis results for the total metals fractions by the addition of the suffix "S" (for "soluble") to the client sample identification (CW-5BS, for example). The word "Dissolved" was also added by the evaluator to the results forms for the dissolved sample analyses in Attachment A to emphasize the distinction, as well as for consistency with previous data sets generated in support of this project.

Based on the evaluation, the following qualifiers were applied:

- Results for total lead in MW-201 and for dissolved lead in MW-202S were qualified as estimated (J) based on an unacceptably high recovery in the associated CRDL standard.
- Results for lead in MW-201, MW-201S, MW-202, MW-202S, MW-103A, MW-103AS, CW-5B, CW-5BS, CW-7A, CW-7AS, MW-102A, MW-102AS, DUP-L+RR, DUP-L+RRS, and MW-104AS were qualified as estimated (J, UJ) due to negative responses in the associated calibration blanks.
- Results for total and dissolved lead in all samples were qualified as estimated (J, UJ) based on unacceptably low recoveries in the matrix spike analyses.
- The result for total arsenic in MW-103A was qualified as estimated (J) due to an unacceptably high post-digestion spike recovery.
- Results for total lead in MW-202, CW-5B, CW-7A, MW-104A, MW-102A, and DUP-L+RR, and for dissolved lead in MW-202S, CW-5BS, CW-7AS, MW-102AS, and DUP-L+RRS were qualified as estimated (J, UJ) due to unacceptably low post-digestion spike recoveries.
- Results for dissolved arsenic in MW-102AS and DUP-L+RRS were qualified as estimated (J) based on poor reproducibility in the field duplicate analyses.
- Results for total and dissolved lead in MW-202 were qualified as estimated (J) due to discrepancies between the measured total and dissolved concentrations.

All "B," "N," and "W" qualifiers applied by the laboratory were removed by the evaluator.

These qualifiers are reflected on the Inorganic Analysis Data Sheets (Form Is) included as Attachment A to this report. Region I qualifier definitions are also provided in this attachment. Where a sample result was qualified as estimated for more than one reason (as listed above), the qualifier was applied once and no further action was taken.

Mr. Dan Garrigan

Evaluation of STL/Illinois Lot No. 9A08G369, STL/Connecticut Case No. 1470A (Metals in GW)

August 31, 2000

Page 3

At the discretion of the data user, the following points may warrant attention by laboratory and field personnel with regard to this data package and/or prior to future sampling events at this site:

- Documentation of sample preservation in the field on all chain of custody records.
- Documentation of cooler temperatures and sample pHs on laboratory receipt on all chain of custody records.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation and Sample Integrity: Copies of the two applicable field chain of custody (COC) records were not included in this data package. A "Subcontracting Requisition Form" was provided and apparently serves as the COC for the shipment of the samples from STL/Connecticut to STL/Illinois. The samples were shipped on 8/9/00 and received on 8/10/00. No shipping information is recorded on the Requisition Form, and no copy of a courier airbill is included in the data package.

No physical preservation or cooler temperature information was recorded by either laboratory on the Requisition Form. An acceptable temperature (4°C) on laboratory receipt was documented on the field COC for the samples collected on 7/12/00 (as reviewed in associated data packages). On the field COC for samples collected 7/13/00, "7°" was documented by the laboratory. It was assumed that this was also a centigrade measurement. Although slightly above the generally accepted temperature range of 4°C±2°C, the exceedance is very minor and no action was taken on this basis.

Acidification of the samples for dissolved metals analysis with nitric acid was not recorded by the sampler on either of the field COCs and was not recorded by either laboratory on the Requisition Form. Documentation of sample pHs was, however, present on the STL/Illinois digestion logs. All samples were verified to be at pH<2 at the time of preparation for analysis, confirming appropriate preservation.

Field filtration of the samples for dissolved metals analysis was documented on both field COC records.

Total and dissolved metals analyses were performed 8/11/00 through 8/18/00, well within the 6-month holding time specified by the QAPP.

Calibrations: Based on the Analysis Run Logs for each instrument and analysis date (Contract Laboratory Program [CLP] Form XIV), initial and continuing calibration verification (ICV/CCV) standards were run at appropriate frequencies throughout all reported GFAA (graphite furnace atomic absorption) analysis series. All ICV/CCV recoveries documented on the accompanying Initial and Continuing Calibration Verification Reports (CLP Form II) were acceptable (QC 90-110%).

Mr. Dan Garrigan

Evaluation of STL/Illinois Lot No. 9A08G369, STL/Connecticut Case No. 1470A (Metals in GW)

August 31, 2000

Page 4

Linearity check results (i.e., correlation coefficients for the initial multipoint calibrations) for arsenic and lead by GFAA were found in the raw data provided in the data package and were acceptable (>0.995) at the start of each analysis series.

Contract required detection limit (CRDL) standards were run at the beginning of each analysis series. Recoveries were acceptable (80-120%) with the exception of lead (152%) in the 8/11/00 CRDL standard. The high recovery suggests that low measurements of lead may be biased high or represent false positives. Results for total lead in MW-201 and for dissolved lead in MW-202S were qualified as estimated (J) on this basis.

Blanks: Based on the run logs, initial and continuing calibration blanks (ICB/CCBs) were run at appropriate frequencies throughout the reported sample analysis series. No arsenic or lead was reported above the applicable instrument detection limit (IDL) in any of the ICB/CCBs run by GFAA. Lead (-1.8 to -1.2 µg/L) was reported below the negative IDL in many of the CCBs run during all three analysis series, and arsenic was reported below the negative IDL (-1.6 and -1.9 µg/L) in two of the CCBs run during the 8/15/00 GFAA series. Results for lead in MW-201, MW-201S, MW-202, MW-202S, MW-103A, MW-103AS, CW-5B, CW-5BS, CW-7A, CW-7AS, MW-102A, MW-102AS, DUP-L+RR, DUP-L+RRS, and MW-104AS were qualified as estimated (J, UJ) on this basis. Results for arsenic in the sample analyses associated with the affected CCBs on 8/15/00 were greater than the action limit for qualification (two times the absolute value of the associated blank response), therefore no action was taken on this basis.

A water matrix preparation blank (PBW) was digested with the samples (per the preparation logs, CLP Form XIII, and the laboratory's digestion logs); no arsenic or lead contamination was reported in the PBW.

Laboratory Control Samples: Laboratory control sample results were reported in the data package. Acceptable (QC 80-120%) recoveries were reported for both arsenic and lead.

Duplicate Analysis: Sample CW-5B was analyzed in duplicate for total arsenic, total lead, dissolved arsenic, and dissolved lead. No total or dissolved lead or arsenic was detected above the IDL in any of the paired sample analyses, therefore no quantitative evaluation of precision could be made for these analytes using these data.

Matrix Spike Analysis: Sample CW-5B was prepared and analyzed as a matrix spike for total arsenic, total lead, dissolved arsenic, and dissolved lead. Reported percent recoveries for total and dissolved arsenic (98.0% and 104.4%) were acceptable (QC 75-125%). Recoveries for total and dissolved lead, however, were unacceptably low (70.4% and 44.0%). Results for total and dissolved lead in all samples were qualified as estimated (J, UJ) on this basis. The "N" qualifiers applied by the laboratory to all sample results for lead were removed by the evaluator.

Mr. Dan Garrigan

Evaluation of STL/Illinois Lot No. 9A08G369, STL/Connecticut Case No. 1470A (Metals in GW)

August 31, 2000

Page 5

Graphite Furnace Atomic Absorption (GFAA) QC: Post-digestion spike (PDS) recoveries for arsenic and lead were reported on the analysis run logs and were acceptable (QC 85-115%) with the exception of total arsenic in MW-103A (127.9%), CW-5B (124.6%), MW-104A (115.6%), and DUP-L+RR (115.9%), for dissolved arsenic in MW-103AS (138.2%), CW-5BS (121.5%), and DUP-L+RRS (118.1%), for total lead in MW-202 (66.5%), CW-5B (49.5%), CW-7A (75.6%), MW-104A (63.1%), MW-102A (71.7%), and DUP-L+RR (72.8%), and for dissolved lead in MW-202S (81.5%), CW-5BS (41.2%), CW-7AS (77.6%), MW-102AS (72.2%), and DUP-L+RRS (64.0%).

MW-104A, DUP-L+RR and DUP-L+RRS were rerun at five- or ten-fold dilutions for arsenic. The diluted analyses gave acceptable PDS recoveries and the diluted analysis results were reported, therefore no further action was warranted with respect to these sample results.

The positive result for arsenic in MW-103A was qualified as estimated (J) based on the unacceptably high PDS recovery. Arsenic was not detected above the IDL in MW-103S, CW-5B, and CW-5BS and the PDS recoveries were too high; therefore, no qualifiers were applied on this basis.

Results for total lead in MW-202, CW-5B, CW-7A, MW-104A, MW-102A, and DUP-L+RR and for dissolved lead in MW-202S, CW-5BS, CW-7AS, MW-102AS, and DUP-L+RRS were qualified as estimated (J, UJ) because the PDS recoveries were all unacceptably low.

Analysis by the Method of Standard Additions (MSA) was not required for any of the above samples. All "W" qualifiers applied by the laboratory were removed by the validator.

Sample Results: Results greater than or equal to the laboratory-specified IDLs were appropriately reported on the Inorganic Analysis Data Sheets for both analytes in each sample.

All undetected sample results for arsenic and lead were correctly adjusted (multiplied by 0.5) to reflect the concentration factor resulting from processing 100 mL of sample to a final volume of 50 mL, which was clearly documented on the laboratory's digestion logs. Positive sample results were not similarly verified as part of this evaluation.

IDLs for undetected analytes were also appropriately adjusted to reflect sample dilutions, where applicable. Positive sample results were not similarly verified as part of this evaluation.

Form X documentation of laboratory-specified IDLs established on 7/15/00 for arsenic (1.5 µg/L) and lead (1.2 µg/L) on AA4, for arsenic (1.3 µg/L) on AA3, and for lead (1.2 µg/L) on AA6 was provided in the data package. An IDL of 1.5 µg/L was reported by the laboratory for arsenic in all samples; this is acceptable, as it represents the higher of the two IDLs established on the two instruments used for these analyses.



Mr. Dan Garrigan

Evaluation of STL/Illinois Lot No. 9A08G369, STL/Connecticut Case No. 1470A (Metals in GW)

August 31, 2000

Page 6

Field Duplicates: Sample DUP-L+RR was identified as a field duplicate of MW-102A based on information provided by the client. Paired results for total arsenic (26.3 RPD) were acceptable. No total or dissolved lead was reported in either sample, therefore no quantitative evaluation of precision could be performed for this analyte using these data. An unacceptably high RPD was obtained for dissolved arsenic (39.5%; Region I Validation QC \leq 30%). Results for dissolved arsenic in MW-102AS and DUP-L+RRS were qualified as estimated (J) based on this poor reproducibility.

Total versus Dissolved Concentrations: Comparison of total versus dissolved results for lead and arsenic in the ground water samples revealed one case where the dissolved concentration was higher than the total concentration: in MW-202, a low concentration of dissolved lead (0.82 μ g/L) was reported while total lead was not detected above the IDL (0.60 U). Since the positive result was less than twice the sample-specific IDL, this discrepancy most likely reflects the increased variability typically observed at concentrations near the detection limit of the method. Results for total and dissolved lead in MW-202 were qualified as estimated (J, UJ) on this basis.

Documentation: Data package documentation was inconsistent with the specifications of the QAPP; specifically, a Level IV package was provided when a Level III package was required. The presence of the raw data documentation does not adversely affect the reported sample results, but, pursuant to the validation specifications for this project, it was not reviewed by Trillium as part of this evaluation except where otherwise noted above.

Client-assigned sample identifications were modified by the laboratory. Hyphens were eliminated from MW-201, MW-202, CW-5B, and CW-7A; both hyphens and the "A" suffix were eliminated from MW-102A, MW-103A, and MW-104A; and DUP-L+RR was shortened to DUPLR. The laboratory-abbreviated identifications were used consistently throughout the data package. The correct identifications, as found on the COC records, are used throughout this report and have been added to the attached Form Is by the validator.

Please let me know if you have any questions regarding this data package review report.

Sincerely,

Carol A. Erikson
Quality Assessment Manager

CAE/psn

July 2000\GWMetals (95406)

ATTACHMENT A

**INORGANIC ANALYSIS DATA SHEETS
STL/Illinois Lot No. 9A08G369, STL/Connecticut Case No.1470A**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.**
- J - The associated numerical value is an estimated quantity.**
- R - The data are unusable (Note: Analyte may or may not be present).**
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.**

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-201
MW201

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____ SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER _____ Lab Sample ID: 9A08G369-001

Level (low/med): LOW _____ Date Received: 08/10/00

% Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----------|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.75 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 2.5 | | <i>NJ</i> | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

Color Before: COLORLESS _____ Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS _____ Clarity After: CLEAR _____ Artifacts: _____

Comments:
MW-201 _____

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-201 Dissolved
MW201S

Lab Name: STL_CHICAGO Contract: _____

CAE 8/31/00

Lab Code: STL Case No.: _____ SAS No.: _____

SDG No.: U08369

Matrix (soil/water): WATER

Lab Sample ID: 9A08G369-002

Level (low/med): LOW

Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|----|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.75 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.60 | U | UJ | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
MW-201 SOLUBLE

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-202
MW202

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____

SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER

Lab Sample ID: 9A08G369-003

Level (low/med): LOW__

Date Received: 08/10/00

% Solids: __0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L__

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|--------------|------------|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.75 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.60 | U | <i>UJW</i> | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

Color Before: COLORLESS Clarity Before: CLEAR__ Texture: _____

Color After: COLORLESS Clarity After: CLEAR__ Artifacts: _____

Comments:
MW-202 _____

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-202 Dissolved
MW202S

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369 *08/31/00*

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-004

Level (low/med): LOW Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|--------------|-----------|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.75 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.82 | F | <i>JW</i> | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

08/31/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
MW-202 SOLUBLE

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-103A
MW103

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-005

Level (low/med): LOW Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|----------|------------|----------------------|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.80 | <i>✓</i> | <i>AJ</i> | F <i>CAE 8/31/00</i> |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.60 | <i>✓</i> | <i>AUJ</i> | F <i>CAE 8/31/00</i> |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
MW-103A

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-103A Dissolved
MW103S

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-006

Level (low/med): LOW Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|----------|----------|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.75 | U | <i>W</i> | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.60 | <i>U</i> | <i>W</i> | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

CAE 8/31/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

MW-103A SOLUBLE

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CW-7A
CW7A

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____

SDG No.: U08369

CAE 8/31/00

Matrix (soil/water): WATER

Lab Sample ID: 9A08G369-007

Level (low/med): LOW

Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|----|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 9.4 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 3.0 | | UJ | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
CW-7A

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CW-7A Dissolved
CW7AS

Lab Name: STL_CHICAGO Contract: _____

CE 8/31/00

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-008

Level (low/med): LOW Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 9.1 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 3.0 | ✓ | WJA | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CE 8/31/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
CW-7A SOLUBLE

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CW-5B
CW5B

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____ SDG No.: U08369

Matrix (soil/water): WATER _____ Lab Sample ID: 9A08G369-009

Level (low/med): LOW _____ Date Received: 08/10/00

% Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|------|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.75 | U | ✓ | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.60 | ✓ | UJ/A | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

CAE 8/31/00

Color Before: COLORLESS _____ Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS _____ Clarity After: CLEAR _____ Artifacts: _____

Comments:

CW-5B

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CW-5B Dissolved
CW5BS

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-010

Level (low/med): LOW Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|------|----------------------|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.75 | U | / | F <i>CAE 8/31/00</i> |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.60 | / | UJWA | F <i>CAE 8/31/00</i> |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
CW-5B SOLUBLE _____

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-104A
MW104

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-011

Level (low/med): LOW Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|-----------|---|----------------------|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 94.7 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 4.2 | <i>JW</i> | | F <i>CAE 8/31/00</i> |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: BROWN Clarity Before: CLOUDY Texture: _____

Color After: YELLOW Clarity After: CLEAR Artifacts: _____

Comments:
MW-104A

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-104A Dissolved
MW104S

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-012

Level (low/med): LOW Date Received: 08/10/00

* Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 86.9 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 3.0 | ✓ | *UJ | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
MW-104A SOLUBLE

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-102A
MW102

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____ SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER _____ Lab Sample ID: 9A08G369-013

Level (low/med): LOW _____ Date Received: 08/10/00

% Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|------------|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 18.5 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 3.0 | <i>WJW</i> | | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

Color Before: COLORLESS _____ Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS _____ Clarity After: CLEAR _____ Artifacts: _____

Comments:

MW-102A

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-102A Dissolved
MW102S

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369

CAE 8/31/00

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-014

Level (low/med): LOW Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-----|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 17.6 | | J | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 3.0 | | USA | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

CAE 8/31/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
MW-102A SOLUBLE

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

DUP-L+RR
DUPLR

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____ SDG No.: U08369 *CAE 8/31/00*

Matrix (soil/water): WATER _____ Lab Sample ID: 9A08G369-015

Level (low/med): LOW _____ Date Received: 08/10/00

% Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-------------|----------------------|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 14.2 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.60 | | <i>UJ W</i> | F <i>CAE 8/31/00</i> |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS _____ Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS _____ Clarity After: CLEAR _____ Artifacts: _____

Comments:
DUP-LRR _____

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

DUP-LRR Dissolved
DUPLRS

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U08369

CAE 8/31/00

Matrix (soil/water): WATER Lab Sample ID: 9A08G369-016

Level (low/med): LOW Date Received: 08/10/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|-------|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 11.8 | | J | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | 0.60 | | UT/ND | F |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CAE 8/31/00

CAE 8/31/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
DUP-LRR SOLUBLE



TRILLIUM INC.
Consultants in Environmental Chemistry

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October 5, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review Report - STL Case No. 1470A - Total Iron in Ground Water

Dear Dan:

I have completed my evaluation of the inorganics analysis data prepared by the Severn Trent Laboratory (STL) in Monroe, Connecticut, for eight ground water (GW) samples from the L&RR Site. These data were reported in a single data package, under Case No. 1470A, which was received by Trillium for review on October 5, 2000. The following GW samples were reported:

MW-201
MW-103A
CW-5B
MW-102A

MW-202
CW-7A
MW-104A
DUP-L+RR

This data package includes re-analysis data for all eight samples. A data package received on August 28, 2000, presented the original analysis results. Review of the original data revealed an excessively large discrepancy between the results for the field duplicate pair (MW-102A and DUP-L+RR). Historically, the field duplicate pair has shown very good reproducibility, and paired results for all other analyzed parameters did match quite well. At the request of the evaluator, the laboratory rechecked their calculations and verified that the correct values were reported based on the raw data. Subsequently, the laboratory ran "quick and dirty" reruns of the field duplicate samples, and found the paired results to be very consistent. Therefore, re-analysis of all eight samples for total iron was requested by the evaluator, due to the possibility that one or more additional sample results could be wrong if two samples had been switched at some point during the original preparations. This did, in fact, turn out to be the case. The following table summarizes the original and re-analysis results for total iron:

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Mr. Dan Garrigan
Evaluation of STL Case No. 1470A (Total Iron in GW)
October 5, 2000
Page 2

| Sample | Original Result (µg/L) | Re-Analysis Result (µg/L) |
|----------|------------------------|---------------------------|
| MW-201 | 351 | 1530 |
| MW-202 | 321 | 7790 |
| MW-103A | 47700 | 13.8 |
| CW-7A | 15800 | 15200 |
| CW-5B | 3780 | 3010 |
| MW-104A | 82600 | 89100 |
| MW-102A | 33.8 | 37200 |
| DUP-L+RR | 43900 | 36600 |

From these data, it appears that samples MW-103A and MW-102A were switched at some point in the original preparation and analysis process for total iron. The re-analysis results are also consistent with results obtained for total iron in the April 2000 samples collected from this site. Therefore, only the re-analysis results were reviewed and discussed in the remainder of this report. The original analysis results (pages 78-105 of the August 25, 2000 data package) have been clipped separately and marked "Do Not Use" for clarity.

Analyses were performed according to EPA Method 6010B, which is found in the Third Edition of SW-846, Update III. My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced method, and the USEPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses* (6/88; Rev. 2/89). Qualifiers consistent with those defined in the Region I document were applied as necessary and appropriate.

No metals analysis raw data were provided in the data package; a Level III evaluation of the summary forms as presented was performed, as required by the QAPP.

Based on the evaluation, the following action was taken:

- The result for total iron in MW-103A was qualified as estimated (J) because it was less than twice the instrument detection limit. The "B" qualifier applied by the laboratory was removed by the evaluator.

Mr. Dan Garrigan
Evaluation of STL Case No. 1470A (Total Iron in GW)
October 5, 2000
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This action is reflected on the Inorganic Analysis Data Sheet (Form I) for MW-103A included in Attachment A to this report. Region I qualifier definitions are also provided in this attachment.

At the discretion of the data user, the following points may warrant attention by laboratory and field personnel with regard to this data package and/or prior to future sampling events at this site:

- Documentation of sample preservation in the field on all chain of custody records.
- Documentation of complete cooler temperatures and sample pHs on laboratory receipt directly on all chain of custody records.
- Use of updated (within three months of sample analyses) instrument detection limits in association with every data package.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation and Sample Integrity: Copies of the applicable chain of custody (COC) records were not included in the 10/4/00 data package, however these records were available in the 8/25/00 data package (which also included results for other analysis parameters). The following discussion refers to review of those documents.

The two applicable COC records documented sample collection dates of 7/12/00 and 7/13/00. An acceptable temperature (4°C) on laboratory receipt was documented on the COC for the samples collected on 7/12/00. On the COC for samples collected 7/13/00, "7°" was documented by the laboratory; it was assumed that this was also a centigrade measurement. Although slightly above the generally accepted temperature range of 4°C±2°C, the exceedance is very minor and no action was taken on this basis.

Acidification of the samples for total iron analysis with nitric acid was not recorded by the sampler on either of the COCs and documentation of sample pHs on laboratory receipt or at the time of preparation for analysis was also not included in the data package. At the evaluator's request, the laboratory provided their sample receiving preservative records for all samples, which verified that the sample containers preserved with nitric acid were at pH<2 on receipt from the field in all cases. No further action was taken on this basis.

The re-analyses for total iron were performed on 9/25/00, well within the 6-month holding time specified by the QAPP.

Calibrations: Based on the Analysis Run Log (Contract Laboratory Program [CLP] Form XIV), initial and continuing calibration verification (ICV/CCV) standards were run at appropriate frequencies throughout the ICP (inductively coupled plasma) analysis series. All ICV/CCV

Mr. Dan Garrigan

Evaluation of STL Case No. 1470A (Total Iron in GW)

October 5, 2000

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recoveries documented on the accompanying Initial and Continuing Calibration Verification Reports (CLP Form II) were acceptable (QC 90-110%).

Blanks: Based on the run log, initial and continuing calibration blanks (ICB/CCBs) were run at appropriate frequencies throughout the reported sample analysis series. Iron was not detected above the instrument detection limit (IDL) in any of the ICB/CCBs run during the ICP analysis series.

A water matrix preparation blank (PBW) was digested with the samples (per the preparation log, CLP Form XIII). Iron was not detected in the PBW.

ICP Interference Check Sample: No interference check sample results were reported in the data package, although entries for these analyses were found on the run log.

Laboratory Control Samples: Laboratory control sample results were reported in the data package. An acceptable (QC 80-120%) recovery was reported for iron (91.0%).

Duplicate Analysis: Sample CW-5B was analyzed in duplicate for total iron, and excellent reproducibility was demonstrated (2.9 relative percent difference [RPD]).

Matrix Spike Analysis: Sample CW-5B was prepared and analyzed as a matrix spike for total iron. The reported percent recovery (80.9%) was acceptable (QC 75-125%).

ICP Serial Dilution: An acceptable serial dilution analysis result (QC \leq 15% difference, per Region I validation guidance) for iron in MW-104A (11.9%) was reported in the data package.

Sample Results: Results greater than or equal to the laboratory-specified IDL were appropriately reported on the Inorganic Analysis Data Sheets for total iron in each sample.

Form X documentation of the laboratory-specified IDL established for iron on 4/15/00 (7.2 $\mu\text{g/L}$) was provided in the data package. IDLs should be established quarterly and this is much more than three months prior to the sample analyses. The laboratory should be instructed to establish and provide documentation of updated IDLs in association with every data package.

The result for total iron in MW-103A was greater than the applicable IDL but below the laboratory-specified reporting limit (RL) and was correctly reported by the laboratory with a "B" qualifier. As concentrations approach the IDL the accuracy of the measurement decreases; values closer to the RL, however, are probably quite accurate. Therefore, a guideline of 2xIDL was used to determine whether the reported result warranted qualification; specifically, sample results below the respective RL, less than 2xIDL and not otherwise qualified warrant qualification as estimated. The result for total iron in MW-103A was qualified as estimated (J) on this basis, and the "B" qualifier applied by the laboratory was removed by the validator.

Mr. Dan Garrigan
Evaluation of STL Case No. 1470A (Total Iron in GW)
October 5, 2000
Page 5

Field Duplicates: Sample DUP-L+RR was identified as a field duplicate of MW-102A based on information provided by the client. Paired results for total iron showed excellent reproducibility, with an RPD of 1.6%.

Documentation: Data package documentation was consistent with the specifications of the QAPP; specifically, a Level III package containing summary forms only with no supporting raw data was provided.

Documentation of sample pHs on laboratory receipt was not provided in the data package as received for review. This information was verbally requested by the evaluator on 8/30/00 (in association with the original data package); sample receiving preservation logs were received from the laboratory via facsimile on 9/2/00. These records were inserted into the 8/25/00 data package by the evaluator as pages 55a and 55b.

The narrative included in the 10/4/00 data package is identical to the narrative in the 8/25/00 data package. The statement, "Iron failed the controls for spike recovery analysis of sample CW-5B resulting in one 'N' flag." is not applicable to the 10/4/00 data package and was deleted by the evaluator.

No interference check sample results were reported, although the run log indicates that these quality control analyses were performed.

The preparation and analysis summary form on pages 10-11 of the 10/4/00 data package implies that samples were prepared and analyzed on 9/25/00 for total and dissolved arsenic and lead as well as for total iron. Only total iron reruns were requested, and only total iron data were reported in the data package, although it is certainly possible that the additional preparations were unnecessarily performed by the laboratory.

Please let me know if you have any questions regarding this data package review report.

Sincerely,



Carol A. Erikson
Quality Assessment Manager

CAE/esc
July 2000\GWTtotalFe (95406)

ATTACHMENT A

**INORGANIC ANALYSIS DATA SHEETS
Laboratory Case No. 1470A**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.**
- J - The associated numerical value is an estimated quantity.**
- R - The data are unusable (Note: Analyte may or may not be present).**
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.**

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-201

Lab Name: STL

Contract: _____

Lab Code: STL Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix (soil/water): WATER

Lab Sample ID: T001470A-01

Level (low/med): LOW

Date Received: 07/13/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | | | | NR |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | 1530 | | | P |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| 57-12-5 | Cyanide | | | | NR |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

Total Metals

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-202

Lab Name: STL

Contract: _____

Lab Code: STL Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix (soil/water): WATER

Lab Sample ID: T001470A-02

Level (low/med): LOW

Date Received: 07/13/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | | | | NR |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | 7790 | | | P |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| 57-12-5 | Cyanide | | | | NR |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

Total Metals

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-103A

Lab Name: STL

Contract: _____

Lab Code: STL Case No.: 1470A

SAS No.: _____

SDG No.: A1470Matrix (soil/water): WATERLab Sample ID: T001470A-03Level (low/med): LOWDate Received: 07/13/00% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---------|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | | | | NR |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | 13.8 | # | J | P |
| 7439-92-1 | Lead | | | CAUTION | NR |
| 7439-95-4 | Magnesium | | | 10/5/00 | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| 57-12-5 | Cyanide | | | | NR |

Color Before: COLORLESSClarity Before: CLEAR

Texture: _____

Color After: COLORLESSClarity After: CLEAR

Artifacts: _____

Comments:

Total Metals

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CW-7A

Lab Name: STL

Contract: _____

Lab Code: STL Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix (soil/water): WATER

Lab Sample ID: T001470A-04

Level (low/med): LOW

Date Received: 07/13/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | | | | NR |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | 15200 | | | F |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| 57-12-5 | Cyanide | | | | NR |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

Total Metals

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CW-5B

Lab Name: STL

Contract: _____

Lab Code: STL Case No.: 1470A

SAS No.: _____

SDG No.: A1470Matrix (soil/water): WATERLab Sample ID: T001470A-06Level (low/med): LOWDate Received: 07/14/00% Solids: 0.0Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | | | | NR |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | 3010 | | | P |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| 57-12-5 | Cyanide | | | | NR |

Color Before: COLORLESSClarity Before: CLEAR

Texture: _____

Color After: COLORLESSClarity After: CLEAR

Artifacts: _____

Comments:

Total Metals

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-104A

Lab Name: STL

Contract: _____

Lab Code: STL Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix (soil/water): WATER

Lab Sample ID: T001470A-07

Level (low/med): LOW

Date Received: 07/14/00

* Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | | | | NR |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | 89100 | | | P |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| 57-12-5 | Cyanide | | | | NR |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

Total Metals

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

MW-102A

Lab Name: STL

Contract: _____

Lab Code: STL Case No.: 1470A

SAS No.: _____

SDG No.: A1470

Matrix (soil/water): WATER

Lab Sample ID: T001470A-08

Level (low/med): LOW

Date Received: 07/14/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | | | | NR |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | 37200 | | | P |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| 57-12-5 | Cyanide | | | | NR |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:

Total Metals

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

| |
|----------|
| DUP-L+RR |
|----------|

Lab Name: STL

Contract: _____

Lab Code: STL Case No.: 1470A

SAS No.: _____

SDG No.: A1470Matrix (soil/water): WATERLab Sample ID: T001470A-09Level (low/med): LOWDate Received: 07/14/00% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | | | | NR |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | 36600 | | | P |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| 57-12-5 | Cyanide | | | | NR |

Color Before: COLORLESSClarity Before: CLEAR

Texture: _____

Color After: COLORLESSClarity After: CLEAR

Artifacts: _____

Comments:

Total Metals



TRILLIUM INC.
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September 14, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review Report - STL/Connecticut Case No. 1470A - Ammonia Chloride, and BOD in Ground Water

Dear Dan:

I have completed my evaluation of the wet chemistry analysis data prepared by the Severn Trent Laboratory in Shelton, Connecticut (STL/Connecticut), for eight ground water samples from the L&RR Site, which were reported in a single data package under STL Case No. 1470A. The data package was received by Trillium for review on August 28, 2000, with additional requested documentation provided on September 2 and 11, 2000. The following samples were reported:

| | |
|---------|----------|
| MW-201 | MW-202 |
| MW-103A | CW-7A |
| CW-5B | MW-104A |
| MW-102A | DUP-L+RR |

Analyses were performed according to EPA Methods 325.2 (chloride), 350.1 (ammonia), and 405.1 (BOD); these three methods are found in *Methods for Chemical Analysis of Water and Wastes* (EPA-600/4-79-020, 3/83). My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced methods, and, to the extent applicable, the USEPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses* (6/88; Rev. 2/89). Qualifiers consistent with those defined in the Region I document were applied as necessary and appropriate.

No raw data were provided in the data package as initially submitted for review. As specified by the QAPP, only a Level III evaluation of the summary results was required. However, for these parameters, the raw data pages also serve as analysis run logs and are necessary to verify calibrations when spectrophotometric methods are used as well as generally helpful for verifying that all samples were analyzed. Raw data documentation has also been provided in previous data packages generated

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Mr. Dan Garrigan

Evaluation of STL/Connecticut Case No. 1470A (Wet Chemistry Parameters in GW)

September 14, 2000

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in support of this project. Additional documentation was therefore requested from and received by the laboratory as discussed throughout this report.

Based on the evaluation, results for BOD in all samples collected on 7/12/00 (MW-201, MW-202, MW-103A, and CW-7A) were qualified as estimated (J, UJ) because incubation was started after the 48 hour holding time for this parameter had expired. All other sample results were determined to be valid as reported; no qualifiers were applied.

At the discretion of the data user, the following points may warrant attention by field and laboratory personnel with regard to this data package and prior to future sampling events at this site:

- Documentation of sample preservation performed in the field on all chain of custody records.
- Documentation of complete cooler temperatures and sample pHs at the time of laboratory receipt on all chain of custody records.
- Inclusion of raw data as needed to document calibrations and analysis of all samples.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation and Sample Integrity: Copies of the two applicable chain of custody (COC) records were included in the data package, documenting sample collection dates of 7/12/00 and 7/13/00. An acceptable temperature (4°C) on laboratory receipt was documented on the COC for the samples collected on 7/12/00. On the COC for samples collected 7/13/00, "7°" was documented by the laboratory. It was assumed that this was also a centigrade measurement. Although slightly above the generally accepted temperature range of 4°C ±2°C, the exceedance is very minor and no action was taken on this basis.

Acidification of the samples for ammonia analysis with sulfuric acid was recorded by the sampler on only one of the COCs and documentation of sample pHs on laboratory receipt or at the time of analysis was not included in the data package. At the evaluator's request, STL/Connecticut provided their sample receiving preservative records for all samples, which verified that the sample containers preserved with sulfuric acid were at pH<2 on receipt from the field in all cases. No further action was taken on this basis.

Mr. Dan Garrigan

Evaluation of STL/Connecticut Case No. 1470A (Wet Chemistry Parameters in GW)

September 14, 2000

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All sample analyses for chloride and ammonia were performed within the 28-day holding time specified by the QAPP for these two parameters. BOD incubation was initiated on 7/15/00 for all samples; this is within the required 48-hour holding time for those samples collected on 7/13/00 but exceeds the holding time for those samples collected on 7/12/00. Results for BOD in the four samples collected on 7/12/00 (MW-201, MW-202, MW-103A, and CW-7A) were qualified as estimated (J, UJ) on this basis.

Calibrations: Chloride analyses were performed according to EPA 325.2. No documentation of an initial calibration for chloride was provided in the package initially received for review. At the request of the evaluator, copies of the applicable instrument print-outs, including the initial calibration for chloride, were provided by the laboratory via FedEx on 9/11/00. Based on this documentation, an initial calibration (3.0-75 mg/L) was run for chloride on 7/23/00 (immediately prior to sample analyses) and acceptable correlation coefficients (>0.995) were reported for two segments of that curve. Initial and continuing calibration verification standards were run at appropriate frequencies throughout the chloride analysis series, and all recoveries were acceptable (QC 85-115%) as reported on the summary forms provided for review.

Ammonia analyses were performed according to EPA 350.1. No documentation of an initial calibration for ammonia was provided in the package initially received for review. At the request of the evaluator, copies of the applicable instrument print-outs, including the initial calibration for ammonia, were provided by the laboratory via FedEx on 9/11/00. Based on this documentation, an initial calibration (0.04-4.0 mg/L) was run for ammonia on 7/20/00 (immediately prior to sample analyses) and acceptable correlation coefficients (>0.995) were reported for two segments of that curve. Initial and continuing calibration verification standards were run at appropriate frequencies throughout the ammonia analysis series, and all recoveries were acceptable (QC 85-115%) as reported on the summary forms provided for review.

Initial calibration is not applicable to the BOD analyses, which were performed by titration.

Blanks: Method blank results were reported for ammonia, chloride, and BOD on summary forms in the data package; no contamination was reported above the applicable reporting limit in any of these blanks.

Initial and continuing calibration blanks (ICB/CCBs) were run at the appropriate frequencies for both chloride and ammonia and no contamination was detected in any of these blanks, based on the summary forms provided for review.

Mr. Dan Garrigan

Evaluation of STL/Connecticut Case No. 1470A (Wet Chemistry Parameters in GW)

September 14, 2000

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Laboratory Control Samples: Laboratory control samples (LCSs) were analyzed in association with the sample analyses for ammonia and chloride. Acceptable recoveries (87.1% and 103.7%) were reported in both cases.

Duplicate Analysis: Unspiked duplicate analyses of CW-5B were performed for chloride and ammonia. Neither chloride nor ammonia was detected above the reporting limit in either sample analysis; therefore, no evaluation of precision could be made using these data.

Results for an unspiked duplicate analysis of a non-project sample were reported for BOD. These data have no relevance to the site samples included in this data set and were not further reviewed.

Matrix Spike Analysis: Sample CW-5B was prepared and analyzed as a matrix spike for ammonia and chloride. Reported percent recoveries (86.4% and 112.1%) were acceptable (QC 75-125%) for both spikes.

Spiked sample results for a non-project sample were reported for BOD. These data have no relevance to the site samples included in this data set and were not further reviewed.

Sample Results: Sample results greater than or equal to the laboratory-specified reporting limits (RLs) for ammonia, chloride, and BOD were reported on "Wet Chem Analysis Data Sheets" for all samples. Non-detected analytes were reported as less than the applicable RL in all cases. Based on the raw data/run logs provided by the laboratory, some samples were appropriately re-analyzed at dilutions for some parameters.

Field Duplicates: Sample DUP-L+RR was identified as a field duplicate of MW-102A. Positive paired results for chloride (5.4 RPD), ammonia (1.5 RPD), and BOD (22.9 RPD) showed acceptable reproducibility (QC \leq 30%).

Documentation: Data package documentation was consistent with the specifications of the QAPP; specifically, a Level III package containing summary forms only was provided. However, no documentation of initial or continuing calibration standards for chloride and ammonia was provided, although this information has been provided in previous data packages generated in support of this project. For wet chemistry parameters in general, the raw data pages serve as run logs and are helpful for verifying that all samples were, in fact, analyzed. For those parameters run by spectrophotometric methods, the raw data pages are necessary to evaluate calibrations and calibration blanks, which are not routinely summarized elsewhere. Additional documentation for chloride, ammonia, and BOD was therefore requested from and provided by the laboratory on 9/11/00, as discussed previously in this report. The BOD documentation was added to the end of the data package as pages 125-127; the

Mr. Dan Garrigan

Evaluation of STL/Connecticut Case No. 1470A (Wet Chemistry Parameters in GW)

September 14, 2000

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ammonia documentation was added as pages 128-142; and the chloride documentation was added as pages 143-150.

COC records provided in the data package included all reported samples and were properly completed except that sample preservation was recorded on only one of the COCs and improper corrections were noted on both documents. In the future, all corrections should be made by drawing a single line through the incorrect information, inserting the correct information, and initialing and dating the change. "Write-overs" are not legally defensible.

Documentation of sample pHs on laboratory receipt (or at the time of analysis) was not provided in the data package as received for review. This information was verbally requested by the validator on 8/30/00; sample receiving preservation logs were received from the laboratory via facsimile on 9/2/00. These records were inserted into the data package by the evaluator as pages 55a and 55b.

Please let me know if you have any questions regarding this data package review report.

Sincerely,



Carol A. Erikson
Quality Assessment Manager

CAE/esc

July 2000\GWWetChem (95406)

ATTACHMENT A

**WET CHEM ANALYSIS DATA SHEETS (Form I)
STL/Connecticut Case No. 1470A**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.**
- J - The associated numerical value is an estimated quantity.**
- R - The data are unusable (Note: Analyte may or may not be present).**
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.**



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September 11, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review - STL/Illinois Lot No. 9A07G171, STL/Connecticut Case No. 1470A - COD in Ground Waters

Dear Dan:

I have completed my evaluation of the chemical oxygen demand (COD) analysis data prepared by the Severn Trent Laboratory (STL) in University Park, Illinois (under subcontract to STL/Connecticut), for eight ground water samples from the L&RR site. The data were reported in a single data package under STL/Illinois Lot No. 9A07G171, STL/Connecticut Case No. 1470A, which was received by Trillium for review on August 28, 2000 and included the following samples:

| | |
|---------|----------|
| MW-201 | MW-202 |
| MW-103A | CW-7A |
| CW-5B | MW-104A |
| MW-102A | DUP-L+RR |

Analyses were performed according to Hach Method 8000, which is a closed reflux digestion with titration assay, as referenced by the laboratory. My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced method, and, to the extent applicable, the USEPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses* (6/88; Rev. 2/89).

Based on the evaluation, all sample results for COD were determined to be valid as reported. No qualifiers were applied.

At the discretion of the data user, the following points may warrant attention by field and laboratory personnel with regard to this data package and prior to future sampling events at this site:

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Evaluation of STL/Illinois Lot No. 9A07G171, STL/Connecticut Case No. 1470A (COD in GW)

September 11, 2000

Page 2

- Documentation of sample preservation performed in the field on all chain of custody records.
- Documentation of complete cooler temperatures and sample pHs at the time of laboratory receipt on all applicable chain of custody records.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation, Sample Integrity: Copies of the two applicable field chain of custody (COC) records were not included in the data package. A "Subcontracting Requisition Form" was provided, which apparently serves as the COC for the shipment of samples from STL/Connecticut to STL/Illinois. The samples were shipped on 7/19/00 and received on 7/20/00; a copy of a FedEx label is also provided in the data package, however the airbill number is not recorded on the Requisition Form and the airbill is not dated. Therefore, the airbill cannot be reliably linked to the shipment of these samples.

The notation "Temp 2.2" is present in the "Special Instructions" portion of the Requisition Form, but it is not clear when this entry was made or who made it. For the purposes of this evaluation, it was assumed that this was recorded by STL/Illinois on receipt of the samples and that it means that the cooler was at the acceptable temperature of 2.2°C when it arrived.

Acidification of the samples for COD analysis with sulfuric acid was recorded by the sampler on only one of the field COCs (as reviewed in associated data packages) and was not recorded by STL/Connecticut at all on the Requisition Form. Documentation of sample pHs was also not documented by either laboratory anywhere else in this data package. At the request of the validator, STL/Connecticut provided their sample receiving preservative records for all samples, which verified that the sample containers preserved with sulfuric acid were at pH<2 on receipt from the field in all cases. For the purposes of this evaluation, it was assumed that the pHs of the samples for COD analysis did not change during transport from STL/Connecticut to STL/Illinois and no further action was taken on this basis.

All sample analyses for COD were performed within the 28-day holding time specified by the QAPP for this parameter.

Calibrations: Calibration is not applicable to the titrimetric method used for these analyses.

Blanks: Low- and high-level method blanks (MBs) containing no COD above the reporting limit were reported in the data package.

Laboratory Control Samples: Low- and high-level laboratory control samples (LCSs) were analyzed in association with the sample analyses for COD. Acceptable recoveries (114% and 94.0%) were reported in both cases.

Mr. Dan Garrigan

Evaluation of STL/Illinois Lot No. 9A07G171, STL/Connecticut Case No. 1470A (COD in GW)

September 11, 2000

Page 3

Duplicate Analysis: No unspiked laboratory duplicate analysis was performed on any of the samples in this data package.

Matrix Spike Analysis: Sample CW-5B was prepared and analyzed as a matrix spike/matrix spike duplicate (MS/MSD) pair. Reported COD recoveries were acceptable (113% and 102%) and reproducible (relative percent difference, RPD, 9.8%).

Sample Results: Sample results greater than or equal to the laboratory-specified reporting limit (5 mg/L) for COD were reported on summary forms for all samples. MW-104A was analyzed as a high-level sample; all remaining samples were analyzed as low-level samples.

Field Duplicates: Sample DUP-L+RR was identified as a field duplicate of MW-102A. COD was not detected above the RL in either sample, therefore no quantitative evaluation of precision could be performed using these data.

Documentation: Data package documentation was inconsistent with the specifications of the QAPP, in that full raw data documentation was provided when a Level III package containing summary forms only was required. However, for wet chemistry parameters in general, the raw data pages serve as run logs and are helpful for verifying that all samples were, in fact, analyzed. Therefore, the data package documentation was appropriate as provided.

Field COC records were not provided in the data package, and documentation of sample pHs on receipt at STL/Illinois was not found in the data package as received for review. The field COC records were available for review in other data packages associated with this data set, and no action was taken on this basis. Chemical preservation of the samples for COD analysis was verified based on the preservative logs provided by STL/Connecticut, which verified that they were at pH<2 on arrival at that laboratory. It was assumed that the pHs did not change during shipment to STL/Illinois, and no further action was taken on this basis.

Sample results were reported by STL/Illinois using the STL/Connecticut identifications. COC identifications were added to the results forms in Attachment A for the sake of clarity.

Please let me know if you have any questions regarding this data package review report.

Sincerely,



Carol A. Erikson
Quality Assessment Manager

ATTACHMENT A

**COD in Ground Waters
STL/Illinois Lot No. 9A07G171, STL/Connecticut Case No. 1470A**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.**
- J - The associated value is an estimated quantity.**
- R - The data are unusable. (Note: Analyte may or may not be present).**
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.**

To: STL Connecticut
Severn Trent Laboratories
128 Long Hill Cross Road
Shelton, CT 06484

Date: Friday July 28th, 2000

RE: 1470A-01 MW-201 *caeg/30/00*
Project # 00000-000-000-0000
Lab ID: 9A07G171-001
Sample Date: 07/12/00
Date Received: 07/20/00

Attn: Paul Hobart

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|--------|-----------------|
| COD | 5 | u mg/L | 5 |

To: STL Connecticut
Severn Trent Laboratories
128 Long Hill Cross Road
Shelton, CT 06484

Date: Friday July 28th, 2000

RE: 1470A-02 MW-202
Project # 00000-000-000-0000 CAE
Lab ID: 9A07G171-002 8/30/00
Sample Date: 07/12/00
Date Received: 07/20/00

Attn: Paul Hobart

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|--------|-----------------|
| COD | 5 | u mg/L | 5 |

To: STL Connecticut
Severn Trent Laboratories
128 Long Hill Cross Road
Shelton, CT 06484

Date: Friday July 28th, 2000

RE: 1470A-03 MW-103A *CAE 8/30/00*
Project # 00000-000-000-0000
Lab ID: 9A07G171-003
Sample Date: 07/12/00
Date Received: 07/20/00

Attn: Paul Hobart

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|--------|-----------------|
| COD | 5 | u mg/L | 5 |

To: STL Connecticut
Severn Trent Laboratories
128 Long Hill Cross Road
Shelton, CT 06484

Date: Friday July 28th, 2000

RE: 1470A-04 CW-7A *CAE 8/30/00*
Project # 00000-000-000-0000
Lab ID: 9A07G171-004
Sample Date: 07/12/00
Date Received: 07/20/00

Attn: Paul Hobart

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|--------|-----------------|
| COD | 5 | u mg/L | 5 |

To: STL Connecticut
Severn Trent Laboratories
128 Long Hill Cross Road
Shelton, CT 06484

Date: Friday July 28th, 2000

RE: 1470A-06 CW-5B CAE 8/30/00
Project # 00000-000-000-0000
Lab ID: 9A07G171-005
Sample Date: 07/13/00
Date Received: 07/20/00

Attn: Paul Hobart

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|--------|-----------------|
| COD | 5 | u mg/L | 5 |

To: STL Connecticut
Severn Trent Laboratories
128 Long Hill Cross Road
Shelton, CT 06484

Date: Friday July 28th, 2000

RE: 1470A-07 MW-104A CAE 8/30/00
Project # 00000-000-000-0000
Lab ID: 9A07G171-006
Sample Date: 07/13/00
Date Received: 07/20/00

Attn: Paul Hobart

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|-------|-----------------|
| COD | 220 | mg/L | 50 |

To: STL Connecticut
Severn Trent Laboratories
128 Long Hill Cross Road
Shelton, CT 06484

Date: Friday July 28th, 2000

RE: 1470A-08 MW-102A CAE8/30/00
Project # 00000-000-000-0000
Lab ID: 9A07G171-007
Sample Date: 07/13/00
Date Received: 07/20/00

Attn: Paul Hobart

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|--------|-----------------|
| COD | 5 | u mg/L | 5 |

To: STL Connecticut
Severn Trent Laboratories
128 Long Hill Cross Road
Shelton, CT 06484

Attn: Paul Hobart

Date: Friday July 28th, 2000

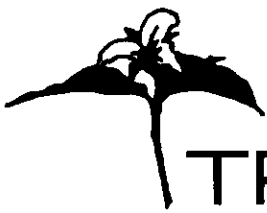
RE: 1470A-09 *DUP-L+RR* *CAE8/30/00*
Project # 00000-000-000-0000
Lab ID: 9A07G171-008
Sample Date: 07/13/00
Date Received: 07/20/00

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|--------|-----------------|
| COD | 5 | u mg/L | 5 |

Appendix D

**Groundwater Data Validation
Reports
Compuchem Environmental
Laboratories**



TRILLIUM INC.
Consultants in Environmental Chemistry

356 FARRAGUT CROSSING DR.
KNOXVILLE, TN 37922

(423) 966-8880

FAX (423) 966-8885

cerikson@trilliuminc.com

August 31, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review Report - CompuChem Case No. Q1458, SDG R1458, VOAs in Ground Water

Dear Dan:

I have completed my evaluation of the volatile organics analysis data prepared by CompuChem Environmental (CompuChem) for one ground water (GW) sample and one trip blank (TB) from the L&RR Site, which were reported in a single data package under Case No. Q1458, SDG R1458. The data package was received by Trillium for review on August 11, 2000, and includes results for the following samples:

TB

MW-104A

Analyses were performed according to EPA Method 8260B (SW-846, Third Edition, Update III). My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced method, and the *Region I, EPA-New England Data Validation Functional Guidelines for Evaluating Environmental Analyses* (12/96). Qualifiers consistent with those defined in the Region I document were applied as necessary and appropriate.

Although the QAPP specifies that Level III data packages are required for all analyses except the 5-year sampling events, full raw data deliverables (i.e., quantitation reports, chromatograms, mass spectra, etc.) were provided for the samples, calibration standards, blanks, spikes and instrument performance checks. ***These raw data were not reviewed by Trillium as part of this evaluation.*** As specified in Section 9.2 of the QAPP (page 23), a Level III evaluation, limited to an overview of the information provided on the summary forms, was performed.

Based on the evaluation, the following qualifiers were applied:

- The result for 2-butanone in TB was rejected (R) based on unacceptably low relative response factors (RRFs) in the associated calibration standards.

HOME OFFICE:

28 GRACE'S DRIVE • COATESVILLE, PA19320 • (610) 383-7233 • FAX (610) 383-7907

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Mr. Dan Garrigan

Evaluation of CompuChem Case No. Q1458, R1458 (VOCs in GW)

August 31, 2000

Page 2 of 6

- Results for acetone in MW-104A and TB and for 2-butanone in MW-104A were qualified as estimated (J) based on unacceptably low RRFs in the associated calibration standards.
- Results for dichlorodifluoromethane, acetone, 2-butanone, 1,3-dichlorobenzene, 1,4-dichlorobenzene, 1,2-dichlorobenzene, 1,2,4-trichlorobenzene, naphthalene, and 1,2,3-trichlorobenzene in MW-104A were qualified as estimated (J, UJ) based on unacceptable percent relative standard deviation (%RSD) values in the associated initial calibration.
- Results for acetone, 1,3-dichlorobenzene, 1,4-dichlorobenzene, 1,2-dichlorobenzene, 1,2-dibromo-3-chloropropane, 1,2,4-trichlorobenzene, naphthalene, and 1,2,3-trichlorobenzene in TB were qualified as estimated (J, UJ) based on unacceptable %RSD values in the associated initial calibration.
- Results for acetone in MW-104A and TB were qualified as less than the reported value or less than the sample-specific reporting limit, whichever was greater (U), based on associated method blank contamination.
- Results for methylene chloride, 1,3-dichlorobenzene, and 1,2,4-trichlorobenzene in MW-104A were qualified as less than the reported value or less than the sample-specific RL, whichever was greater (U), based on associated TB contamination.
- Results for all detected target analytes in TB were qualified as estimated (J) due to unacceptably high recoveries for three of four system monitoring compounds in the reported analysis.
- Results for chloroethane, 1,1-dichloroethane, benzene, ethylbenzene, m/p-xylenes, o-xylene, 1,2,4-trimethylbenzene, 1,4-dichlorobenzene, naphthalene, and total xylenes in MW-104A were qualified as estimated (J) because they exceeded the upper limit of the established calibration range. No quantitatively accurate analysis results are available for these compounds, and the data user is cautioned that the accuracy of the reported values for these compounds cannot be verified. In particular, those results exceeding the upper limit of the calibration range by more than 20% (i.e., values greater than 30 µg/L) should be considered minimum concentrations.
- All "B" and "E" qualifiers appropriately applied by the laboratory were removed by the validator.

Mr. Dan Garrigan

Evaluation of CompuChem Case No. Q1458, R1458 (VOCs in GW)

August 31, 2000

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These qualifiers are reflected on the affected Organic Analysis Data Sheets (Form I) included as Attachment A to this report. Region I qualifier definitions are also provided in this attachment. Where a sample result was qualified as estimated for more than one reason (as listed above), the qualifier was applied once and no further action was taken.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation, Sample Integrity: A copy of the field chain of custody (COC) record was included in the data package, documenting a sample collection date of 7/13/00 for both samples. Volatile organics analyses were performed on 7/27/00, just within the required 14-day holding time. An acceptable ($4^{\circ}\text{C} \pm 2^{\circ}\text{C}$) cooler temperature (3°C) on laboratory receipt was recorded on the COC. Acidification of the samples for volatiles analysis with hydrochloric acid was documented by the sampler on the COC; although sample pHs on laboratory receipt were not documented on the COC, verification of successful acidification of both samples (pH=1) was confirmed on the Water Batch Sheet prepared by the laboratory and included in the data package.

GC/MS Instrument Performance Check: Form V was provided for the single bromofluorobenzene (BFB) performance check run on instrument F50051 in association with the reported sample analyses. Reported mass abundances were acceptable for this performance check.

Calibration: Form VI was provided for the initial calibration (IC) performed on 7/27/00 which established a calibration range of 0.5-25 $\mu\text{g/L}$ for most target analytes (acetone, 2-butanone, and 4-methyl-2-pentanone were in the standards at five times these concentrations). Reported average relative response factors (RRFs) were below the minimum acceptable value (0.05) for acetone (0.014), 2-butanone (0.004), and 1,2,3-trichloropropane (0.046). The result for 2-butanone in TB was rejected (R) on this basis. Acetone was detected in both samples and 2-butanone was detected in MW-104A; therefore, results for acetone in MW-104A and TB and for 2-butanone in MW-104A were qualified as estimated (J) based on the low average RRFs.

The average RRF for 1,2,3-trichloropropane was only slightly below the minimum acceptable value and an acceptable RRF was reported for this compound in the continuing calibration standard. Therefore, based on professional judgment, results for 1,2,3-trichloropropane were not qualified on this basis.

Percent relative standard deviations (%RSDs) were less than the maximum method-specified acceptance limit (15%) except for dichlorodifluoromethane (16.9%), chloromethane (18.7%), acetone (33.1%), 2-butanone (17.9%), 4-methyl-2-pentanone (18.4%), bromoform (18.5%), 1,2,3-trichloropropane (18.7%), 1,3-dichlorobenzene (20.7%), 1,4-dichlorobenzene (20.1%), 1,2-dichlorobenzene (19.7%), 1,2-dibromo-3-chloropropane (19.9%), 1,2,4-trichlorobenzene (25.7%), hexachlorobutadiene (24.6%), naphthalene (30.9%), and 1,2,3-trichlorobenzene (33.9%).

Mr. Dan Garrigan

Evaluation of CompuChem Case No. Q1458, R1458 (VOCs in GW)

August 31, 2000

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In evaluating this IC, the laboratory exercised an option included in Method 8000B (which contains calibration requirements for Method 8260B) that allows calculation of the mean %RSD value for all of the analytes in the calibration standards. If the mean %RSD is less than 15%, then (1) the IC may still be considered acceptable even if one or more individual %RSDs exceeds 15% and (2) the mean RRFs from IC may be used for calculation of all sample results, rather than having to use regression analyses to generate calibration curves for individual target analytes with %RSDs greater than 15%. This calculation was documented in the data package, and a mean %RSD of 14.5% was reported by the laboratory. This is, however, a contractual/method compliance issue; since the evaluation effort focuses on the individual analyte responses in determining whether or not associated sample results require qualification, the use of the mean %RSD by the laboratory to evaluate the IC has no direct bearing on the data evaluation effort.

For chloromethane, 4-methyl-2-pentanone, bromoform, 1,2,3-trichloropropane, and hexachlorobutadiene, no positive results were reported in the two samples and the %RSDs did not exceed the less-restrictive validation criterion of 30%. Therefore, no qualifiers were applied to the sample results for these compounds on this basis. Results for dichlorodifluoromethane, acetone, 2-butanone, 1,3-dichlorobenzene, 1,4-dichlorobenzene, 1,2-dichlorobenzene, 1,2,4-trichlorobenzene, naphthalene, and 1,2,3-trichlorobenzene in MW-104A and for acetone, 1,3-dichlorobenzene, 1,4-dichlorobenzene, 1,2-dichlorobenzene, 1,2-dibromo-3-chloropropane, 1,2,4-trichlorobenzene, naphthalene, and 1,2,3-trichlorobenzene in TB were qualified as estimated (J, UJ) based on the unacceptable %RSD values.

Form VII was provided for one continuing calibration (CC) standard run in association with these samples. In fact, the samples were run immediately following the IC and this "CC" was the mid-level standard from the IC. RRFs were acceptable (greater than 0.05) in the CC except for acetone (0.012) and 2-butanone (0.004). Results for acetone and 2-butanone were previously qualified based on similar responses in the IC; no additional action was taken based on the CC results.

All reported percent differences (%Ds) from the IC were less than the maximum method-specified acceptance criterion of 20%.

Blanks: Results for one method blank (MB) associated with the site sample analyses were provided in the data report. Acetone (2 µg/L) was reported in this MB. Results for acetone in MW-104A and TB were qualified as less than the reported value or less than the sample-specific reporting limit, whichever was greater (U), based on the associated MB contamination. In both cases, the sample result was less than ten times the MB concentration. All "B" qualifiers appropriately applied by the laboratory were removed by the validator.

One trip blank was submitted with this data set; after qualifications based on MB contamination, methylene chloride (2 µg/L), toluene (0.2 µg/L), 1,3-dichlorobenzene (0.2 µg/L), 1,4-

Mr. Dan Garrigan

Evaluation of CompuChem Case No. Q1458, R1458 (VOCs in GW)

August 31, 2000

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dichlorobenzene (0.2 µg/L), 1,2-dichlorobenzene (0.3 µg/L), n-butylbenzene (0.2 µg/L), 1,2-dibromo-3-chloropropane (0.2 µg/L), 1,2,4-trichlorobenzene (0.4 µg/L), and naphthalene (0.5 µg/L) were reported in the TB. Results for methylene chloride, 1,3-dichlorobenzene, and 1,2,4-trichlorobenzene in MW-104A were qualified as less than the reported value or less than the sample-specific RL, whichever was greater (U), based on the associated TB contamination. The remaining compounds listed above were either not detected in MW-104A or the result was greater than the action limit for qualification (five or ten times the TB concentration) and no further action was warranted on this basis.

System Monitoring Compounds: Reported recoveries (Form II) of the four system monitoring compounds (SMCs) were acceptable for MW-104A, the MB, and the laboratory control sample. Recoveries for three of the four SMCs in TB were too high (131-142%). Results for all detected target analytes in TB were qualified as estimated (J) on this basis.

Matrix Spike/Matrix Spike Duplicate: No matrix spike analysis was performed on MW-104A.

Laboratory Control Sample: Results for a laboratory control sample were reported on a Form III-equivalent in the data package. Acceptable recoveries (102-120%) were reported for the five spiked target analytes listed on this form (1,1-dichloroethene, benzene, trichloroethene, toluene, and chlorobenzene).

Field Duplicate: No field duplicate pair was included in this data package.

Internal Standard Responses: Internal standard (IS) areas and retention times (RTs) were reported on Form VIII for both site samples reported in this data set and the related QC analyses; all areas and RTs were within acceptable limits.

Sample Results: Results for MW-104A and TB were appropriately reported on Form Is, each of which indicated the appropriate analysis date and a sample analysis volume of 25 mL.

Results for chloroethane, 1,1-dichloroethane, benzene, ethylbenzene, m/p-xylenes, o-xylene, 1,2,4-trimethylbenzene, 1,4-dichlorobenzene, naphthalene, and total xylenes in the undiluted analysis of MW-104A exceeded the upper limit of the calibration range and were therefore qualified as estimated (J); the "E" qualifiers appropriately applied by the laboratory to these results were removed by the evaluator. According to the narrative, a diluted rerun analysis was performed by the laboratory but was lost due to an instrument failure. No further analyses were performed because the holding time had expired. Therefore, only the undiluted analysis results are available for MW-104A. The data user is cautioned that the accuracy of the reported values for those compounds exceeding the calibration range cannot be verified. In particular, those results exceeding the upper limit of the calibration range by more than 20% (i.e., values greater than 30 µg/L) should be considered minimum concentrations.

Mr. Dan Garrigan

Evaluation of CompuChem Case No. Q1458, R1458 (VOCs in GW)

August 31, 2000

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Reporting limits for target analytes that were not detected were correctly reported as equivalent to the lowest concentration IC standard analyzed.

Documentation: Data package documentation was inconsistent with the specifications of the QAPP; specifically, a Level IV package was provided when a Level III package was required. The fact that the additional raw data documentation was included certainly does not adversely affect the reported sample results, but, pursuant to the validation specifications for this project, it was not reviewed by Trillium as part of this evaluation.

On the field COC, the date accompanying the laboratory receipt signature was incomplete (no year was included). This COC documentation issue does not directly affect the validity of the reported sample results, but it could be problematic if the data were used in litigation.

Please let me know if you have any questions regarding this data package review report.

Sincerely,



Carol A. Erikson

Quality Assessment Manager

CAE/ekd

July 2000\voacc (95406)

ATTACHMENT A

**ORGANIC ANALYSIS DATA SHEETS (Form I)
Case No. Q1458, SDG R1458
Volatiles in Water**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected. The associated numerical value is the sample quantitation limit. The sample quantitation limit accounts for sample-specific dilution factors and percent solids corrections or sample sizes that deviate from those required by the method.**
- J - The associated numerical value is an estimated quantity.**
- R - The data are unusable (analyte may or may not be present). Resampling and reanalysis is necessary for verification. The R replaces the numerical value or sample quantitation limit.**
- UJ - The analyte was analyzed for, but was not detected. The sample quantitation limit is an estimated quantity.**

FORM 1
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-104A

Lab Name: COMPUCHEM

Contract:

Lab Code: LIBRTY

Case No.:

SAS No.:

SDG No.: R1458

Matrix: (soil/water) WATER

Lab Sample ID: R1458-1

Sample wt/vol: 25 (g/ml) ML

Lab File ID: R1458-1A51

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/27/00

GC Column: EQUITY624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS: (ug/L or ug/Kg) UG/L Q

| | | | |
|----------|--------------------------|-----|----|
| 75-71-8 | Dichlorodifluoromethane | 0.2 | J |
| 74-87-3 | Chloromethane | 0.5 | U |
| 75-01-4 | Vinyl Chloride | 1 | |
| 74-83-9 | Bromomethane | 0.5 | U |
| 75-00-3 | Chloroethane | 44 | J |
| 75-69-4 | Trichlorofluoromethane | 0.5 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.5 | U |
| 67-64-1 | Acetone | 16 | UJ |
| 75-09-2 | Methylene Chloride | 4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 2 | |
| 75-34-3 | 1,1-Dichloroethane | 41 | J |
| 594-20-7 | 2,2-Dichloropropane | 0.5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | 0.9 | |
| 78-93-3 | 2-butanone | 10 | J |
| 74-97-5 | Bromochloromethane | 0.5 | U |
| 67-66-3 | Chloroform | 0.5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.5 | U |
| 56-23-5 | Carbon Tetrachloride | 0.5 | U |
| 563-58-6 | 1,1-dichloropropene | 0.5 | U |
| 71-43-2 | Benzene | 25 | J |
| 107-06-2 | 1,2-Dichloroethane | 9 | |
| 79-01-6 | Trichloroethene | 0.5 | U |
| 78-87-5 | 1,2-Dichloropropane | 10 | |
| 74-95-3 | Dibromomethane | 0.5 | U |
| 75-27-4 | Bromodichloromethane | 0.5 | U |
| 108-10-1 | 4-Methyl-2-pentanone | 3 | U |
| 108-88-3 | Toluene | 12 | |
| 79-00-5 | 1,1,2-Trichloroethane | 0.5 | U |
| 127-18-4 | Tetrachloroethene | 0.5 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.5 | U |
| 124-48-1 | Dibromochloromethane | 0.5 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.5 | U |
| 108-90-7 | Chlorobenzene | 10 | |

ca Erikson 8/31/00

FORM I VOA

FORM 1
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-104A

Lab Name: COMPUCHEM

Contract:

Lab Code: LIBRTY

Case No.:

SAS No.:

SDG No.: R1458

Matrix: (soil/water) WATER

Lab Sample ID: R1458-1

Sample wt/vol: 25 (g/ml) ML

Lab File ID: R1458-1A51

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/27/00

GC Column: EQUITY624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L

CAS NO.

COMPOUND

Q

| | | | |
|-----------|-----------------------------|-----|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.5 | U |
| 100-41-4 | Ethylbenzene | 110 | J |
| 108-38-3 | m,p-Xylene | 130 | J |
| 95-47-6 | o-Xylene | 60 | J |
| 100-42-5 | Styrene | 0.5 | U |
| 75-25-2 | Bromoform | 0.5 | U |
| 98-82-8 | Isopropyl Benzene | 24 | |
| 108-86-1 | Bromobenzene | 0.4 | J |
| 96-18-4 | 1,2,3-Trichloropropane | 0.5 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.5 | U |
| 103-65-1 | n-Propyl Benzene | 5 | |
| 95-49-8 | 2-Chlorotoluene | 0.5 | U |
| 106-43-4 | 4-Chlorotoluene | 0.5 | U |
| 108-67-8 | 1,3,5-Trimethyl Benzene | 13 | |
| 98-06-6 | tert-Butyl Benzene | 0.5 | U |
| 95-63-6 | 1,2,4-Trimethyl Benzene | 40 | J |
| 135-98-8 | sec-Butyl Benzene | 0.5 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.9 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 27 | J |
| 99-87-6 | p-Isopropyl Toluene | 12 | |
| 95-50-1 | 1,2-Dichlorobenzene | 6 | J |
| 104-51-8 | n-Butyl Benzene | 0.5 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.5 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1 | U |
| 87-68-3 | Hexachlorobutadiene | 0.5 | U |
| 91-20-3 | Naphthalene | 58 | J |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.5 | U |
| 540-59-0 | 1,2-Dichloroethene (total) | 3 | |
| 1330-20-7 | Xylene (total) | 200 | J |

20/15/8
M. Erikson 8/31/02

FORM I VOA

FORM 1
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB

Lab Name: COMPUCHEM

Contract:

Lab Code: LIBRTY

Case No.:

SAS No.:

SDG No.: R1458

Matrix: (soil/water) WATER

Lab Sample ID: R1458-2

Sample wt/vol: 25 (g/ml) ML

Lab File ID: R1458-2A51

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/27/00

GC Column: EQUITY624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

| | | | |
|----------|--------------------------|-----|---|
| 75-71-8 | Dichlorodifluoromethane | 0.5 | U |
| 74-87-3 | Chloromethane | 0.5 | U |
| 75-01-4 | Vinyl Chloride | 0.5 | U |
| 74-83-9 | Bromomethane | 0.5 | U |
| 75-00-3 | Chloroethane | 0.5 | U |
| 75-69-4 | Trichlorofluoromethane | 0.5 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.5 | U |
| 67-64-1 | Acetone | 3 | U |
| 75-09-2 | Methylene Chloride | 2 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.5 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.5 | U |
| 594-20-7 | 2,2-Dichloropropane | 0.5 | U |
| 156-59-2 | cis-1,2-Dichloroethene | 0.5 | U |
| 78-93-3 | 2-butanone | 3 | U |
| 74-97-5 | Bromochloromethane | 0.5 | U |
| 67-66-3 | Chloroform | 0.5 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.5 | U |
| 56-23-5 | Carbon Tetrachloride | 0.5 | U |
| 563-58-6 | 1,1-dichloropropene | 0.5 | U |
| 71-43-2 | Benzene | 0.5 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.5 | U |
| 79-01-6 | Trichloroethene | 0.5 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.5 | U |
| 74-95-3 | Dibromomethane | 0.5 | U |
| 75-27-4 | Bromodichloromethane | 0.5 | U |
| 108-10-1 | 4-Methyl-2-pentanone | 3 | U |
| 108-88-3 | Toluene | 0.2 | J |
| 79-00-5 | 1,1,2-Trichloroethane | 0.5 | U |
| 127-18-4 | Tetrachloroethene | 0.5 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.5 | U |
| 124-48-1 | Dibromochloromethane | 0.5 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.5 | U |
| 108-90-7 | Chlorobenzene | 0.5 | U |

3 2 3 4
 H
 R
 CA Erikson 8/31/00

FORM I VOA

FORM 1
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB

Lab Name: COMPUCHEM

Contract:

Lab Code: LIBRTY

Case No.:

SAS No.:

SDG No.: R1458

Matrix: (soil/water) WATER

Lab Sample ID: R1458-2

Sample wt/vol: 25 (g/ml) ML

Lab File ID: R1458-2A51

Level: (low/med) LOW

Date Received: 07/14/00

% Moisture: not dec. _____

Date Analyzed: 07/27/00

GC Column: EQUITY624 ID: 0.53 (mm)

Dilution Factor: 1.0

Soil Extract Volume: _____ (uL)

Soil Aliquot Volume: _____ (uL)

CAS NO. COMPOUND CONCENTRATION UNITS:
(ug/L or ug/Kg) UG/L Q

| | | | |
|----------------|-----------------------------|-----|-----|
| 630-20-6----- | 1,1,1,2-Tetrachloroethane | 0.5 | U |
| 100-41-4----- | Ethylbenzene | 0.5 | U |
| 108-38-3----- | m,p-Xylene | 1 | U |
| 95-47-6----- | o-Xylene | 0.5 | U |
| 100-42-5----- | Styrene | 0.5 | U |
| 75-25-2----- | Bromoform | 0.5 | U |
| 98-82-8----- | Isopropyl Benzene | 0.5 | U |
| 108-86-1----- | Bromobenzene | 0.5 | U |
| 96-18-4----- | 1,2,3-Trichloropropane | 0.5 | U |
| 79-34-5----- | 1,1,2,2-Tetrachloroethane | 0.5 | U |
| 103-65-1----- | n-Propyl Benzene | 0.5 | U |
| 95-49-8----- | 2-Chlorotoluene | 0.5 | U |
| 106-43-4----- | 4-Chlorotoluene | 0.5 | U |
| 108-67-8----- | 1,3,5-Trimethyl Benzene | 0.5 | U |
| 98-06-6----- | tert-Butyl Benzene | 0.5 | U |
| 95-63-6----- | 1,2,4-Trimethyl Benzene | 0.5 | U |
| 135-98-8----- | sec-Butyl Benzene | 0.5 | U |
| 541-73-1----- | 1,3-Dichlorobenzene | 0.2 | J |
| 106-46-7----- | 1,4-Dichlorobenzene | 0.2 | J |
| 99-87-6----- | p-Isopropyl Toluene | 0.5 | U |
| 95-50-1----- | 1,2-Dichlorobenzene | 0.3 | J |
| 104-51-8----- | n-Butyl Benzene | 0.2 | J |
| 96-12-8----- | 1,2-Dibromo-3-Chloropropane | 0.2 | J |
| 120-82-1----- | 1,2,4-Trichlorobenzene | 0.4 | J |
| 87-68-3----- | Hexachlorobutadiene | 0.5 | U |
| 91-20-3----- | Naphthalene | 0.5 | J |
| 87-61-6----- | 1,2,3-Trichlorobenzene | 0.5 | XUJ |
| 540-59-0----- | 1,2-Dichloroethene (total) | 0.5 | U |
| 1330-20-7----- | Xylene (total) | 0.5 | U |

CAE
8/31/00



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September 11, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review Report - CompuChem Case No. Q1458, SDGs R1458 and S1458 - Total and Dissolved Arsenic and Lead and Total Iron in Ground Water

Dear Dan:

I have completed my evaluation of the inorganics analysis data prepared by CompuChem Environmental for one ground water (GW) sample from the L&RR Site, which was reported in two data packages under Case No. Q1458, SDGs R1458 and S1458. The packages were received by Trillium for review on August 11, 2000 and September 6, 2000, respectively. Total and dissolved metals analysis results for sample MW-104A were reported.

Analyses were performed according to EPA SW-846 (Third Edition)-Update III Methods for total and dissolved arsenic, total and dissolved lead, and total iron; all elements were analyzed using trace inductively coupled plasma (ICP) instrumentation (EPA 6010B). My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced methods, and the USEPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses* (6/88; Rev. 2/89). Qualifiers consistent with those defined in the Region I document were applied as necessary and appropriate.

Although the QAPP specifies that Level III data packages are required for all analyses except the 5-year sampling events, full raw data deliverables (i.e., ICP instrument print-outs) were provided for all of the metals analyses. *These raw data were not reviewed by Trillium as part of this evaluation.* As specified in Section 9.2 of the QAPP (page 23), a Level III evaluation, limited to an overview of the information provided on the summary forms, was performed.

Based on the evaluation, results for total and dissolved lead in MW-104A were qualified as estimated (J) based on an unacceptably high recovery in the associated low range standard. All results for arsenic and iron were determined to be valid as reported; no other qualifiers were applied.

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Mr. Dan Garrigan

Evaluation of CompuChem Case No. Q1458, SDGs R1458 and S1458 (Metals)

September 11, 2000

Page 2

These qualifiers are noted on the Inorganic Analysis Data Sheets (Form I) for the reported sample analyses in Attachment A to this report; Region I qualifier definitions are also provided in this attachment.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation and Sample Integrity: A copy of the applicable field chain of custody (COC) record was included in the data package for SDG R1458 but was not found in the data package for SDG S1458; a copy of the COC from R1458 was made by the evaluator and inserted into S1458 as page 175a to ensure that complete documentation is available for future reference.

The site sample was collected on 7/13/00. Total and dissolved metals analyses were performed on 7/29/00, well within the 6-month holding time specified by the QAPP.

An acceptable ($4^{\circ}\text{C} \pm 2^{\circ}\text{C}$) cooler temperature (3°C) on laboratory receipt was recorded on the COC. The "Filtered/Unfiltered" column on the COC indicated that both filtered and unfiltered sample containers were provided for analysis, and acidification with nitric acid was also noted by the sampler on the COC. Verification of successful acidification was not recorded by the laboratory on the COC, but was found ($\text{pH} < 2$) for both the total and dissolved metals samples on the laboratory receiving log and on the laboratory preparation logs.

Results from analysis of the dissolved metals sample fraction were distinguished from analysis results for the total metals fraction only by reporting the samples in separate SDGs and data packages. The word "Dissolved" was added by the evaluator to the dissolved metals analysis results form (Form I) attached to this report to clarify the distinction; similarly, the word "Total" was added to the total metals analysis results form.

Calibrations: Based on the Analytical Run Logs (Contract Laboratory Program, CLP, Form XIV) in both data packages, initial and continuing calibration verification (ICV/CCV) standards were run at the appropriate frequencies throughout the single analysis series during which MW-104A was run for both total and dissolved constituents. All ICV/CCV recoveries documented on the accompanying CLP Form IIAs were acceptable (QC 90-110%).

Results for a "low range standard (LRS)," containing arsenic at $10\ \mu\text{g/L}$ and lead at $3\ \mu\text{g/L}$, were reported on Form IIB in each data package; iron (at $100\ \mu\text{g/L}$) was also reported in SDG R1458. Recoveries for arsenic (82.5%) and iron (99.9%) were acceptable (QC 80-120%), but an unacceptably high recovery was reported for lead (125.0%). Since the recovery for lead was too high, suggesting a high bias at low concentrations, and the detected total and dissolved lead concentrations were less than four times the reporting limit (RL) of $3\ \mu\text{g/L}$, the results for total and dissolved lead in MW-104A were qualified as estimated (J).

Mr. Dan Garrigan

Evaluation of CompuChem Case No. Q1458, SDGs R1458 and S1458 (Metals)

September 11, 2000

Page 3

Blanks: Initial and continuing calibration blanks (ICB/CCBs) were run at the appropriate frequencies throughout the sample analysis series. Arsenic, iron, and lead were not reported above the applicable IDLs or below the negative IDLs in any of the ICB/CCBs.

Two water matrix preparation blanks (PBW) were analyzed with the site sample (CLP Form XIII), one each with the total and dissolved fractions; no arsenic or lead was reported above the applicable IDLs (or below the negative IDLs) in either PBW, but iron (47.510 µg/L) was reported in the PBW associated with SDG R1458. The sample result for iron was much greater than the action limit for qualification based on the blank contamination (five times the blank concentration), therefore no action was warranted on this basis.

Interference Check Sample: Recoveries reported on CLP Form IV in each data package for all target elements in the interference check sample analyses performed at the start of the analytical series were acceptable (QC 80-120%).

Laboratory Control Samples (LCS): Acceptable recoveries (QC 80-120%) for total arsenic, iron, and lead were reported on the Laboratory Control Sample Report (CLP Form VII) in the data package for SDG R1458 (101.8-104.2%). Acceptable recoveries for dissolved arsenic and lead were reported on the LCS Report in the data package for SDG S1458 (100.8% and 102.3%).

Duplicate Analysis: MW-104A was run as an unspiked duplicate pair for dissolved arsenic and lead. Excellent reproducibility was demonstrated for arsenic (0.2 relative percent difference [RPD]). The RPD for lead (21.3%) slightly exceeded the method-specified acceptance criterion of 20%, and the result was qualified "N" by the laboratory. Since the concentration of dissolved lead in the sample was less than the CRDL and the paired results agreed within ±CRDL, no qualifiers were warranted on this basis and the "N" qualifier was removed.

Matrix Spike Analysis: MW-104A was prepared and analyzed as a matrix spike/matrix spike duplicate pair for dissolved arsenic and lead. Recoveries (93.5-103.7%) were acceptable (QC 75-125%) and reproducible (RPDs 1.5% and 2.6%).

ICP Serial Dilution: Serial dilution analysis was performed on MW-104A for dissolved arsenic and lead. Both target analytes were less than 50xIDL in the undiluted sample analysis, therefore no meaningful information was obtained from this quality control analysis and no further evaluation was made.

Sample Results: Results greater than or equal to the laboratory-specified IDLs were appropriately reported on the Inorganic Analysis Data Sheets (Form I) for the individual sample analytes.

Form X was provided in both data packages, documenting IDLs established on 7/15/00. All non-detected blank results were reported to the appropriate IDLs, and all sample analyses were documented on the run logs (Form XIV) in the two data packages.

Mr. Dan Garrigan

Evaluation of CompuChem Case No. Q1458, SDGs R1458 and S1458 (Metals)

September 11, 2000

Page 4

Field Duplicates: No field duplicate sample was submitted with this data set.

Total versus Dissolved Concentrations: Comparison of total versus dissolved results for arsenic and lead in MW-102A revealed no cases where the dissolved concentration was greater than the total concentration.

Documentation: Data package documentation was inconsistent with the specifications of the QAPP; specifically, a Level IV package was provided when a Level III package was required. The presence of this additional raw data documentation certainly does not adversely affect the reported sample results, but, pursuant to the validation specifications for this project, it was not reviewed by Trillium as part of this evaluation.

The field COC record was not included in the data package for SDG S1458. A copy of the COC record found in SDG R1458 was made by the evaluator and inserted into S1458 as page 175a to ensure that complete documentation is available for future reference.

On the field COC record, the date accompanying the laboratory receipt signature is incomplete; no year is included. In addition, sample pHs on laboratory receipt were not recorded on the COC and the project-specific analytes for analysis were not specified by the sampler on the COC.

Total iron results were not reported in the data package for SDG R1458 as originally received for review. At the request of the evaluator, the laboratory provided revised summary forms (Forms I, IIA, III, IV, VII, X, XII, and XIV and the LRS Summary Form) on 9/6/00 via FedEx. These forms were page-numbered and inserted into the R1458 data package by the evaluator, *replacing* the pages originally provided by the laboratory.

Although the laboratory was sure they had sent it, the data package for SDG S1458 was not received by Trillium with the rest of the data packages associated with this data set. At the evaluator's request, a second copy of this data package (which was, in fact, dated August 4, 2000) was received by Trillium on September 6, 2000.

Please let me know if you have any questions regarding this report.

Sincerely,



Carol A. Erikson
Quality Assessment Manager

CAE/esc

July 2000\Metalscc (95406)

ATTACHMENT A

**INORGANIC ANALYSIS DATA SHEETS
CompuChem Case No. Q1458, SDGs R1458 and S1458**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.**
- J - The associated numerical value is an estimated quantity.**
- R - The data are unusable (Note: Analyte may or may not be present).**
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.**

SW-846 METALS

1

INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW-104A

TOTAL

CAE 9/8/00

Lab Name: CompuChem

Contract: _____

Lab Code: LIBRTY

Case No.: Q1458

SAS No.: _____

SDG No.: R1458

Matrix (soil/water): WATER

Lab Sample ID: R1458-1

Level (low/med): LOW

Date Received: 07/14/00

Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|---------|---------------|---|---|---|
| 7440-38-2 | Arsenic | 113 | | | P |
| 7439-89-6 | Iron | 99200 | | | P |
| 7439-92-1 | Lead | 11.3 | | J | P |

CAE 9/8/00

Color Before: BROWN

Clarity Before: CLOUDY

Texture: _____

Color After: BROWN

Clarity After: CLOUDY

Artifacts: _____

Comments: _____

SW-846 METALS

1

INORGANIC ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW-104A

DISSOLVED

CAE 9/8/00

Lab Name: CompuChem

Contract: _____

Lab Code: LIBRTY

Case No.: Q1458

SAS No.: _____

SDG No.: S1458

Matrix (soil/water): WATER

Lab Sample ID: S1458-1

Level (low/med): LOW

Date Received: 07/14/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|---------|---------------|--------------|--------------|-----|
| 7440-38-2 | Arsenic | 113 | | | P |
| 7439-92-1 | Lead | 2.3 | B | X | J P |

CAE 9/8/00

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments: _____



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September 1, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review Report - CompuChem Case No. R1458 - Chloride and Ammonia in Ground Water and TestAmerica Project No. 200001 - BOD and COD in Ground Water

Dear Dan:

I have completed my evaluation of the wet chemistry analysis data prepared by CompuChem Environmental (chloride and ammonia) and TestAmerica (BOD and COD) for one ground water (GW) sample from the L&RR Site. These data were reported in two data packages under CompuChem Case R1458 and TestAmerica Project No. 200001, which were received by Trillium for review on August 11, 2000. Results for sample MW-104A were reported in both data packages.

Analyses were performed according to EPA Methods 325.2 (chloride), 410.4 (COD), 350.1 (ammonia), and 405.1 (BOD); all four methods are found in *Methods for Chemical Analysis of Water and Wastes* (EPA-600/4-79-020, 3/83). My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced methods, and, to the extent applicable, the USEPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses* (6/88; Rev. 2/89). Qualifiers consistent with those defined in the Region I document were applied as necessary and appropriate.

Based on the evaluation, the result for BOD in MW-104A was qualified as estimated (J) because incubation was initiated past the specified holding time. No additional qualifiers were applied; the remaining sample results were determined to be valid as reported based on the available data reports.

This qualifier is reflected on the affected laboratory reporting form included as Attachment A to this report. Region I qualifier definitions are also provided in this attachment.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation and Sample Integrity: A copy of the field chain of custody (COC) record was included in the CompuChem data package and a copy of the COC record transferring the BOD and COD samples to TestAmerica was included in the TestAmerica report. Both COCs documented a sample collection date of 7/13/00. An acceptable cooler temperature (3°C) was documented on the field COC by CompuChem on receipt of the samples. No cooler temperature was recorded by TestAmerica on the subcontract COC,

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Mr. Dan Garrigan

Evaluation CompuChem Case No. R1458 and TestAmerica Project No. 200001 (Wet Chemistry)

September 1, 2000

Page 2 of 3

however an acceptable temperature of 2 °C was recorded on their cooler receipt form, which was also provided for review.

Chemical preservation of the sample containers for ammonia and COD analysis using sulfuric acid was documented on both COC records. Sample pH on laboratory receipt was not recorded on the COCs by either laboratory, but CompuChem's receiving log indicated that both sample containers were at pH < 2 and TestAmerica's cooler receipt form noted that "correct preservatives" were used. It was therefore assumed that the sample was properly preserved at the time of collection.

Analyses for chloride were performed on 8/3/00, ammonia analyses were performed on 7/29/00, and COD analyses were run on 7/19/00; these are all within the 28-day holding time specified by the QAPP. BOD incubation was initiated on 7/15/00 at 22:00, which is 10 hours past the required 48-hour holding time. The reported result for BOD in MW-104A was qualified as estimated (J) on this basis. Five-day BOD readings were appropriately taken on 7/20/00.

Calibrations: Based on the run log provided in the CompuChem data package, initial and continuing calibration verification (ICV/CCV) standards were run at the appropriate frequencies throughout the analysis series for ammonia. All ICV/CCV recoveries for this analysis parameter were acceptable (QC 85-115%). Summary results for an eight-point calibration curve run on 7/29/00 were also found in the data package; an acceptable correlation coefficient of 0.9998 (QC \geq 0.995) was reported for the best-fit quadratic equation.

Based on the run log provided in the CompuChem data package, ICV/CCV standards were run at the appropriate frequencies throughout the analysis series for chloride. All ICV/CCV recoveries for this analysis parameter were acceptable (QC 85-115%). Summary results for a nine-point calibration curve run on 8/3/00 were also found in the data package; an acceptable correlation coefficient of 0.9991 was reported for the best-fit quadratic equation.

Based on the run log provided in the TestAmerica data package, ICV/CCV standards were run at the appropriate frequencies throughout the analysis series for COD. All ICV/CCV recoveries for this analysis parameter were acceptable (QC 85-115%). Summary results for a six-point calibration curve run on 7/19/00 were also found in the data package; an acceptable correlation coefficient of 0.9997 was reported for the best-fit linear regression.

Calibration, per se, is not applicable to BOD measurements.

Blanks: Initial and continuing calibration blanks (ICB/CCBs) were run at the appropriate frequencies throughout the ammonia, chloride, and COD analysis series; no target analytes were detected above the reporting limit (RL) in any of these blanks.

A water matrix preparation blank (PBW) was also prepared with the sample for each parameter; no ammonia, chloride, BOD, or COD was reported in the PBW analyses.



Mr. Dan Garrigan

Evaluation CompuChem Case No. R1458 and TestAmerica Project No. 200001 (Wet Chemistry)

September 1, 2000

Page 3 of 3

Laboratory Control Samples: Acceptable recoveries (90.6-103%) for all four analysis parameters were reported for the laboratory control samples run in association with the site sample.

Duplicate Analysis: MW-104A was run in duplicate for COD, and excellent reproducibility (relative percent difference 1.7%) was demonstrated.

No laboratory duplicate analyses were reported for ammonia or chloride. Duplicate analysis results using a non-project sample were reported for BOD, but these data are not relevant to MW-104A and were not reviewed.

Matrix Spike Analysis: MW-104A was not prepared and analyzed as a matrix spike for any of the four wet chemistry analysis parameters.

Spike recovery results for a non-project sample were reported for COD, but these data are not relevant to MW-104A and were not reviewed.

Sample Results: Sample results for each analysis parameter were correctly transcribed from the laboratory run logs and appropriately reported on summary forms.

Field Duplicates: No field duplicate was submitted with this site sample.

Documentation: Data package documentation was complete and acceptable for all four analysis parameters. On the field COC, the date accompanying the laboratory receipt signature was incomplete (no year was included). On the subcontract COC, the dates accompanying both "Relinquished by" signatures are incomplete (no year is included for either date) and there is no signature indicating receipt at TestAmerica in Nashville, TN, where the analyses were performed. These COC documentation issues do not directly affect the validity of the reported sample results, but they could be problematic if the data were used in litigation.

Please let me know if you have any questions regarding this data package review report.

Sincerely,

Carol A. Erikson

Quality Assessment Manager

CAE/das

July 2000\wchemcc (95406)

ATTACHMENT A

LABORATORY REPORTING FORMS
CompuChem Case No. R1458 and TestAmerica Project No. 200001

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.**
- J - The associated numerical value is an estimated quantity.**
- R - The data are unusable (Note: Analyte may or may not be present).**
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.**

AMMONIA ANALYSIS

SUMMARY REPORT

| ITEM NO. | SAMPLE IDENTIFIER | COMPUCHEM NUMBER | RESULT (mg/L) | REPORTING LIMIT (mg/L) |
|----------|-------------------|------------------|---------------|------------------------|
| 1. | MW-104A | R1458-1 | 15.8 | 0.1 |

BRL = BELOW REPORTING LIMIT

Reviewed by/ID#: W. George 12405 Date: 8/3/00

CHLORIDE ANALYSIS

SUMMARY REPORT

| ITEM NO. | SAMPLE IDENTIFIER | COMPUCHEM NUMBER | RESULT (mg/L) | REPORTING LIMIT (mg/L) |
|----------|-------------------|------------------|---------------|------------------------|
| 1. | MW-104A | R1458-1 | 386 | 3 |

BRL = BELOW REPORTING LIMIT

Reviewed by/ID#: PW George 12405 Date: 8/3/00

5
RUS
8/11/00

Test America

INCORPORATED

ANALYTICAL REPORT

COMPUCHEM 2303
 DIANE BYRD
 501 MADISON AVENUE
 CARY, NC 27513

Lab Number: 00-A99172
 Sample ID: MW-104A
 Sample Type: Water
 Site ID:

Project:
 Project Name: L&RR
 Sampler:

Date Collected: 7/13/00
 Time Collected: 12:50
 Date Received: 7/15/00
 Time Received: 9:00

| Analyte | Result | Units | Report Limit | Quan Limit | Dil Factor | Analysis Date | Analysis Time | Analyst | Method | Batch |
|---------------------------|--------|----------|--------------|------------|------------|---------------|---------------|------------|--------|-------|
| BOD Set Up | | | | | | 7/15/00 | 22:00 | | | |
| BOD 5 Day | 27.0 | J mg/l | 2.0 | 2.0 | 1 | 7/20/00 | 20:30 | M.Shockley | 405.1 | 6571 |
| *MISCELLANEOUS CHEMISTRY* | | | | | | | | | | |
| pH | 6.50 | pH Units | | | 1 | 7/18/00 | 11:40 | K. Stewart | 150.1 | 6405 |
| Chemical Oxygen Demand | 118. | mg/l | 15.0 | 3.00 | 5 | 7/19/00 | 15:32 | K.Kenney | 410.4 | 8294 |
| | | | | | | 7/23/00 | | | | |

BOD: Analyzed out of hold

ND - Not detected at the report limit.

Sample report continued . . .

TABLE 2
DATA QUALIFIER DEFINITIONS BY MATRIX

GROUNDWATER AND SURFACE WATER

Volatile Organics in Groundwater and Surface Waters
STL Job Nos. 20980-84697, 84717 and 84719

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected. The associated numerical value is the sample quantitation limit. The sample quantitation limit accounts for the sample-specific dilution factors and percent solids corrections or sample sizes that deviate from those required by the method.
 - J - The associated numerical value is an estimated quantity.
 - R - The data are unusable (analyte may or may not be present). Resampling and reanalysis is necessary for verification. The R replaces the numerical value or sample quantitation limit.
 - UJ - The material was analyzed for, but was not detected. The sample quantitation limit is an estimated quantity.
-

TOTAL AND DISSOLVED METALS IN GROUNDWATER AND SURFACE WATERS

STL Job Nos. 20980-84697, 84717 and 84719

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.
 - J - The associated value is an estimated quantity.
 - R - The data are unusable (Note: Analyte may or may not be present).
 - UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.
-

AMMONIA, CHLORIDE, BOD AND COD IN GROUNDWATER AND SURFACE WATERS

Laboratory Job No. 083967-10

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.
- J - The associated numerical value is an estimated quantity.
- R - The data are unusable (Note: Analyte may or may not be present).
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.

TABLE 2 Continued
DATA QUALIFIER DEFINITIONS BY MATRIX

GROUNDWATER
Compuchem Laboratories
Laboratory Job No. 33003
Volatile Organics in Groundwater

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected. The associated numerical value is the sample quantitation limit. The sample quantitation limit accounts for the sample-specific dilution factors and percent solids corrections or sample sizes that deviate from those required by the method.
 - J - The associated numerical value is an estimated quantity.
 - R - The data are unusable (compound may or may not be present). Resampling and reanalysis is necessary for verification. The R replaces the numerical value or the sample quantitation limit.
 - UJ - The material was analyzed for, but was not detected. The sample quantitation limit is an estimated quantity.
-

TOTAL AND DISSOLVED METALS IN GROUNDWATER

Compuchem Laboratories
Laboratory Job No. 33003
Inorganic Analysis in Groundwater

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.
 - J - The associated numerical value is an estimated quantity.
 - R - The data are unusable (Note: Analyte may or may not be present).
 - UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.
-

CHLORIDE, BOD, COD AND AMMONIA IN GROUNDWATER

Compuchem Laboratories
Wet Chemistry In Groundwater
Laboratory Job No. 33003/00017

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.
- J - The associated numerical value is an estimated quantity.
- R - The data are unusable (Note: Analyte may or may not be present).
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.

Appendix E

**Surface Water Data Validation
Reports
Severn Trent Laboratories**



TRILLIUM INC.
Consultants in Environmental Chemistry

356 FARRAGUT CROSSING DR.
KNOXVILLE, TN 37922
(423) 966-8880
FAX (423) 966-8885
cerikson@trilliuminc.com

June 13, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review Report - STL Lot No. 9A04G680 - Total and Dissolved Arsenic in Surface Water

Dear Dan:

I have completed my evaluation of the inorganics analysis data prepared by the Severn Trent Laboratory (STL) in University Park, Illinois, for seven surface water (SW) samples from the L&RR Site, which were reported in a single data package under Lot No. 9A04G680. The data package was received by Trillium for review on April 28, 2000, and the following samples were reported:

| | | |
|-------|-------|-------|
| LCH-5 | LCH-4 | SW-16 |
| SW-10 | SW-8 | SW-5 |
| DUP | | |

Analyses were performed according to EPA Method 7060A (total and dissolved arsenic), which is found in the Third Edition of SW-846. My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced method, and the USEPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses* (6/88; Rev. 2/89); professional judgment was applied as necessary. Qualifiers consistent with those defined in the Region I document were applied where appropriate.

As required by the QAPP, a Level III data package was provided for these metals analyses. As specified in Section 9.2 of the QAPP (page 23), a Level III evaluation, limited to an overview of the information provided on the summary forms, was performed.

Results from analysis of the dissolved metals sample fractions were distinguished from analysis results for the total metals fractions by the addition of the suffix "S" (for "soluble") to the client sample identification (SW-8S, for example). The word "Dissolved" was also added by the evaluator

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Mr. Dan Garrigan

Evaluation of STL Lot No. 9A04G680 (Total and Dissolved Arsenic in SW)

June 13, 2000

Page 2

to the results forms for the dissolved sample analyses to emphasize the distinction, and for consistency with previous data sets generated in support of this project.

Based on the evaluation, the following qualifiers were applied:

- Results for total and dissolved arsenic in LCH-4 and DUP were qualified as estimated (J) due to poor reproducibility in the field duplicate analyses. Total and dissolved arsenic in DUP were similarly qualified due to a severe discrepancy between the measured total and dissolved concentrations and professional judgment. The result for dissolved arsenic in DUP was also estimated based on an unacceptably low post-digestion spike recovery.
- Results for total and dissolved arsenic in SW-5 were qualified as estimated (J, UJ) due to a minor discrepancy between the measured total and dissolved concentrations.

These qualifiers are reflected on the Inorganic Analysis Data Sheets (Form Is) included as Attachment A to this report. Region I qualifier definitions are also provided in this attachment.

At the discretion of the data user, the following points may warrant attention by laboratory personnel with regard to this data package and/or prior to future sampling events at this site:

- Documentation of sample preservation performed in the field on the chain of custody record.
- Documentation of field filtration for dissolved metals analysis on the chain of custody record.
- Documentation of cooler temperatures and sample pHs at the time of laboratory receipt on the chain of custody record.
- Documentation of linearity check data (correlation coefficient, at a minimum) for all analytes run by GFAA techniques.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation and Sample Integrity: A copy of the field chain of custody (COC) record was included in the data package, documenting a sample collection date of 4/7/00 for the surface waters. Acceptable cooler temperatures (3.6°C and 2.6°C) on laboratory receipt were

Mr. Dan Garrigan

Evaluation of STL Lot No. 9A04G680 (Total and Dissolved Arsenic in SW)

June 13, 2000

Page 3

recorded on the COC. Acidification of the samples with nitric acid, however, was not noted by the sampler on the COC and no verification of acidification was documented in the data package by the laboratory. At the request of the evaluator, the laboratory provided a revised COC for the surface water samples, with a checkmark in the "Preserved" space added. In addition, "pH<2" was noted for all samples on the laboratory digestion logs for arsenic, which were also provided by the laboratory at the request of the evaluator. These records verify successful preservation of the samples for metals analysis.

The revised COC was inserted into the data package for G680 as page 199a and the laboratory digestion logs were inserted as pages 82a and 82b by the evaluator.

Field-filtration of the samples intended for dissolved metals analysis was also not recorded on the COC. For the purposes of this review, it was assumed that the samples *were* field-filtered for dissolved metals analyses (prior to acidification) based on procedures followed during previous sampling events.

Metals analyses were performed 4/14/00 through 4/24/00, well within the 6-month holding time specified by the QAPP.

Calibrations: Based on the run logs (Form XIV), initial and continuing calibration verification (ICV/CCV) standards were run at the appropriate frequencies throughout the three reported GFAA (graphite furnace atomic absorption) analysis series. All ICV/CCV recoveries documented on the accompanying Form II were acceptable (QC 90-110%).

Analysis of linearity check standards for arsenic was evident from the run log entries, however linearity results were not documented in the data package. Therefore, no evaluation of this portion of the data could be made. At the discretion of the data user, the laboratory should be requested to provide this information (at a minimum, the applicable correlation coefficients) for inclusion in the data package, to ensure that complete documentation is available for future reference. For the purposes of this validation effort, it was assumed that acceptable correlation coefficients (>0.995) were achieved for arsenic prior to the reported sample analyses in each series.

Contract required detection limit (CRDL) standards, containing arsenic at 10 µg/L, were also run at the start of each analysis series. Acceptable (QC 80-120%) recoveries were reported for all three reported CRDL standards (100.7-115.7%).

Blanks: Initial and continuing calibration blanks (ICB/CCBs) were also run at the appropriate frequencies throughout each sample analysis series. Per Form III, no arsenic was detected above the instrument detection limit (IDL) in any of the ICB/CCBs. Arsenic was reported at a concentration

Mr. Dan Garrigan

Evaluation of STL Lot No. 9A04G680 (Total and Dissolved Arsenic in SW)

June 13, 2000

Page 4

below the negative IDL in one CCB (-1.9 µg/L) run on 4/24/00. No sample results were affected by this negative response.

A water matrix preparation blank (PBW) was digested with the samples (per the preparation log); no arsenic was detected in the PBW.

Laboratory Control Samples: An acceptable recovery for arsenic (97.5%) was reported on Form VII in the data package.

Duplicate Analysis: Sample LCH-5 was analyzed in duplicate for total arsenic. A slightly high relative percent difference (RPD) of 20.2% (QC ≤20%) was reported on Form VI for the paired results, but the arsenic concentrations in both analyses were less than five times the CRDL specified on Form X and the paired values were well within ±CRDL. Therefore, no action was warranted on this basis.

Matrix Spike Analysis: Sample LCH-5 was prepared and analyzed as a matrix spike/matrix spike duplicate (MS/MSD) pair for total arsenic. Acceptable percent recoveries (96.4% and 94.2%) were reported on Form V for these analyses (QC 75-125%), which also demonstrated excellent reproducibility (2.3 RPD).

No separate spiked analysis was done for dissolved arsenic.

Graphite Furnace Atomic Absorption (GFAA) QC: Post-digestion spike (PDS) recoveries for arsenic were reported on the analysis run logs and were acceptable (QC 85-115%) with the exception of arsenic in DUPS (51.3%), despite a 1:4 dilution performed by the laboratory. The result for arsenic in DUPS was qualified as estimated (J) on this basis.

Unacceptable PDS recoveries were also obtained for several samples that were re-analyzed at various dilutions. In these cases, the diluted analyses gave acceptable PDS recoveries and the diluted analysis results were reported, therefore no further action was warranted on this basis.

Sample Results: Results greater than or equal to the laboratory-specified IDL were appropriately reported on the Inorganic Analysis Data Sheets (Form Is) for total and dissolved arsenic in each sample.

CLP Form X was provided in the data package, indicating an IDL of 1.7 µg/L established for arsenic on 4/15/00.

Mr. Dan Garrigan

Evaluation of STL Lot No. 9A04G680 (Total and Dissolved Arsenic in SW)

June 13, 2000

Page 5

All undetected sample results for arsenic were correctly adjusted (multiplied by 0.5) to reflect the concentration factor resulting from processing 100 mL of sample to a final volume of 50 mL, which was clearly documented on the laboratory digestion logs. Positive sample results cannot be similarly verified without the supporting raw data.

IDLs for undetected analytes were also appropriately adjusted to reflect sample dilutions, where applicable. Positive sample results cannot be similarly verified without the supporting raw data.

Field Duplicates: Sample DUP was identified as a field duplicate of LCH-4 based on information provided by the laboratory. Paired results for total arsenic (69.3 RPD) and dissolved arsenic (44.4 RPD) exceeded the QAPP-specified acceptance limit of 30%. Although the arsenic concentrations in both samples were less than five times the CRDL specified on Form X, the paired values were also outside the alternate criterion for low concentrations (\pm CRDL). Therefore, results for total and dissolved arsenic in LCH-4 and DUP were qualified as estimated (J) on this basis.

Total versus Dissolved Concentrations: Comparison of validated total versus validated dissolved results for arsenic in the seven surface water samples revealed the following cases where the dissolved concentration was higher than the total concentration:

| | | ug/L | | |
|--------|----|--------|-----------|------|
| Sample | | Total | Dissolved | %D |
| SW-5 | As | 0.85 U | 1.1 | -- |
| DUP | As | 3.2 J | 8.8 J | 175% |

For SW-5, the positive dissolved arsenic result is only slightly greater than the reporting limit, and these results most likely reflect the increased variability inherent at concentrations at or near the lower measurement limit. Therefore, results for total and dissolved arsenic in SW-5 were qualified as estimated (J, UJ) on this basis.

For DUP, the percent difference (%D) is unusually high. Validation guidelines recommend that results for which the %D exceeds 50% be rejected, however several additional pieces of data are available in this case that allow a less drastic conclusion. DUP is a field duplicate of LCH-4. Although the paired results for total and dissolved arsenic in these two samples showed poor reproducibility, the total versus dissolved comparison for LCH-4 was acceptable. In addition, post-digestion spike recoveries were acceptable for analyses of total and dissolved arsenic in LCH-4 and for total arsenic in DUP; only the post-digestion spike recovery for (dissolved) arsenic in DUPs was unacceptable (too low). Finally, all four of the arsenic results for these two samples are less than ten times the



Mr. Dan Garrigan

Evaluation of STL Lot No. 9A04G680 (Total and Dissolved Arsenic in SW)

June 13, 2000

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CRDL reported on Form X in the data package. It may be that the discrepancies simply reflect the increased variability typically observed at low concentrations. Or, the low post-digestion spike recovery may suggest that the reported concentration for dissolved arsenic in DUPS is incorrectly biased high. Or, it is possible that one or more of the results is incorrectly reported, which cannot be evaluated in this level review. Therefore, based on the generally low analyte concentrations, review of all the available data, and professional judgment, results for total and dissolved arsenic in DUP were not rejected but were qualified as estimated (J).

Documentation: Data package documentation was consistent with the specifications of the QAPP; specifically, a Level III package was provided, as required. However, summary forms documenting linearity checks (initial multipoint calibration correlation coefficients) for GFAA analytes were missing from the data packages provided for review. At the discretion of the data user, the laboratory may be requested to provide the missing linearity information to ensure that accurate and complete documentation is available for future reference.

Documentation of sample pHs on laboratory receipt was incomplete in the data package as received for review. Clarifications and corrections, as appropriate, were verbally requested by the validator on 6/6/00; responses were received from the laboratory via FedEx on 6/9/00. Corrected documents (revised COC and digestion logs) were inserted into the data package as previously discussed.

A revised COC for the surface water samples, with a checkmark in the "Preserved" space added, was provided by the laboratory, but this new entry was not initialed or dated. While this has no direct effect on the technical validity of the reported sample data, it could be problematic if the data are used for litigation.

Please let me know if you have any questions regarding this data package review report.

Sincerely,

Carol A. Erikson

Quality Assessment Manager

CAE/ekd

95406\Apr00\swmetals

ATTACHMENT A

**INORGANIC ANALYSIS DATA SHEETS (FORM Is)
Laboratory Lot No. 9A04G680**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.**
- J - The associated numerical value is an estimated quantity.**
- R - The data are unusable (Note: Analyte may or may not be present).**
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.**

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

LCH-5

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U04680

Matrix (soil/water): WATER Lab Sample ID: 9A04G680-001

Level (low/med): LOW Date Received: 04/08/00

Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 4.4 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: YELLOW Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

LCH-5S
DISSOLVED
CSC 6/13/00

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____

SDG No.: U04680

Matrix (soil/water): WATER

Lab Sample ID: 9A04G680-002

Level (low/med): LOW

Date Received: 04/08/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.85 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
SOLUBLE

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

LCH-4

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____ SDG No.: U04680

Matrix (soil/water): WATER Lab Sample ID: 9A04G680-003

Level (low/med): LOW _____ Date Received: 04/08/00

% Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 6.6 | | J | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Case 6/13/00

Color Before: COLORLESS Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS Clarity After: CLEAR _____ Artifacts: _____

Comments:

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

LCH-4S
DISSOLVED

CASE 6/13/00

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____

SDG No.: U04680

Matrix (soil/water): WATER

Lab Sample ID: 9A04G680-004

Level (low/med): LOW _____

Date Received: 04/08/00

Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L _____

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 5.6 | | J | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

CASE 6/13/00

Color Before: COLORLESS Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS Clarity After: CLEAR _____ Artifacts: _____

Comments:

SOLUBLE _____

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

SW-16

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U04680

Matrix (soil/water): WATER Lab Sample ID: 9A04G680-005

Level (low/med): LOW Date Received: 04/08/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.85 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

SW-16S
DISSOLVED
CAE 6/13/00

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____ SDG No.: U04680

Matrix (soil/water): WATER _____ Lab Sample ID: 9A04G680-006

Level (low/med): LOW _____ Date Received: 04/08/00

Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.85 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS _____ Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS _____ Clarity After: CLEAR _____ Artifacts: _____

Comments:
SOLUBLE _____

INORGANIC ANALYSES DATA SHEET

SW-10

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____ SDG No.: U04680

Matrix (soil/water): WATER _____ Lab Sample ID: 9A04G680-007

Level (low/med): LOW _____ Date Received: 04/08/00

% Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.85 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS _____ Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS _____ Clarity After: CLEAR _____ Artifacts: _____

Comments:

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

SW-10S
DISSOLVED

02E 6/13/00

Lab Name: STL_CHICAGO

Contract: _____

Lab Code: STL

Case No.: _____

SAS No.: _____

SDG No.: U04680

Matrix (soil/water): WATER

Lab Sample ID: 9A04G680-008

Level (low/med): LOW

Date Received: 04/08/00

Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.85 | U | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS

Clarity Before: CLEAR

Texture: _____

Color After: COLORLESS

Clarity After: CLEAR

Artifacts: _____

Comments:
SOLUBLE

INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

SW-8

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____ SDG No.: U04680

Matrix (soil/water): WATER Lab Sample ID: 9A04G680-009

Level (low/med): LOW _____ Date Received: 04/08/00

Solids: _____ 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 23.2 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: YELLOW _____ Clarity Before: CLEAR _____ Texture: _____

Color After: COLORLESS _____ Clarity After: CLEAR _____ Artifacts: _____

Comments:

INORGANIC ANALYSES DATA SHEET

SW-8S
DISSOLVED

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____

CAE 6/13/00
SDG No.: U04680

Matrix (soil/water): WATER

Lab Sample ID: 9A04G680-010

Level (low/med): LOW

Date Received: 04/08/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 11.0 | | | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

SOLUBLE

U.S. EPA - CLP

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

SW-5

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: SAS No.: SDG No.: U04680

Matrix (soil/water): WATER Lab Sample ID: 9A04G680-011

Level (low/med): LOW Date Received: 04/08/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|----|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 0.85 | ✓ | UJ | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

0226/13/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

INORGANIC ANALYSES DATA SHEET

SW-5S
DISSOLVED

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U04680 *COE 6/13/00*

Matrix (soil/water): WATER Lab Sample ID: 9A04G680-012

Level (low/med): LOW Date Received: 04/08/00

Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 1.1 | | J | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

COE 6/13/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
SOLUBLE

1
INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

DUP

Lab Name: STL_CHICAGO Contract: _____

Lab Code: STL Case No.: _____ SAS No.: _____ SDG No.: U04680

Matrix (soil/water): WATER Lab Sample ID: 9A04G680-013

Level (low/med): LOW Date Received: 04/08/00

% Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 3.2 | | J | F |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

04/08/00

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:

INORGANIC ANALYSES DATA SHEET

DUPS
DISSOLVED

CAE 6/13/00

Lab Name: STL_CHICAGO _____ Contract: _____

Lab Code: STL _____ Case No.: _____ SAS No.: _____

SDG No.: U04680

Matrix (soil/water): WATER

Lab Sample ID: 9A04G680-014

Level (low/med): LOW

Date Received: 04/08/00

Solids: 0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L

| CAS No. | Analyte | Concentration | C | Q | M |
|-----------|-----------|---------------|---|----|---------------|
| 7429-90-5 | Aluminum | | | | NR |
| 7440-36-0 | Antimony | | | | NR |
| 7440-38-2 | Arsenic | 8.8 | | JY | F CAE 6/13/00 |
| 7440-39-3 | Barium | | | | NR |
| 7440-41-7 | Beryllium | | | | NR |
| 7440-43-9 | Cadmium | | | | NR |
| 7440-70-2 | Calcium | | | | NR |
| 7440-47-3 | Chromium | | | | NR |
| 7440-48-4 | Cobalt | | | | NR |
| 7440-50-8 | Copper | | | | NR |
| 7439-89-6 | Iron | | | | NR |
| 7439-92-1 | Lead | | | | NR |
| 7439-95-4 | Magnesium | | | | NR |
| 7439-96-5 | Manganese | | | | NR |
| 7439-97-6 | Mercury | | | | NR |
| 7440-02-0 | Nickel | | | | NR |
| 7440-09-7 | Potassium | | | | NR |
| 7782-49-2 | Selenium | | | | NR |
| 7440-22-4 | Silver | | | | NR |
| 7440-23-5 | Sodium | | | | NR |
| 7440-28-0 | Thallium | | | | NR |
| 7440-62-2 | Vanadium | | | | NR |
| 7440-66-6 | Zinc | | | | NR |
| | Cyanide | | | | NR |

Color Before: COLORLESS Clarity Before: CLEAR Texture: _____

Color After: COLORLESS Clarity After: CLEAR Artifacts: _____

Comments:
SOLUBLE _____



TRILLIUM INC.
Consultants in Environmental Chemistry

356 FARRAGUT CROSSING DR.
KNOXVILLE, TN 37922

(423) 966-8880
FAX (423) 966-8885
cerikson@trilliuminc.com

June 13, 2000

Mr. Dan Garrigan
O&M, Inc.
450 Montbrook Lane
Knoxville, TN 37919

RE: Data Package Review - STL Lot No. 9A04G680 - Chloride in Surface Waters

Dear Dan:

I have completed my evaluation of the chloride analysis data prepared by the Severn Trent Laboratory (STL) in University Park, Illinois, for seven surface water samples from the L&RR site. The data were reported in a single data package under Lot No. 9A04G680, which was received by Trillium for review on April 28, 2000 and included the following samples:

| | | |
|-------|-------|-------|
| LCH-5 | LCH-4 | SW-16 |
| SW-10 | SW-8 | SW-5 |
| DUP | | |

Analyses were performed according to EPA Method 325.2 (EPA-600/4-79-020, March 1983). My evaluation was based on the specifications of the project-specific Quality Assurance Project Plan (QAPP, 9/96), the referenced method, and, to the extent applicable, the USEPA Region I *Laboratory Data Validation Functional Guidelines for Evaluating Inorganics Analyses* (6/88; Rev. 2/89).

Based on the evaluation, the following qualifiers were applied:

- Results for chloride in LCH-5 and DUP were qualified as estimated (J) due to poor reproducibility in the field duplicate analyses.

These qualifiers are reflected on the laboratory's Inorganic Data Reports included as Attachment A to this report. Region I qualifier definitions are also provided in this attachment.

At the discretion of the data user, the following points may warrant attention by field and laboratory personnel with regard to this data package and prior to future sampling events at this site:

HOME OFFICE:

28 GRACE'S DRIVE • COATESVILLE, PA19320 • (610) 383-7233 • FAX (610) 383-7907

OFFICES IN:

LOUISIANA • MARYLAND • NEW JERSEY • NORTH CAROLINA • PENNSYLVANIA • TENNESSEE • TEXAS

Mr. Dan Garrigan

Evaluation of STL Lot No. 9A04G680 (Chloride in SW)

June 13, 2000

Page 2

- Documentation of sample preservation performed in the field on all applicable chain of custody records.
- Documentation of cooler temperatures and sample pHs at the time of laboratory receipt on all applicable chain of custody records.
- Inclusion of raw data as needed to document calibrations and analysis of all samples.

Specific details regarding the review and evaluation of these data are discussed below:

Holding Times, Preservation, Sample Integrity: A copy of the field chain of custody (COC) record was included in the data package, documenting a sample collection date of 4/7/00 for the surface waters. Acceptable cooler temperatures (3.6°C and 2.6°C) on laboratory receipt were recorded on the COC. Acidification of the samples for chloride analysis, however, was not noted by the sampler on the COC and no verification of acidification was documented in the data package by the laboratory. At the request of the evaluator, the laboratory provided a revised COC for the surface water samples, with a checkmark in the "Preserved" space added.

The revised COC was inserted into the data package for G680 as page 199a.

Sample pHs were still not found on the additional documentation provided by the laboratory. For the purposes of this evaluation, it was assumed that samples for chloride analyses were properly preserved at the time of collection, and no further action was taken on this basis.

All sample analyses for chloride were performed within the 28-day holding time specified by the QAPP for this parameter.

Calibrations: Although a spectrophotometric method was referenced for chloride, no documentation of an initial calibration curve or calibration verification standards was provided in the data package. A statement in the narrative says, "The initial and continuing calibration verification standards and blanks were in control," but is not specific as to which method is referred to. At the request of the evaluator, copies of the applicable instrument print-outs and analyst log book pages as well as a summary data table for chloride were provided by the laboratory via FedEx on 6/9/00. Based on this documentation, an initial calibration (1.0-30 mg/L) was run for chloride on 4/12/00 (immediately prior to sample analyses) and an acceptable correlation coefficient (>0.995) was reported for that curve. Initial and continuing calibration verification standards were run at appropriate frequencies throughout the chloride analysis series, and all recoveries were acceptable (QC 85-115%).

The chloride raw data documentation was inserted into the data package for G680 as pages 12a through 12i by the evaluator.

Mr. Dan Garrigan

Evaluation of STL Lot No. 9A04G680 (Chloride in SW)

June 13, 2000

Page 3

Blanks: A method blank (MB) containing no chloride above the reporting limit was reported in the data package.

Calibration blanks were also run at the appropriate frequencies throughout the analysis series, based on the raw data documentation provided by the laboratory as discussed above. No chloride was detected above the reporting limit in any of the calibration blanks.

Laboratory Control Samples: A laboratory control sample (LCS) was analyzed in association with the sample analyses for chloride, and an acceptable recovery (103%) was reported.

Duplicate Analysis: No laboratory duplicate analysis was performed on any of the samples in this data package.

Matrix Spike Analysis: No matrix spike analysis was performed on any of the samples in this data package.

Sample Results: Sample results greater than or equal to the laboratory-specified reporting limits for chloride were reported on summary forms for all samples. Based on the raw data provided by the laboratory, some samples were appropriately re-analyzed at dilutions for some parameters

Field Duplicates: Sample DUP was identified as a field duplicate of LCH-4. Paired results for chloride showed poor reproducibility, with a relative percent difference of 47.7%. Results for chloride in LCH-4 and DUP were qualified as estimated (J) on this basis.

Documentation: Data package documentation was consistent with the specifications of the QAPP; specifically, a Level III package containing summary forms only was provided. However, no documentation of initial or continuing calibration standards for chloride was provided, although this information has been provided in previous data packages generated in support of this project. For wet chemistry parameters in general, the raw data pages serve as run logs and are helpful for verifying that all samples were, in fact, analyzed. For those parameters run by spectrophotometric methods, the raw data pages are necessary to evaluate calibrations and calibration blanks, which are not routinely summarized elsewhere. Additional documentation for chloride was therefore requested from and provided by the laboratory as discussed previously in this report.

Documentation of sample pHs and cooler temperatures on laboratory receipt was also incomplete in the data packages as received for review. Clarifications and corrections, as appropriate, were verbally requested by the validator on 6/6/00; responses were received from the laboratory via FedEx on 6/9/00. Corrected documents were inserted into the applicable data packages as discussed previously in this report.

Mr. Dan Garrigan
Evaluation of STL Lot No. 9A04G680 (Chloride in SW)
June 13, 2000
Page 4

A revised COC for the surface water samples, with a checkmark in the "Preserved" space added, was provided by the laboratory, but this new entry was not initialed or dated. While this has no direct effect on the technical validity of the reported sample data, it could be problematic if the data are used for litigation.

Please let me know if you have any questions regarding this data package review report.

Sincerely,



Carol A. Erikson
Quality Assessment Manager

CAE/hrs
95406\Apr00\swchloride

ATTACHMENT A

**Chloride in Surface Waters
STL Lot No. 9A04G680**

EPA Region I Qualifier Definitions:

- U - The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit.**
- J - The associated value is an estimated quantity.**
- R - The data are unusable. (Note: Analyte may or may not be present).**
- UJ - The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.**

To: O&M-L&RR Landfill
301 Gallaher View Road
Suite 230
Knoxville, TN 37919

Attn: Mr. Daniel Garrigan

Date: Thursday April 20th, 2000

RE: LCH-5
Project # 00000-000-000-0000
Lab ID: 9A04G680-001
Sample Date: 04/07/00
Date Received: 04/08/00

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|-------|-----------------|
| Chloride | 16.3 | mg/L | 1.0 |

To: O&M-L&RR Landfill
301 Gallaher View Road
Suite 230
Knoxville, TN 37919

Attn: Mr. Daniel Garrigan

Date: Thursday April 20th, 2000

RE: LCH-4
Project # 00000-000-000-0000
Lab ID: 9A04G680-003
Sample Date: 04/07/00
Date Received: 04/08/00

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|------------------------|-------|-----------------|
| Chloride | 80.7 J | mg/L | 5.0 |
| | <i>caution 6/13/00</i> | | |

To: O&M-L&RR Landfill
301 Gallaher View Road
Suite 230
Knoxville, TN 37919

Attn: Mr. Daniel Garrigan

Date: Thursday April 20th, 2000

RE: SW-16
Project # 00000-000-000-0000
Lab ID: 9A04G680-005
Sample Date: 04/07/00
Date Received: 04/08/00

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|-------|--------------------|
| Chloride | 29.4 | mg/L | 1.0 |

To: O&M-L&RR Landfill
301 Gallaher View Road
Suite 230
Knoxville, TN 37919

Attn: Mr. Daniel Garrigan

Date: Thursday April 20th, 2000

RE: SW-10
Project # 00000-000-000-0000
Lab ID: 9A04G680-007
Sample Date: 04/07/00
Date Received: 04/08/00

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|-------|--------------------|
| Chloride | 30.2 | mg/L | 2.0 |

To: O&M-L&RR Landfill
301 Gallaher View Road
Suite 230
Knoxville, TN 37919

Attn: Mr. Daniel Garrigan

Date: Thursday April 20th, 2000

RE: SW-8
Project # 00000-000-000-0000
Lab ID: 9A04G680-009
Sample Date: 04/07/00
Date Received: 04/08/00

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|-------|-----------------|
| Chloride | 58.3 | mg/L | 5.0 |

To: O&M-L&RR Landfill
301 Gallaher View Road
Suite 230
Knoxville, TN 37919

Attn: Mr. Daniel Garrigan

Date: Thursday April 20th, 2000

RE: SW-5
Project # 00000-000-000-0000
Lab ID: 9A04G680-011
Sample Date: 04/07/00
Date Received: 04/08/00

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|--------|-------|--------------------|
| Chloride | 10.5 | mg/L | 1.0 |

To: O&M-L&RR Landfill
301 Gallaher View Road
Suite 230
Knoxville, TN 37919

Attn: Mr. Daniel Garrigan

Date: Thursday April 20th, 2000

RE: DUP
Project # 00000-000-000-0000
Lab ID: 9A04G680-013
Sample Date: 04/07/00
Date Received: 04/08/00

Inorganic Data Report

| Parameters | Result | Units | Reporting Limit |
|------------|-------------|-------|-----------------|
| Chloride | 49.6 J | mg/L | 2.0 |
| | CAE 4/13/00 | | |

Appendix F

**Ambient Air Sampling Report
AirRecon, Inc.**

July 11, 2000

Mr. Dan Garrigan
Operations and Maintenance, Inc.
301 Gallaher View Rd.
Suite 230
Knoxville, TN 37919

**RE: L&RR Landfill
North Smithfield, RI
AirRECON Project 311-00105-00-000**

Dear Mr. Garrigan:

Enclosed, please find copies of the laboratory results from the test program performed at the L&RR landfill located in North Smithfield RI, and associated data from the test program.

Five TO-14 samples were collected around the perimeter of the landfill. The sampling duration was twenty-two (22) hours. The shortened sampling duration was used to insure samples did not go "ambient" and that enough pressure was left in the samples during shipment back to the laboratory.

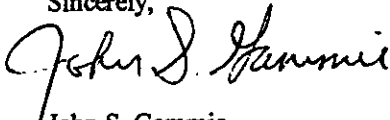
In addition, one (1) USEPA Method 18 grab sample was also collected at the header to the flare. This sample was collected over a four-hour period.

All sampling was performed without incident, and by USEPA methodologies.

If you have any questions or require additional information, please do not hesitate to contact me at (860) 721-7728.

It was a pleasure being of service to you. Should you require additional work please feel free to contact me.

Sincerely,



John S. Gammie
Sr. Environmental Specialist
AirRECON a division of LFR

Cc. Andrew McNeel/ AirRECON/ NJ
Thor Helgasin/ DeMaximus, Inc.

Sampling Data on June 8-9 2000 at the Smithfield, RI Landfill TO 14 Samples

Weather Conditions June 8, 2000 Barometric Pressure 30.13, dewpoint 11, west wind (0-4)mph, temp 75-80F

| Location | start time | date | finish time | date | tank # | orifice# |
|-------------------------|------------|--------|-------------|--------|--------|---------------|
| AS-1 near gate entrance | 1120am | 6/8/00 | 0920am | 6/9/00 | 332 | 05967 (1) |
| AS-1 dup | 1120am | 6/8/00 | 0920am | 6/9/00 | 271 | 05033 (95-03) |
| AS-2 | 1124am | 6/8/00 | 0924am | 6/9/00 | 2701 | 05483 (98-17) |
| AS-3 | 1129am | 6/8/00 | 0929am | 6/9/00 | 2507 | 05995 (97-03) |
| AS-4 closest to flare | 1133am | 6/8/00 | 0933am | 6/9/00 | 913 | 05493 (96-06) |

Weather Conditions 6-9-2000 Barometric Pressure 29.97, dew point 17, west wind (0-9) mph temp 75-80F

Method 18 Sampling data

| Start time | date | finish time | date |
|------------|--------|-------------|--------|
| 0953am | 6/8/00 | 1353pm | 6/8/00 |

WORK ORDER #: 0006175

Work Order Summary

CLIENT: Tom Brown
LFR Levine-Fricke
5 Johnson Drive
P.O. Box 130
Raritan, NJ 08869

BILL TO: Tom Brown
LFR Levine-Fricke
5 Johnson Drive
P.O. Box 130
Raritan, NJ 08869

PHONE: 908-526-1000

P.O. #

FAX: 908-526-7886

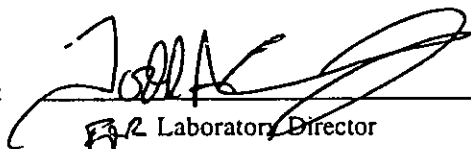
PROJECT # 311-00105 OIM

DATE RECEIVED: 6/12/00

DATE COMPLETED: 6/22/00

| <u>FRACTION #</u> | <u>NAME</u> | <u>TEST</u> | <u>RECEIPT YAC./PRES.</u> |
|-------------------|-------------|-------------|-------------------------------|
| 01A | AS-1 | TO-14 | 6.5 "Hg |
| 02A | AS-1 Dup | TO-14 | 6.0 "Hg |
| 03A | AS-2 | TO-14 | 5.0 "Hg |
| 04A | AS-3 | TO-14 | 6.0 "Hg |
| 05A | AS-4 | TO-14 | 6.0 "Hg |
| 06A | BLANK | TO-14 | 28.5 "Hg |
| 07A | Lab Blank | TO-14 | NA |
| 07B | Lab Blank | TO-14 | NA |

CERTIFIED BY:


LFR Laboratory Director

DATE:

6/29/00

Certification numbers: CA ELAP - 1149, NY ELAP - 11291, UT ELAP - E-217, AZ ELAP - AZ0567

180 BLUE RAVINE ROAD, SUITE B FOLSOM, CA - 95630
(916) 985-1000 . (800) 985-5955 . FAX (916) 985-1020

LABORATORY NARRATIVE

TO-14

LFR Levine-Fricke
Workorder# 0006175

Six 6 Liter Summa Canister samples were received on June 12, 2000. The laboratory performed analysis via EPA Method TO-14 using GC/MS in the full scan mode. The method involves concentrating up to 0.5 liters of air. The concentrated aliquot is then flash vaporized and swept through a water management system to remove water vapor. Following dehumidification, the sample passes directly into the GC/MS for analysis. See the data sheets for the reporting limits for each compound.

During the five point calibration, two low-level standards are used. The low-level standard for TO-14 compounds is spiked at 0.5 ppbv and represents the reporting limit for these compounds. The low-level standard for the non-TO-14 compounds is spiked at 2.0 ppbv and represents the reporting limit for these compounds. The TO-14 compounds are present in both standards but are excluded from reporting in the 2.0 ppbv standard since a lower level is already included in the curve.

Method modifications taken to run these samples include:

| <i>Requirement</i> | <i>TO-14</i> | <i>ATL Modifications</i> |
|--|---------------------------------|--|
| Internal standard retention times. | Not specified. | Within 0.50 minutes of most recent daily CCV internal standards |
| Internal standard recoveries. | Not specified. | Within 40% of the daily CCV internal standard area for blanks and samples. |
| Internal standard retention times. | Not specified. | Within 0.50 minutes of most recent daily CCV internal standards |
| Internal calibration criteria. | Not specified. | RSD of 30% or less for standard compounds, 40% or less for non-standard and polar compounds |
| Continuing calibration verification criteria | Not specified. | 70 - 130% for at least 90% of standard compounds, 60 - 140% for at least 80% of non-standard and polar compounds |
| Response factor for quantitation. | Average response factor (ICAL). | Average response factor (ICAL). |

Receiving Notes

The chain of custody information for sample BLANK did not match the entry on the sample tag. The client was notified and the information on the chain of custody was used to process and report the sample.

Analytical Notes

Methylene Chloride was detected in the laboratory blank analyzed on 06/20/00 at less than 5X the reporting limit. Associated sample results are reported as qualified.

Definition of Data Qualifying Flags

Seven qualifiers may have been used on the data analysis sheets and indicates as follows:

B - Compound present in laboratory blank greater than reporting limit(background subtraction not performed).

J - Estimated value.

E - Exceeds instrument calibration range.

S - Saturated peak.

Q - Exceeds quality control limits.

U - Compound analyzed for but not detected above the reporting limit.

N - The identification is based on presumptive evidence.

AIR TOXICS LTD.

SAMPLE NAME : AS-1

ID#: 0006175-01A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | c061919 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.71 | Date of Analysis: | 6/19/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|---------------------------|----------------------|-----------------------|------------------|-------------------|
| Freon 12 | 0.86 | 4.3 | 1.0 | 5.3 |
| Freon 114 | 0.86 | 6.1 | Not Detected | Not Detected |
| Chloromethane | 0.86 | 1.8 | 1.5 | 3.1 |
| Vinyl Chloride | 0.86 | 2.2 | Not Detected | Not Detected |
| Bromomethane | 0.86 | 3.4 | Not Detected | Not Detected |
| Chloroethane | 0.86 | 2.3 | Not Detected | Not Detected |
| Freon 11 | 0.86 | 4.9 | Not Detected | Not Detected |
| 1,1-Dichloroethene | 0.86 | 3.4 | Not Detected | Not Detected |
| Freon 113 | 0.86 | 6.7 | Not Detected | Not Detected |
| Methylene Chloride | 0.86 | 3.0 | 1.4 | 5.0 |
| 1,1-Dichloroethane | 0.86 | 3.5 | Not Detected | Not Detected |
| cis-1,2-Dichloroethene | 0.86 | 3.4 | Not Detected | Not Detected |
| Chloroform | 0.86 | 4.2 | Not Detected | Not Detected |
| 1,1,1-Trichloroethane | 0.86 | 4.7 | Not Detected | Not Detected |
| Carbon Tetrachloride | 0.86 | 5.5 | Not Detected | Not Detected |
| Benzene | 0.86 | 2.8 | Not Detected | Not Detected |
| 1,2-Dichloroethane | 0.86 | 3.5 | Not Detected | Not Detected |
| Trichloroethene | 0.86 | 4.7 | Not Detected | Not Detected |
| 1,2-Dichloropropane | 0.86 | 4.0 | Not Detected | Not Detected |
| cis-1,3-Dichloropropene | 0.86 | 3.9 | Not Detected | Not Detected |
| Toluene | 0.86 | 3.3 | Not Detected | Not Detected |
| trans-1,3-Dichloropropene | 0.86 | 3.9 | Not Detected | Not Detected |
| 1,1,2-Trichloroethane | 0.86 | 4.7 | Not Detected | Not Detected |
| Tetrachloroethene | 0.86 | 5.9 | Not Detected | Not Detected |
| Ethylene Dibromide | 0.86 | 6.7 | Not Detected | Not Detected |
| Chlorobenzene | 0.86 | 4.0 | Not Detected | Not Detected |
| Ethyl Benzene | 0.86 | 3.8 | Not Detected | Not Detected |
| m,p-Xylene | 0.86 | 3.8 | Not Detected | Not Detected |
| o-Xylene | 0.86 | 3.8 | Not Detected | Not Detected |
| Styrene | 0.86 | 3.7 | Not Detected | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.86 | 6.0 | Not Detected | Not Detected |
| 1,3,5-Trimethylbenzene | 0.86 | 4.3 | Not Detected | Not Detected |
| 1,2,4-Trimethylbenzene | 0.86 | 4.3 | Not Detected | Not Detected |
| 1,3-Dichlorobenzene | 0.86 | 5.2 | Not Detected | Not Detected |
| 1,4-Dichlorobenzene | 0.86 | 5.2 | Not Detected | Not Detected |
| Chlorotoluene | 0.86 | 4.5 | Not Detected | Not Detected |
| 1,2-Dichlorobenzene | 0.86 | 5.2 | Not Detected | Not Detected |
| 1,2,4-Trichlorobenzene | 0.86 | 6.4 | Not Detected | Not Detected |
| Hexachlorobutadiene | 0.86 | 9.3 | Not Detected | Not Detected |
| Propylene | 3.4 | 6.0 | Not Detected | Not Detected |
| 1,3-Butadiene | 3.4 | 7.7 | Not Detected | Not Detected |
| Acetone | 3.4 | 8.2 | 16 | 39 |

AIR TOXICS LTD.

SAMPLE NAME : AS-1

ID#: 0006175-01A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | 0061919 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1/1 | Date of Analysis: | 6/19/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|----------------------------------|-------------------|--------------------|---------------|----------------|
| Carbon Disulfide | 3.4 | 11 | Not Detected | Not Detected |
| 2-Propanol | 3.4 | 8.5 | Not Detected | Not Detected |
| trans-1,2-Dichloroethene | 3.4 | 14 | Not Detected | Not Detected |
| Vinyl Acetate | 3.4 | 12 | Not Detected | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 3.4 | 10 | Not Detected | Not Detected |
| Hexane | 3.4 | 12 | Not Detected | Not Detected |
| Tetrahydrofuran | 3.4 | 10 | Not Detected | Not Detected |
| Cyclohexane | 3.4 | 12 | Not Detected | Not Detected |
| 1,4-Dioxane | 3.4 | 12 | Not Detected | Not Detected |
| Bromodichloromethane | 3.4 | 23 | Not Detected | Not Detected |
| 4-Methyl-2-pentanone | 3.4 | 14 | Not Detected | Not Detected |
| 2-Hexanone | 3.4 | 14 | Not Detected | Not Detected |
| Dibromochloromethane | 3.4 | 30 | Not Detected | Not Detected |
| Bromoform | 3.4 | 36 | Not Detected | Not Detected |
| 4-Ethyltoluene | 3.4 | 17 | Not Detected | Not Detected |
| Ethanol | 3.4 | 6.5 | Not Detected | Not Detected |
| Methyl tert-Butyl Ether | 3.4 | 12 | Not Detected | Not Detected |
| Heptane | 3.4 | 14 | Not Detected | Not Detected |

Container Type: 6 Liter Summa Canister

| Surrogates | % Recovery | Method Limits |
|-----------------------|------------|---------------|
| 1,2-Dichloroethane-d4 | 104 | 70-130 |
| Toluene-d8 | 105 | 70-130 |
| 4-Bromofluorobenzene | 94 | 70-130 |

AIR TOXICS LTD.

SAMPLE NAME : AS-1 Dup

ID#: 0006175-02A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | 6061920 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.68 | Date of Analysis: | 6/19/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|---------------------------|----------------------|-----------------------|------------------|-------------------|
| Freon 12 | 0.84 | 4.2 | Not Detected | Not Detected |
| Freon 114 | 0.84 | 6.0 | Not Detected | Not Detected |
| Chloromethane | 0.84 | 1.8 | Not Detected | Not Detected |
| Vinyl Chloride | 0.84 | 2.2 | Not Detected | Not Detected |
| Bromomethane | 0.84 | 3.3 | Not Detected | Not Detected |
| Chloroethane | 0.84 | 2.2 | Not Detected | Not Detected |
| Freon 11 | 0.84 | 4.8 | Not Detected | Not Detected |
| 1,1-Dichloroethene | 0.84 | 3.4 | Not Detected | Not Detected |
| Freon 113 | 0.84 | 6.5 | Not Detected | Not Detected |
| Methylene Chloride | 0.84 | 3.0 | 1.4 | 5.1 |
| 1,1-Dichloroethane | 0.84 | 3.4 | Not Detected | Not Detected |
| cis-1,2-Dichloroethene | 0.84 | 3.4 | Not Detected | Not Detected |
| Chloroform | 0.84 | 4.2 | Not Detected | Not Detected |
| 1,1,1-Trichloroethane | 0.84 | 4.6 | Not Detected | Not Detected |
| Carbon Tetrachloride | 0.84 | 5.4 | Not Detected | Not Detected |
| Benzene | 0.84 | 2.7 | Not Detected | Not Detected |
| 1,2-Dichloroethane | 0.84 | 3.4 | Not Detected | Not Detected |
| Trichloroethene | 0.84 | 4.6 | Not Detected | Not Detected |
| 1,2-Dichloropropane | 0.84 | 3.9 | Not Detected | Not Detected |
| cis-1,3-Dichloropropene | 0.84 | 3.9 | Not Detected | Not Detected |
| Toluene | 0.84 | 3.2 | 4.9 | 19 |
| trans-1,3-Dichloropropene | 0.84 | 3.9 | Not Detected | Not Detected |
| 1,1,2-Trichloroethane | 0.84 | 4.6 | Not Detected | Not Detected |
| Tetrachloroethene | 0.84 | 5.8 | Not Detected | Not Detected |
| Ethylene Dibromide | 0.84 | 6.6 | Not Detected | Not Detected |
| Chlorobenzene | 0.84 | 3.9 | Not Detected | Not Detected |
| Ethyl Benzene | 0.84 | 3.7 | 3.3 | 15 |
| m,p-Xylene | 0.84 | 3.7 | 23 | 100 |
| o-Xylene | 0.84 | 3.7 | 3.8 | 16 |
| Styrene | 0.84 | 3.6 | Not Detected | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.84 | 5.9 | Not Detected | Not Detected |
| 1,3,5-Trimethylbenzene | 0.84 | 4.2 | 2.1 | 10 |
| 1,2,4-Trimethylbenzene | 0.84 | 4.2 | 1.0 | 5.2 |
| 1,3-Dichlorobenzene | 0.84 | 5.1 | Not Detected | Not Detected |
| 1,4-Dichlorobenzene | 0.84 | 5.1 | 0.99 | 6.0 |
| Chlorotoluene | 0.84 | 4.4 | Not Detected | Not Detected |
| 1,2-Dichlorobenzene | 0.84 | 5.1 | 0.94 | 5.8 |
| 1,2,4-Trichlorobenzene | 0.84 | 6.3 | Not Detected | Not Detected |
| Hexachlorobutadiene | 0.84 | 9.1 | Not Detected | Not Detected |
| Propylene | 3.4 | 5.9 | Not Detected | Not Detected |
| 1,3-Butadiene | 3.4 | 7.6 | Not Detected | Not Detected |
| Acetone | 3.4 | 8.1 | 36 | 87 |

AIR TOXICS LTD.

SAMPLE NAME : AS-1 Dup

ID#: 0006175-02A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | C061920 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.68 | Date of Analysis: | 6/19/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|----------------------------------|-------------------|--------------------|---------------|----------------|
| Carbon Disulfide | 3.4 | 11 | 4.1 | 13 |
| 2-Propanol | 3.4 | 8.4 | Not Detected | Not Detected |
| trans-1,2-Dichloroethene | 3.4 | 14 | Not Detected | Not Detected |
| Vinyl Acetate | 3.4 | 12 | Not Detected | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 3.4 | 10 | 5.9 | 18 |
| Hexane | 3.4 | 12 | Not Detected | Not Detected |
| Tetrahydrofuran | 3.4 | 10 | Not Detected | Not Detected |
| Cyclohexane | 3.4 | 12 | Not Detected | Not Detected |
| 1,4-Dioxane | 3.4 | 12 | Not Detected | Not Detected |
| Bromodichloromethane | 3.4 | 23 | Not Detected | Not Detected |
| 4-Methyl-2-pentanone | 3.4 | 14 | Not Detected | Not Detected |
| 2-Hexanone | 3.4 | 14 | Not Detected | Not Detected |
| Dibromochloromethane | 3.4 | 29 | Not Detected | Not Detected |
| Bromoform | 3.4 | 35 | Not Detected | Not Detected |
| 4-Ethyltoluene | 3.4 | 17 | Not Detected | Not Detected |
| Ethanol | 3.4 | 6.4 | Not Detected | Not Detected |
| Methyl tert-Butyl Ether | 3.4 | 12 | Not Detected | Not Detected |
| Heptane | 3.4 | 14 | Not Detected | Not Detected |

Container Type: 6 Liter Summa Canister

| Surrogates | % Recovery | Method Limits |
|-----------------------|------------|---------------|
| 1,2-Dichloroethane-d4 | 106 | 70-130 |
| Toluene-d8 | 102 | 70-130 |
| 4-Bromofluorobenzene | 94 | 70-130 |

AIR TOXICS LTD.

SAMPLE NAME : AS-2

ID#: 0006175-03A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|-------------|---------|---------------------|---------|
| File Name: | c061921 | Date of Collection: | 6/8/00 |
| Dil Factor: | 1.61 | Date of Analysis: | 6/19/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|---------------------------|-------------------|--------------------|---------------|----------------|
| Freon 12 | 0.80 | 4.0 | 0.82 | 4.1 |
| Freon 114 | 0.80 | 5.7 | Not Detected | Not Detected |
| Chloromethane | 0.80 | 1.7 | Not Detected | Not Detected |
| Vinyl Chloride | 0.80 | 2.1 | Not Detected | Not Detected |
| Bromomethane | 0.80 | 3.2 | Not Detected | Not Detected |
| Chloroethane | 0.80 | 2.2 | Not Detected | Not Detected |
| Freon 11 | 0.80 | 4.6 | Not Detected | Not Detected |
| 1,1-Dichloroethene | 0.80 | 3.2 | Not Detected | Not Detected |
| Freon 113 | 0.80 | 6.3 | Not Detected | Not Detected |
| Methylene Chloride | 0.80 | 2.8 | 0.89 | 3.1 |
| 1,1-Dichloroethane | 0.80 | 3.3 | Not Detected | Not Detected |
| cis-1,2-Dichloroethene | 0.80 | 3.2 | Not Detected | Not Detected |
| Chloroform | 0.80 | 4.0 | Not Detected | Not Detected |
| 1,1,1-Trichloroethane | 0.80 | 4.5 | Not Detected | Not Detected |
| Carbon Tetrachloride | 0.80 | 5.1 | Not Detected | Not Detected |
| Benzene | 0.80 | 2.6 | Not Detected | Not Detected |
| 1,2-Dichloroethane | 0.80 | 3.3 | Not Detected | Not Detected |
| Trichloroethene | 0.80 | 4.4 | Not Detected | Not Detected |
| 1,2-Dichloropropane | 0.80 | 3.8 | Not Detected | Not Detected |
| cis-1,3-Dichloropropene | 0.80 | 3.7 | Not Detected | Not Detected |
| Toluene | 0.80 | 3.1 | Not Detected | Not Detected |
| trans-1,3-Dichloropropene | 0.80 | 3.7 | Not Detected | Not Detected |
| 1,1,2-Trichloroethane | 0.80 | 4.5 | Not Detected | Not Detected |
| Tetrachloroethene | 0.80 | 5.6 | Not Detected | Not Detected |
| Ethylene Dibromide | 0.80 | 6.3 | Not Detected | Not Detected |
| Chlorobenzene | 0.80 | 3.8 | Not Detected | Not Detected |
| Ethyl Benzene | 0.80 | 3.6 | Not Detected | Not Detected |
| m,p-Xylene | 0.80 | 3.6 | Not Detected | Not Detected |
| o-Xylene | 0.80 | 3.6 | Not Detected | Not Detected |
| Styrene | 0.80 | 3.5 | Not Detected | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.80 | 5.6 | Not Detected | Not Detected |
| 1,3,5-Trimethylbenzene | 0.80 | 4.0 | Not Detected | Not Detected |
| 1,2,4-Trimethylbenzene | 0.80 | 4.0 | Not Detected | Not Detected |
| 1,3-Dichlorobenzene | 0.80 | 4.9 | Not Detected | Not Detected |
| 1,4-Dichlorobenzene | 0.80 | 4.9 | Not Detected | Not Detected |
| Chlorotoluene | 0.80 | 4.2 | Not Detected | Not Detected |
| 1,2-Dichlorobenzene | 0.80 | 4.9 | Not Detected | Not Detected |
| 1,2,4-Trichlorobenzene | 0.80 | 6.1 | Not Detected | Not Detected |
| Hexachlorobutadiene | 0.80 | 8.7 | Not Detected | Not Detected |
| Propylene | 3.2 | 5.6 | Not Detected | Not Detected |
| 1,3-Butadiene | 3.2 | 7.2 | Not Detected | Not Detected |
| Acetone | 3.2 | 7.8 | 23 | 55 |

AIR TOXICS LTD.

SAMPLE NAME : AS-2

ID#: 0006175-03A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | C061921 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.61 | Date of Analysis: | 6/19/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|----------------------------------|----------------------|-----------------------|------------------|-------------------|
| Carbon Disulfide | 3.2 | 10 | Not Detected | Not Detected |
| 2-Propanol | 3.2 | 8.0 | Not Detected | Not Detected |
| trans-1,2-Dichloroethene | 3.2 | 13 | Not Detected | Not Detected |
| Vinyl Acetate | 3.2 | 12 | Not Detected | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 3.2 | 9.6 | 3.3 | 9.9 |
| Hexane | 3.2 | 12 | Not Detected | Not Detected |
| Tetrahydrofuran | 3.2 | 9.6 | Not Detected | Not Detected |
| Cyclohexane | 3.2 | 11 | Not Detected | Not Detected |
| 1,4-Dioxane | 3.2 | 12 | Not Detected | Not Detected |
| Bromodichloromethane | 3.2 | 22 | Not Detected | Not Detected |
| 4-Methyl-2-pentanone | 3.2 | 13 | Not Detected | Not Detected |
| 2-Hexanone | 3.2 | 13 | Not Detected | Not Detected |
| Dibromochloromethane | 3.2 | 28 | Not Detected | Not Detected |
| Bromoform | 3.2 | 34 | Not Detected | Not Detected |
| 4-Ethyltoluene | 3.2 | 16 | Not Detected | Not Detected |
| Ethanol | 3.2 | 6.2 | Not Detected | Not Detected |
| Methyl tert-Butyl Ether | 3.2 | 12 | Not Detected | Not Detected |
| Heptane | 3.2 | 13 | Not Detected | Not Detected |

Container Type: 6 Liter Summa Canister

| Surrogates | % Recovery | Method Limits |
|-----------------------|------------|---------------|
| 1,2-Dichloroethane-d4 | 108 | 70-130 |
| Toluene-d8 | 100 | 70-130 |
| 4-Bromofluorobenzene | 87 | 70-130 |

AIR TOXICS LTD.

SAMPLE NAME : AS-3

ID#: 0006175-04A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | g062019 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.55 | Date of Analysis: | 6/20/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|---------------------------|-------------------|--------------------|---------------|----------------|
| Freon 12 | 0.78 | 3.9 | Not Detected | Not Detected |
| Freon 114 | 0.78 | 5.5 | Not Detected | Not Detected |
| Chloromethane | 0.78 | 1.6 | Not Detected | Not Detected |
| Vinyl Chloride | 0.78 | 2.0 | Not Detected | Not Detected |
| Bromomethane | 0.78 | 3.0 | Not Detected | Not Detected |
| Chloroethane | 0.78 | 2.1 | Not Detected | Not Detected |
| Freon 11 | 0.78 | 4.4 | Not Detected | Not Detected |
| 1,1-Dichloroethene | 0.78 | 3.1 | Not Detected | Not Detected |
| Freon 113 | 0.78 | 6.0 | Not Detected | Not Detected |
| Methylene Chloride | 0.78 | 2.7 | 1.4 B | 4.9 B |
| 1,1-Dichloroethane | 0.78 | 3.2 | Not Detected | Not Detected |
| cis-1,2-Dichloroethene | 0.78 | 3.1 | Not Detected | Not Detected |
| Chloroform | 0.78 | 3.8 | Not Detected | Not Detected |
| 1,1,1-Trichloroethane | 0.78 | 4.3 | Not Detected | Not Detected |
| Carbon Tetrachloride | 0.78 | 5.0 | Not Detected | Not Detected |
| Benzene | 0.78 | 2.5 | Not Detected | Not Detected |
| 1,2-Dichloroethane | 0.78 | 3.2 | Not Detected | Not Detected |
| Trichloroethene | 0.78 | 4.2 | Not Detected | Not Detected |
| 1,2-Dichloropropane | 0.78 | 3.6 | Not Detected | Not Detected |
| cis-1,3-Dichloropropene | 0.78 | 3.6 | Not Detected | Not Detected |
| Toluene | 0.78 | 3.0 | Not Detected | Not Detected |
| trans-1,3-Dichloropropene | 0.78 | 3.6 | Not Detected | Not Detected |
| 1,1,2-Trichloroethane | 0.78 | 4.3 | Not Detected | Not Detected |
| Tetrachloroethene | 0.78 | 5.3 | Not Detected | Not Detected |
| Ethylene Dibromide | 0.78 | 6.0 | Not Detected | Not Detected |
| Chlorobenzene | 0.78 | 3.6 | Not Detected | Not Detected |
| Ethyl Benzene | 0.78 | 3.4 | Not Detected | Not Detected |
| m,p-Xylene | 0.78 | 3.4 | Not Detected | Not Detected |
| o-Xylene | 0.78 | 3.4 | Not Detected | Not Detected |
| Styrene | 0.78 | 3.4 | Not Detected | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.78 | 5.4 | Not Detected | Not Detected |
| 1,3,5-Trimethylbenzene | 0.78 | 3.9 | Not Detected | Not Detected |
| 1,2,4-Trimethylbenzene | 0.78 | 3.9 | Not Detected | Not Detected |
| 1,3-Dichlorobenzene | 0.78 | 4.7 | Not Detected | Not Detected |
| 1,4-Dichlorobenzene | 0.78 | 4.7 | Not Detected | Not Detected |
| Chlorotoluene | 0.78 | 4.1 | Not Detected | Not Detected |
| 1,2-Dichlorobenzene | 0.78 | 4.7 | Not Detected | Not Detected |
| 1,2,4-Trichlorobenzene | 0.78 | 5.8 | Not Detected | Not Detected |
| Hexachlorobutadiene | 0.78 | 8.4 | Not Detected | Not Detected |
| Propylene | 3.1 | 5.4 | Not Detected | Not Detected |
| 1,3-Butadiene | 3.1 | 7.0 | Not Detected | Not Detected |
| Acetone | 3.1 | 7.5 | 6.8 | 16 |

AIR TOXICS LTD.

SAMPLE NAME : AS-3

ID#: 0006175-04A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | g062019 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.55 | Date of Analysis: | 6/20/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|----------------------------------|----------------------|-----------------------|------------------|-------------------|
| Carbon Disulfide | 3.1 | 9.8 | Not Detected | Not Detected |
| 2-Propanol | 3.1 | 7.7 | Not Detected | Not Detected |
| trans-1,2-Dichloroethene | 3.1 | 12 | Not Detected | Not Detected |
| Vinyl Acetate | 3.1 | 11 | Not Detected | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 3.1 | 9.3 | Not Detected | Not Detected |
| Hexane | 3.1 | 11 | Not Detected | Not Detected |
| Tetrahydrofuran | 3.1 | 9.3 | Not Detected | Not Detected |
| Cyclohexane | 3.1 | 11 | Not Detected | Not Detected |
| 1,4-Dioxane | 3.1 | 11 | Not Detected | Not Detected |
| Bromodichloromethane | 3.1 | 21 | Not Detected | Not Detected |
| 4-Methyl-2-pentanone | 3.1 | 13 | Not Detected | Not Detected |
| 2-Hexanone | 3.1 | 13 | Not Detected | Not Detected |
| Dibromochloromethane | 3.1 | 27 | Not Detected | Not Detected |
| Bromoform | 3.1 | 32 | Not Detected | Not Detected |
| 4-Ethyltoluene | 3.1 | 15 | Not Detected | Not Detected |
| Ethanol | 3.1 | 5.9 | Not Detected | Not Detected |
| Methyl tert-Butyl Ether | 3.1 | 11 | Not Detected | Not Detected |
| Heptane | 3.1 | 13 | Not Detected | Not Detected |

B = Compound present in laboratory blank, background subtraction not performed.

Container Type: 6 Liter Summa Canister

| Surrogates | % Recovery | Method Limits |
|-----------------------|------------|------------------|
| 1,2-Dichloroethane-d4 | 104 | 70-130 |
| Toluene-d8 | 101 | 70-130 |
| 4-Bromofluorobenzene | 93 | 70-130 |

AIR TOXICS LTD.

SAMPLE NAME : AS-4

ID#: 0006175-05A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | c061923 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.68 | Date of Analysis: | 6/20/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|---------------------------|----------------------|-----------------------|------------------|-------------------|
| Freon 12 | 0.84 | 4.2 | 0.91 | 4.6 |
| Freon 114 | 0.84 | 6.0 | Not Detected | Not Detected |
| Chloromethane | 0.84 | 1.8 | Not Detected | Not Detected |
| Vinyl Chloride | 0.84 | 2.2 | Not Detected | Not Detected |
| Bromomethane | 0.84 | 3.3 | Not Detected | Not Detected |
| Chloroethane | 0.84 | 2.2 | Not Detected | Not Detected |
| Freon 11 | 0.84 | 4.8 | Not Detected | Not Detected |
| 1,1-Dichloroethene | 0.84 | 3.4 | Not Detected | Not Detected |
| Freon 113 | 0.84 | 6.5 | Not Detected | Not Detected |
| Methylene Chloride | 0.84 | 3.0 | 0.90 | 3.2 |
| 1,1-Dichloroethane | 0.84 | 3.4 | Not Detected | Not Detected |
| cis-1,2-Dichloroethene | 0.84 | 3.4 | Not Detected | Not Detected |
| Chloroform | 0.84 | 4.2 | Not Detected | Not Detected |
| 1,1,1-Trichloroethane | 0.84 | 4.6 | Not Detected | Not Detected |
| Carbon Tetrachloride | 0.84 | 5.4 | Not Detected | Not Detected |
| Benzene | 0.84 | 2.7 | Not Detected | Not Detected |
| 1,2-Dichloroethane | 0.84 | 3.4 | Not Detected | Not Detected |
| Trichloroethene | 0.84 | 4.6 | Not Detected | Not Detected |
| 1,2-Dichloropropane | 0.84 | 3.9 | Not Detected | Not Detected |
| cis-1,3-Dichloropropene | 0.84 | 3.9 | Not Detected | Not Detected |
| Toluene | 0.84 | 3.2 | Not Detected | Not Detected |
| trans-1,3-Dichloropropene | 0.84 | 3.9 | Not Detected | Not Detected |
| 1,1,2-Trichloroethane | 0.84 | 4.6 | Not Detected | Not Detected |
| Tetrachloroethene | 0.84 | 5.8 | Not Detected | Not Detected |
| Ethylene Dibromide | 0.84 | 6.6 | Not Detected | Not Detected |
| Chlorobenzene | 0.84 | 3.9 | Not Detected | Not Detected |
| Ethyl Benzene | 0.84 | 3.7 | Not Detected | Not Detected |
| m,p-Xylene | 0.84 | 3.7 | Not Detected | Not Detected |
| o-Xylene | 0.84 | 3.7 | Not Detected | Not Detected |
| Styrene | 0.84 | 3.6 | Not Detected | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.84 | 5.9 | Not Detected | Not Detected |
| 1,3,5-Trimethylbenzene | 0.84 | 4.2 | Not Detected | Not Detected |
| 1,2,4-Trimethylbenzene | 0.84 | 4.2 | Not Detected | Not Detected |
| 1,3-Dichlorobenzene | 0.84 | 5.1 | Not Detected | Not Detected |
| 1,4-Dichlorobenzene | 0.84 | 5.1 | Not Detected | Not Detected |
| Chlorotoluene | 0.84 | 4.4 | Not Detected | Not Detected |
| 1,2-Dichlorobenzene | 0.84 | 5.1 | Not Detected | Not Detected |
| 1,2,4-Trichlorobenzene | 0.84 | 6.3 | Not Detected | Not Detected |
| Hexachlorobutadiene | 0.84 | 9.1 | Not Detected | Not Detected |
| Propylene | 3.4 | 5.9 | Not Detected | Not Detected |
| 1,3-Butadiene | 3.4 | 7.6 | Not Detected | Not Detected |
| Acetone | 3.4 | 8.1 | 7.5 | 18 |

AIR TOXICS LTD.

SAMPLE NAME : AS-4

ID#: 0006175-05A

EPA METHOD TO-14 GC/MS Full Scan

| | | |
|--------------|---------|----------------------------|
| File Name: | C061923 | Date of Collection: 6/8/00 |
| Dil. Factor: | 1.68 | Date of Analysis: 6/20/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|----------------------------------|----------------------|-----------------------|------------------|-------------------|
| Carbon Disulfide | 3.4 | 11 | Not Detected | Not Detected |
| 2-Propanol | 3.4 | 8.4 | Not Detected | Not Detected |
| trans-1,2-Dichloroethene | 3.4 | 14 | Not Detected | Not Detected |
| Vinyl Acetate | 3.4 | 12 | Not Detected | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 3.4 | 10 | Not Detected | Not Detected |
| Hexane | 3.4 | 12 | Not Detected | Not Detected |
| Tetrahydrofuran | 3.4 | 10 | Not Detected | Not Detected |
| Cyclohexane | 3.4 | 12 | Not Detected | Not Detected |
| 1,4-Dioxane | 3.4 | 12 | Not Detected | Not Detected |
| Bromodichloromethane | 3.4 | 23 | Not Detected | Not Detected |
| 4-Methyl-2-pentanone | 3.4 | 14 | Not Detected | Not Detected |
| 2-Hexanone | 3.4 | 14 | Not Detected | Not Detected |
| Dibromochloromethane | 3.4 | 29 | Not Detected | Not Detected |
| Bromoform | 3.4 | 35 | Not Detected | Not Detected |
| 4-Ethyltoluene | 3.4 | 17 | Not Detected | Not Detected |
| Ethanol | 3.4 | 6.4 | Not Detected | Not Detected |
| Methyl tert-Butyl Ether | 3.4 | 12 | Not Detected | Not Detected |
| Heptane | 3.4 | 14 | Not Detected | Not Detected |

Container Type: 6 Liter Summa Canister

| Surrogates | % Recovery | Method Limits |
|-----------------------|------------|---------------|
| 1,2-Dichloroethane-d4 | 107 | 70-130 |
| Toluene-d8 | 100 | 70-130 |
| 4-Bromofluorobenzene | 92 | 70-130 |

AIR TOXICS LTD.

SAMPLE NAME : Blank

ID#: 0006175-06A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | c061924 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.00 | Date of Analysis: | 6/20/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|---------------------------|-------------------|--------------------|---------------|----------------|
| Freon 12 | 0.50 | 2.5 | Not Detected | Not Detected |
| Freon 114 | 0.50 | 3.6 | Not Detected | Not Detected |
| Chloromethane | 0.50 | 1.0 | Not Detected | Not Detected |
| Vinyl Chloride | 0.50 | 1.3 | Not Detected | Not Detected |
| Bromomethane | 0.50 | 2.0 | Not Detected | Not Detected |
| Chloroethane | 0.50 | 1.3 | Not Detected | Not Detected |
| Freon 11 | 0.50 | 2.8 | Not Detected | Not Detected |
| 1,1-Dichloroethene | 0.50 | 2.0 | Not Detected | Not Detected |
| Freon 113 | 0.50 | 3.9 | Not Detected | Not Detected |
| Methylene Chloride | 0.50 | 1.8 | Not Detected | Not Detected |
| 1,1-Dichloroethane | 0.50 | 2.0 | Not Detected | Not Detected |
| cis-1,2-Dichloroethene | 0.50 | 2.0 | Not Detected | Not Detected |
| Chloroform | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,1,1-Trichloroethane | 0.50 | 2.8 | Not Detected | Not Detected |
| Carbon Tetrachloride | 0.50 | 3.2 | Not Detected | Not Detected |
| Benzene | 0.50 | 1.6 | Not Detected | Not Detected |
| 1,2-Dichloroethane | 0.50 | 2.0 | Not Detected | Not Detected |
| Trichloroethene | 0.50 | 2.7 | Not Detected | Not Detected |
| 1,2-Dichloropropane | 0.50 | 2.3 | Not Detected | Not Detected |
| cis-1,3-Dichloropropene | 0.50 | 2.3 | Not Detected | Not Detected |
| Toluene | 0.50 | 1.9 | Not Detected | Not Detected |
| trans-1,3-Dichloropropene | 0.50 | 2.3 | Not Detected | Not Detected |
| 1,1,2-Trichloroethane | 0.50 | 2.8 | Not Detected | Not Detected |
| Tetrachloroethene | 0.50 | 3.4 | Not Detected | Not Detected |
| Ethylene Dibromide | 0.50 | 3.9 | Not Detected | Not Detected |
| Chlorobenzene | 0.50 | 2.3 | Not Detected | Not Detected |
| Ethyl Benzene | 0.50 | 2.2 | Not Detected | Not Detected |
| m,p-Xylene | 0.50 | 2.2 | Not Detected | Not Detected |
| o-Xylene | 0.50 | 2.2 | Not Detected | Not Detected |
| Styrene | 0.50 | 2.2 | Not Detected | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.50 | 3.5 | Not Detected | Not Detected |
| 1,3,5-Trimethylbenzene | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,2,4-Trimethylbenzene | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,3-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| 1,4-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| Chlorotoluene | 0.50 | 2.6 | Not Detected | Not Detected |
| 1,2-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| 1,2,4-Trichlorobenzene | 0.50 | 3.8 | Not Detected | Not Detected |
| Hexachlorobutadiene | 0.50 | 5.4 | Not Detected | Not Detected |
| Propylene | 2.0 | 3.5 | Not Detected | Not Detected |
| 1,3-Butadiene | 2.0 | 4.5 | Not Detected | Not Detected |
| Acetone | 2.0 | 4.8 | Not Detected | Not Detected |

AIR TOXICS LTD.

SAMPLE NAME : Blank

ID#: 0006175-06A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | C061924 | Date of Collection: | 6/8/00 |
| Dil. Factor: | 1.00 | Date of Analysis: | 6/20/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|----------------------------------|----------------------|-----------------------|------------------|-------------------|
| Carbon Disulfide | 2.0 | 6.3 | Not Detected | Not Detected |
| 2-Propanol | 2.0 | 5.0 | Not Detected | Not Detected |
| trans-1,2-Dichloroethene | 2.0 | 8.0 | Not Detected | Not Detected |
| Vinyl Acetate | 2.0 | 7.2 | Not Detected | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 2.0 | 6.0 | Not Detected | Not Detected |
| Hexane | 2.0 | 7.2 | Not Detected | Not Detected |
| Tetrahydrofuran | 2.0 | 6.0 | Not Detected | Not Detected |
| Cyclohexane | 2.0 | 7.0 | Not Detected | Not Detected |
| 1,4-Dioxane | 2.0 | 7.3 | Not Detected | Not Detected |
| Bromodichloromethane | 2.0 | 14 | Not Detected | Not Detected |
| 4-Methyl-2-pentanone | 2.0 | 8.3 | Not Detected | Not Detected |
| 2-Hexanone | 2.0 | 8.3 | Not Detected | Not Detected |
| Dibromochloromethane | 2.0 | 17 | Not Detected | Not Detected |
| Bromoform | 2.0 | 21 | Not Detected | Not Detected |
| 4-Ethyltoluene | 2.0 | 10 | Not Detected | Not Detected |
| Ethanol | 2.0 | 3.8 | Not Detected | Not Detected |
| Methyl tert-Butyl Ether | 2.0 | 7.3 | Not Detected | Not Detected |
| Heptane | 2.0 | 8.3 | Not Detected | Not Detected |

Container Type: 6 Liter Summa Canister

| Surrogates | % Recovery | Method Limits |
|-----------------------|------------|---------------|
| 1,2-Dichloroethane-d4 | 105 | 70-130 |
| Toluene-d8 | 99 | 70-130 |
| 4-Bromofluorobenzene | 91 | 70-130 |

AIR TOXICS LTD.

SAMPLE NAME : Lab Blank

ID#: 0006175-07A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|---------|---------------------|---------|
| File Name: | 0061909 | Date of Collection: | NA |
| Dil. Factor: | 1.00 | Date of Analysis: | 6/19/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|---------------------------|-------------------|--------------------|---------------|----------------|
| Freon 12 | 0.50 | 2.5 | Not Detected | Not Detected |
| Freon 114 | 0.50 | 3.6 | Not Detected | Not Detected |
| Chloromethane | 0.50 | 1.0 | Not Detected | Not Detected |
| Vinyl Chloride | 0.50 | 1.3 | Not Detected | Not Detected |
| Bromomethane | 0.50 | 2.0 | Not Detected | Not Detected |
| Chloroethane | 0.50 | 1.3 | Not Detected | Not Detected |
| Freon 11 | 0.50 | 2.8 | Not Detected | Not Detected |
| 1,1-Dichloroethene | 0.50 | 2.0 | Not Detected | Not Detected |
| Freon 113 | 0.50 | 3.9 | Not Detected | Not Detected |
| Methylene Chloride | 0.50 | 1.8 | Not Detected | Not Detected |
| 1,1-Dichloroethane | 0.50 | 2.0 | Not Detected | Not Detected |
| cis-1,2-Dichloroethene | 0.50 | 2.0 | Not Detected | Not Detected |
| Chloroform | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,1,1-Trichloroethane | 0.50 | 2.8 | Not Detected | Not Detected |
| Carbon Tetrachloride | 0.50 | 3.2 | Not Detected | Not Detected |
| Benzene | 0.50 | 1.6 | Not Detected | Not Detected |
| 1,2-Dichloroethane | 0.50 | 2.0 | Not Detected | Not Detected |
| Trichloroethene | 0.50 | 2.7 | Not Detected | Not Detected |
| 1,2-Dichloropropane | 0.50 | 2.3 | Not Detected | Not Detected |
| cis-1,3-Dichloropropene | 0.50 | 2.3 | Not Detected | Not Detected |
| Toluene | 0.50 | 1.9 | Not Detected | Not Detected |
| trans-1,3-Dichloropropene | 0.50 | 2.3 | Not Detected | Not Detected |
| 1,1,2-Trichloroethane | 0.50 | 2.8 | Not Detected | Not Detected |
| Tetrachloroethene | 0.50 | 3.4 | Not Detected | Not Detected |
| Ethylene Dibromide | 0.50 | 3.9 | Not Detected | Not Detected |
| Chlorobenzene | 0.50 | 2.3 | Not Detected | Not Detected |
| Ethyl Benzene | 0.50 | 2.2 | Not Detected | Not Detected |
| m,p-Xylene | 0.50 | 2.2 | Not Detected | Not Detected |
| o-Xylene | 0.50 | 2.2 | Not Detected | Not Detected |
| Styrene | 0.50 | 2.2 | Not Detected | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.50 | 3.5 | Not Detected | Not Detected |
| 1,3,5-Trimethylbenzene | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,2,4-Trimethylbenzene | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,3-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| 1,4-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| Chlorotoluene | 0.50 | 2.6 | Not Detected | Not Detected |
| 1,2-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| 1,2,4-Trichlorobenzene | 0.50 | 3.8 | Not Detected | Not Detected |
| Hexachlorobutadiene | 0.50 | 5.4 | Not Detected | Not Detected |
| Propylene | 2.0 | 3.5 | Not Detected | Not Detected |
| 1,3-Butadiene | 2.0 | 4.5 | Not Detected | Not Detected |
| Acetone | 2.0 | 4.8 | Not Detected | Not Detected |

AIR TOXICS LTD.

SAMPLE NAME : Lab Blank

ID#: 0006175-07A

EPA METHOD TO-14 GC/MS Full Scan

| | | | |
|--------------|--------|---------------------|---------|
| File Name: | 061909 | Date of Collection: | NA |
| Dil. Factor: | 1.00 | Date of Analysis: | 6/19/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|----------------------------------|-------------------|--------------------|---------------|----------------|
| Carbon Disulfide | 2.0 | 6.3 | Not Detected | Not Detected |
| 2-Propanol | 2.0 | 5.0 | Not Detected | Not Detected |
| trans-1,2-Dichloroethene | 2.0 | 8.0 | Not Detected | Not Detected |
| Vinyl Acetate | 2.0 | 7.2 | Not Detected | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 2.0 | 6.0 | Not Detected | Not Detected |
| Hexane | 2.0 | 7.2 | Not Detected | Not Detected |
| Tetrahydrofuran | 2.0 | 6.0 | Not Detected | Not Detected |
| Cyclohexane | 2.0 | 7.0 | Not Detected | Not Detected |
| 1,4-Dioxane | 2.0 | 7.3 | Not Detected | Not Detected |
| Bromodichloromethane | 2.0 | 14 | Not Detected | Not Detected |
| 4-Methyl-2-pentanone | 2.0 | 8.3 | Not Detected | Not Detected |
| 2-Hexanone | 2.0 | 8.3 | Not Detected | Not Detected |
| Dibromochloromethane | 2.0 | 17 | Not Detected | Not Detected |
| Bromoform | 2.0 | 21 | Not Detected | Not Detected |
| 4-Ethyltoluene | 2.0 | 10 | Not Detected | Not Detected |
| Ethanol | 2.0 | 3.8 | Not Detected | Not Detected |
| Methyl tert-Butyl Ether | 2.0 | 7.3 | Not Detected | Not Detected |
| Heptane | 2.0 | 8.3 | Not Detected | Not Detected |

Container Type: NA

| Surrogates | % Recovery | Method Limits |
|-----------------------|------------|---------------|
| 1,2-Dichloroethane-d4 | 102 | 70-130 |
| Toluene-d8 | 102 | 70-130 |
| 4-Bromofluorobenzene | 85 | 70-130 |

AIR TOXICS LTD.

SAMPLE NAME : Lab Blank

ID#: 0006175-07B

EPA METHOD TO-14 GC/MS Full Scan

File Name: g062011 Date of Collection: NA
 Dil Factor: 1:100 Date of Analysis: 6/20/00

| Compound | Det. Limit (ppbv) | Det. Limit (ug/m3) | Amount (ppbv) | Amount (ug/m3) |
|---------------------------|-------------------|--------------------|---------------|----------------|
| Freon 12 | 0.50 | 2.5 | Not Detected | Not Detected |
| Freon 114 | 0.50 | 3.6 | Not Detected | Not Detected |
| Chloromethane | 0.50 | 1.0 | Not Detected | Not Detected |
| Vinyl Chloride | 0.50 | 1.3 | Not Detected | Not Detected |
| Bromomethane | 0.50 | 2.0 | Not Detected | Not Detected |
| Chloroethane | 0.50 | 1.3 | Not Detected | Not Detected |
| Freon 11 | 0.50 | 2.8 | Not Detected | Not Detected |
| 1,1-Dichloroethene | 0.50 | 2.0 | Not Detected | Not Detected |
| Freon 113 | 0.50 | 3.9 | Not Detected | Not Detected |
| Methylene Chloride | 0.50 | 1.8 | 0.65 | 2.3 |
| 1,1-Dichloroethane | 0.50 | 2.0 | Not Detected | Not Detected |
| cis-1,2-Dichloroethene | 0.50 | 2.0 | Not Detected | Not Detected |
| Chloroform | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,1,1-Trichloroethane | 0.50 | 2.8 | Not Detected | Not Detected |
| Carbon Tetrachloride | 0.50 | 3.2 | Not Detected | Not Detected |
| Benzene | 0.50 | 1.6 | Not Detected | Not Detected |
| 1,2-Dichloroethane | 0.50 | 2.0 | Not Detected | Not Detected |
| Trichloroethene | 0.50 | 2.7 | Not Detected | Not Detected |
| 1,2-Dichloropropane | 0.50 | 2.3 | Not Detected | Not Detected |
| cis-1,3-Dichloropropene | 0.50 | 2.3 | Not Detected | Not Detected |
| Toluene | 0.50 | 1.9 | Not Detected | Not Detected |
| trans-1,3-Dichloropropene | 0.50 | 2.3 | Not Detected | Not Detected |
| 1,1,2-Trichloroethane | 0.50 | 2.8 | Not Detected | Not Detected |
| Tetrachloroethene | 0.50 | 3.4 | Not Detected | Not Detected |
| Ethylene Dibromide | 0.50 | 3.9 | Not Detected | Not Detected |
| Chlorobenzene | 0.50 | 2.3 | Not Detected | Not Detected |
| Ethyl Benzene | 0.50 | 2.2 | Not Detected | Not Detected |
| m,p-Xylene | 0.50 | 2.2 | Not Detected | Not Detected |
| o-Xylene | 0.50 | 2.2 | Not Detected | Not Detected |
| Styrene | 0.50 | 2.2 | Not Detected | Not Detected |
| 1,1,2,2-tetrachloroethane | 0.50 | 3.5 | Not Detected | Not Detected |
| 1,3,5-Trimethylbenzene | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,2,4-Trimethylbenzene | 0.50 | 2.5 | Not Detected | Not Detected |
| 1,3-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| 1,4-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| Chlorotoluene | 0.50 | 2.6 | Not Detected | Not Detected |
| 1,2-Dichlorobenzene | 0.50 | 3.0 | Not Detected | Not Detected |
| 1,2,4-Trichlorobenzene | 0.50 | 3.8 | Not Detected | Not Detected |
| Hexachlorobutadiene | 0.50 | 5.4 | Not Detected | Not Detected |
| Propylene | 2.0 | 3.5 | Not Detected | Not Detected |
| 1,3-Butadiene | 2.0 | 4.5 | Not Detected | Not Detected |
| Acetone | 2.0 | 4.8 | Not Detected | Not Detected |

AIR TOXICS LTD.

SAMPLE NAME : Lab Blank

ID#: 0006175-07B

EPA METHOD TO-14 GC/MS Full Scan

| | | |
|--------------|---------|---------------------------|
| File Name: | g062011 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 6/20/00 |

| Compound | Det. Limit (ppbv) | Det. Limit (uG/m3) | Amount (ppbv) | Amount (uG/m3) |
|----------------------------------|----------------------|-----------------------|------------------|-------------------|
| Carbon Disulfide | 2.0 | 6.3 | Not Detected | Not Detected |
| 2-Propanol | 2.0 | 5.0 | Not Detected | Not Detected |
| trans-1,2-Dichloroethene | 2.0 | 8.0 | Not Detected | Not Detected |
| Vinyl Acetate | 2.0 | 7.2 | Not Detected | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 2.0 | 6.0 | Not Detected | Not Detected |
| Hexane | 2.0 | 7.2 | Not Detected | Not Detected |
| Tetrahydrofuran | 2.0 | 6.0 | Not Detected | Not Detected |
| Cyclohexane | 2.0 | 7.0 | Not Detected | Not Detected |
| 1,4-Dioxane | 2.0 | 7.3 | Not Detected | Not Detected |
| Bromodichloromethane | 2.0 | 14 | Not Detected | Not Detected |
| 4-Methyl-2-pentanone | 2.0 | 8.3 | Not Detected | Not Detected |
| 2-Hexanone | 2.0 | 8.3 | Not Detected | Not Detected |
| Dibromochloromethane | 2.0 | 17 | Not Detected | Not Detected |
| Bromoform | 2.0 | 21 | Not Detected | Not Detected |
| 4-Ethyltoluene | 2.0 | 10 | Not Detected | Not Detected |
| Ethanol | 2.0 | 3.8 | Not Detected | Not Detected |
| Methyl tert-Butyl Ether | 2.0 | 7.3 | Not Detected | Not Detected |
| Heptane | 2.0 | 8.3 | Not Detected | Not Detected |

Container Type: NA

| Surrogates | % Recovery | Method Limits |
|-----------------------|------------|---------------|
| 1,2-Dichloroethane-d4 | 103 | 70-130 |
| Toluene-d8 | 102 | 70-130 |
| 4-Bromofluorobenzene | 91 | 70-130 |



AIR TOXICS LTD.

AN ENVIRONMENTAL ANALYTICAL LABORATORY

180 BLUE RAVINE ROAD, SUITE B
FOLSOM, CA 95630-4719
(916) 985-1000 FAX: (916) 985-1020

CHAIN-OF-CUSTODY RECORD

Nº 027595

Page 1 of 1

| | | |
|--|--|---|
| Contact Person <u>JOHN GAMMIE</u> Company <u>LFK</u> Address <u>5 JOHNSON DR</u> City <u>RARITAN</u> State <u>NJ</u> Zip <u>08869</u> Phone <u>860 721-7728</u> FAX <u>860 721-7665</u> Collected By: Signature <u>John Gammie</u> | Project Info: P.O. # _____ Project # <u>311-00105</u> Project Name <u>OLM</u> | Turn Around Time: <input checked="" type="checkbox"/> Normal <input type="checkbox"/> Rush _____ Specify _____ |
|--|--|---|

| Lab I.D. | Field Sample I.D. | Date & Time | Analyses Requested | Canister Pressure / Vacuum | | |
|----------|-------------------|---------------|--------------------|----------------------------|-------|---------|
| | | | | Initial | Final | Receipt |
| 0332 | AS-1 | 6/8/00 1042am | TO-14 list | 30.00 | 7.5 | |
| 0271 | AS-1 Dup | 6/8/00 1042am | " | 30.00 | 7.25 | |
| 2701 | AS-2 | 6/8/00 1023am | " | >30.00 | 7.0 | |
| 2507 | AS-3 | 6/8/00 1048am | " | 30.00 | 7.0 | |
| 0913 | AS-4 | 6/8/00 1054am | " | 29.90 | 6.9 | |
| 1034 | Blank | 6/8/00 1100am | " | 30.00 | 30.00 | |
| | | | | | | |
| | | | | | | |
| | | | | | | |

| | | |
|---|---|---|
| Relinquished By: (Signature) <u>John Gammie</u> Date/Time <u>6/9/00 1410hrs</u> Relinquished By: (Signature) _____ Date/Time _____ Relinquished By: (Signature) _____ Date/Time _____ | Print Name <u>JOHN GAMMIE</u> Received By: (Signature) _____ Date/Time _____ Received By: (Signature) _____ Date/Time _____ | Notes: please fax results to me e 860 721-7665 i send report to me e 94 Baneberry Lane Wethersfield, CT 06109 |
|---|---|---|

| | | | | | | | | |
|--------------|--------------|------------|------------|-----------|------------|-----------|-----------------------|--------------|
| Lab Use Only | Shipper Name | Air Bill # | Opened By: | Date/Time | Temp. (°C) | Condition | Custody Seals Intact? | Work Order # |
| | | | | | | | Yes No None N/A | |



AIR TOXICS LTD.
AN ENVIRONMENTAL ANALYTICAL LABORATORY

Sample Transportation Notice
Air Toxics Limited assumes no liability with respect to the collection, handling or shipment of these samples. Client represents and warrants that any sample delivered to LAB will be packaged or accompanied by completed written disclosure of presence of any hazardous substances known or suspected by client. Client further warrants that any sample containing any hazardous substance which is to be delivered to LAB will be packaged, labeled, transported and delivered properly and in accordance with applicable local, State, Federal, national, and international laws, regulations and ordinances of any kind. D.O.T. HAZMAT Hotline (800) 467-4922

180 BLUE RAVINE ROAD, SUITE B
FOLSOM, CA 95630-4719
(916) 985-1000 FAX: (916) 985-1020

CHAIN-OF-CUSTODY RECORD

No. 027595

Page 1 of 1

| | | |
|---|---|---|
| Contact Person <u>JOHN GANNINE</u> Company <u>LFK</u> Address <u>5 JOHNSON DR</u> City <u>RULITMAN</u> State <u>NY</u> Zip <u>02769</u> Phone <u>860 721-2727</u> FAX <u>860 721-7665</u> Collected By: Signature <u>John Gannine</u> | Project info: P.O. # _____ Project # <u>311-00105</u> Project Name <u>DM</u> | Turn Around Time: <input checked="" type="checkbox"/> Normal <input type="checkbox"/> Rush _____ Specify _____ |
|---|---|---|

LF R

| Lab ID | Field Sample I.D. | Date & Time | Analyses Requested | Canister Pressure / Vacuum | | |
|-------------|-------------------|----------------------|--------------------|----------------------------|--------------|---------------|
| | | | | Initial | Final | Receipt |
| <u>0332</u> | <u>AS-1</u> | <u>6/8/00 1042am</u> | <u>TO-14 list</u> | <u>30.00</u> | <u>7.5</u> | <u>6/8/00</u> |
| <u>0271</u> | <u>AS-1 Dup</u> | <u>6/8/00 1042am</u> | <u>"</u> | <u>30.00</u> | <u>7.25</u> | <u>6/8/00</u> |
| <u>2701</u> | <u>AS-2</u> | <u>6/8/00 1023am</u> | <u>"</u> | <u>>30.00</u> | <u>7.0</u> | <u>6/8/00</u> |
| <u>2507</u> | <u>AS-3</u> | <u>6/8/00 1048am</u> | <u>"</u> | <u>30.00</u> | <u>7.0</u> | <u>6/8/00</u> |
| <u>0915</u> | <u>AS-4</u> | <u>6/8/00 1054am</u> | <u>"</u> | <u>29.90</u> | <u>6.9</u> | <u>6/8/00</u> |
| <u>0231</u> | <u>Blank</u> | <u>6/8/00 1100am</u> | <u>"</u> | <u>30.00</u> | <u>30.00</u> | <u>6/8/00</u> |

page 20

| | |
|--|--|
| Relinquished By: (Signature) <u>John Gannine</u> Date/Time <u>6/8/00 14:00</u> Relinquished By: (Signature) _____ Date/Time _____ Relinquished By: (Signature) _____ Date/Time _____ | Received By: (Signature) <u>[Signature]</u> Date/Time <u>6-12-00 907</u> Received By: (Signature) _____ Date/Time _____ |
|--|--|

Notes:
 please Fax results to me e 860 721-7665
 I send report to me e
 94 Bareberry Lane
 Westford Field, CT 06109

| | | | | | |
|---|-----------------------------|---|--|---|-----------------------------|
| Supplier Name: _____ Lab Use Only: <u>Feb 15</u> | Air Eq. # <u>8983295192</u> | Operator: <u>ES</u> Date/Time: <u>6-12-00 207</u> | Temp. (°C): <u>25.2</u> Condition: <u>good</u> | Custody Seals Intact? Yes No <u>None</u> NA | Work Order # <u>0006175</u> |
|---|-----------------------------|---|--|---|-----------------------------|

9885267886

09:13

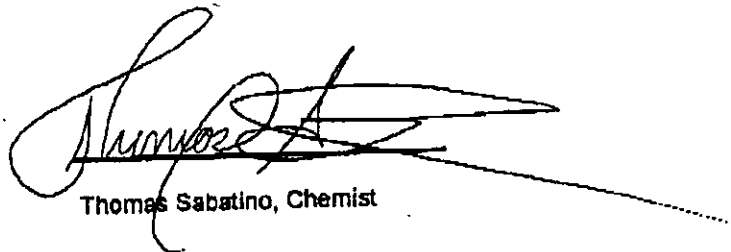
07/06/2000

11

AR06082000, C-7 TNMHCS

SUMMARY ANALYTICAL RESULTS AND CALIBRATION FOR THE ANALYSIS OF
NON-METHANE HYDROCARBONS AS HEPTANE IN GAS SAMPLES
RECEIVED 6/9/00, RUN 6/9/00, PROJECT # AR 311-00105-00-000

| SAMPLE IDENTIFICATION | COMPOUND | CONCENTRATION | UNITS |
|-----------------------|------------------------------|---------------|-------|
| INLET TO FLARE | NON-METHANE HC AS HEPTANE | 6280 | PPMV |



Thomas Sabatino, Chemist

Appendix G

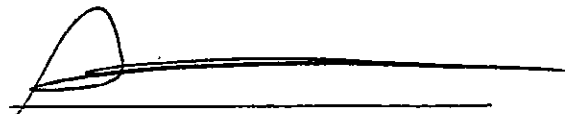
**Method 18 Flare Inlet Gas
Testing Report
AirRecon, Inc.**

**ANALYTICAL RESULTS AND SUPPORTING DOCUMENTS FOR THE ANALYSIS
OF TO-14 VOLATILE ORGANIC COMPOUNDS IN TEDLAR GAS SAMPLING
BAGS, USING GAS CHROMATOGRAPHY/ MASS SPECTROSCOPY DETECTION**

SUBMITTED BY: L-F-R, INC.
5 JOHNSON DRIVE
RARITAN, NEW JERSEY 08869
COLLECTED ON 8/14/2000
SUBMITTED ON 8/15/2000
PROJECT # 311-00142-00-000

ANALYTICAL RESEARCH AND TESTING, INC.
19 ROSS STREET
SOMERVILLE, N.J. 08876

SEPTEMBER 30, 2000



Thomas Sabatino, Director

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SECTION 1.0
ANALYTICAL RESULTS AND SUPPORTING DOCUMENTS FOR THE ANALYSIS
OF TO-14 VOLATILE ORGANIC COMPOUNDS IN TEDLAR GAS SAMPLING
BAGS, USING GAS CHROMATOGRAPHY/ MASS SPECTROSCOPY DETECTION

ANALYSIS OF VOLATILE ORGANIC COMPOUNDS COLLECTED IN TEDLAR SAMPLING
 BAGS, USING USEPA ANALYTICAL METHOD TO-14, MODIFIED, CAPILLARY
 COLUMN GAS CHROMATOGRAPHY/ MASS SPECTROSCOPY, VFRMT14H

| | | | | |
|-------------------|---------------|---------------|-----------|--------|
| CLIENT | L-F-R | TOTAL VOLUME | 60.00 | LITERS |
| SAMPLE # | LAB AIR BLANK | MATRIX | TEDLAR GB | |
| COLLECTED | 8/15/00 | GC/ MS # | VOC30-35 | |
| ARTI # | 00-464B | CC ON COLUMN | 0.50 | CC |
| DATE RECEIVED | 8/15/00 | DATE ANALYZED | 8/15/00 | |
| INITIAL FLOW RATE | 1.00 L/ MIN | ANALYST | T.S. | |
| FINAL FLOW RATE | 1.00 L/ MIN | REPORT DATE | 9/30/00 | |
| SAMPLING TIME | 60 MIN | | | |

| TARGET COMPOUND | NANOGRAMS/ SAMPLE FOUND CORR(2) | | ANALYTE VALUE | | |
|---------------------------|------------------------------------|------|---------------|----------|----|
| | | | MGS/M3 | PPBV (3) | |
| FREON 12 | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 20 |
| METHYLCHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 48 |
| FREON 114 | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 14 |
| VINYLCHELORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 39 |
| METHYLBROMIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 26 |
| ETHYLCHELORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 38 |
| FREON 11 | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 18 |
| VINYLDIENE CHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 25 |
| DICHLOROMETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 29 |
| FREON 113 | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 13 |
| CARBON DISULFIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 32 |
| 1,1-DICHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 25 |
| CIS-1,2-DICHLOROETHYLENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 25 |
| CHLOROFORM | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 20 |
| 1,2-DICHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 25 |
| METHYL CHLOROFORM | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 18 |
| BENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 31 |
| CARBON TETRACHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 16 |
| 1,2-DICHLOROPROPANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 22 |
| TRICHLOROETHYLENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| CIS-1,3-DICHLOROPROPENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 22 |
| TRANS-1,3-DICHLOROPROPENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 22 |
| 1,1,2-TRICHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 18 |
| TOLUENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 27 |
| 1,2-DIBROMOETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 13 |
| TETRACHLOROETHYLENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 13 |
| CHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 22 |
| ETHYLBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 23 |
| 1,3-XYLENE + 1,4-XYLENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 23 |
| STYRENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 24 |
| 1,1,2,2-TETRACHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 15 |
| 1,2-XYLENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 23 |

| TARGET COMPOUND | ANALYTE VALUE | | ANALYTE VALUE | | |
|-----------------------------------|---------------------|----------------|---------------|--------|----|
| | NANOGRAMS/ FOUND | SAMPLE CORR | MGS/M3 | PPBV | |
| 4-ETHYLTOLUENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 20 |
| 1,3,5-TRIMETHYLBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2,4-TRIMETHYLBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| 1,3-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| BENZYLCHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| 1,4-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2,4-TRICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 13 |
| NAPHTHALENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| HEXACHLOROBUTADIENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 9 |
| NON-TARGET COMPOUNDS FOUND | | | | | |
| TRANS-1,2-DICHLOROETHYLENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 25 |

SURROGATE COMPOUND RESPONSE
COMPOUND

| COMPOUND | % RECOVERY | ANALYTE VALUE | | |
|----------------------|------------|--------------------|------------------|------|
| | | NANOGRAMS FOUND | NLS (1) ADDED | |
| 4-BROMOFLUOROBENZENE | 98 | 5.90 | 6.00 | 0.82 |

(1) Milligrams/ M-3 = Milligrams per Cubic Meter Found in Sample. Corrected for Analyte Concentration in Trip Blank

(2) PPBV = Microliters/ cubic meter, Parts per Billion by Volume at 25.0 C Corrected for Blank Value.

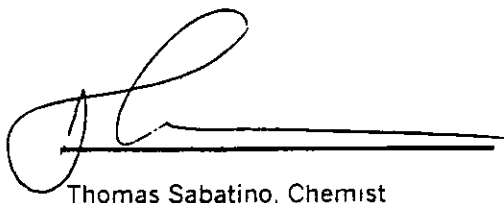
(3) N.D. = Non-Detectable, Less Than Concentration Indicated.

(4) A "J" Indicates an Estimated Value. This Flag is Used Either When Estimating a Concentration for a Tentatively Identified Compound or When a Given Compound is Suspected of Breakthrough and the Reported Value May be Biased Low.

(5) A "B" Indicates When the Analyte is Found in the Field Control as Well as the Sample. It Indicates Probable Contamination and Warns the Data User to Use This Result Cautiously.

TRIP BLANK CORRECTION VALUES

| TARGET COMPOUND | ANALYTE VALUE MILLIGRAMS/ M-3 FOUND |
|---------------------------|---|
| FREON 12 | 0.00 |
| METHYLCHLORIDE | 0.00 |
| FREON 114 | 0.00 |
| VINYLCHLORIDE | 0.00 |
| METHYLBROMIDE | 0.00 |
| ETHYLCHLORIDE | 0.00 |
| FREON 11 | 0.00 |
| ETHYL ETHER | 0.00 |
| VINYLDIENE CHLORIDE | 0.00 |
| DICHLOROMETHANE | 0.00 |
| FREON 113 | 0.00 |
| CARBONDISULFIDE | 0.00 |
| 1,1-DICHLOROETHANE | 0.00 |
| CIS-1,2-DICHLOROETHYLENE | 0.00 |
| CHLOROFORM | 0.00 |
| 1,2-DICHLOROETHANE | 0.00 |
| METHYL CHLOROFORM | 0.00 |
| BENZENE | 0.00 |
| CARBON TETRACHLORIDE | 0.00 |
| 1,2-DICHLOROPROPANE | 0.00 |
| TRICHLOROETHYLENE | 0.00 |
| CIS-1,3-DICHLOROPROPENE | 0.00 |
| TRANS-1,3-DICHLOROPROPENE | 0.00 |
| 1,1,2-TRICHLOROETHANE | 0.00 |
| TOLUENE | 0.00 |
| 1,2-DIBROMOETHANE | 0.00 |
| TETRACHLOROETHYLENE | 0.00 |
| CHLOROBENZENE | 0.00 |
| ETHYLBENZENE | 0.00 |
| 1,3-XYLENE + | |
| 1,4-XYLENE | 0.00 |
| STYRENE | 0.00 |
| 1,1,2,2-TETRACHLOROETHANE | 0.00 |
| 1,2-XYLENE | 0.00 |
| 4-ETHYLTOLUENE | 0.00 |
| 1,3,5-TRIMETHYLBENZENE | 0.00 |
| 1,2,4-TRIMETHYLBENZENE | 0.00 |
| 1,3-DICHLOROBENZENE | 0.00 |
| BENZYLCHLORIDE | 0.00 |
| 1,4-DICHLOROBENZENE | 0.00 |
| 1,2-DICHLOROBENZENE | 0.00 |
| 1,2,4-TRICHLOROBENZENE | 0.00 |
| NAPHTHALENE | 0.00 |
| HEXACHLOROBUTADIENE | 0.00 |



Thomas Sabatino, Chemist

ANALYSIS OF VOLATILE ORGANIC COMPOUNDS COLLECTED IN TEDLAR SAMPLING
 BAGS, USING USEPA ANALYTICAL METHOD TO-14, MODIFIED, CAPILLARY
 COLUMN GAS CHROMATOGRAPHY/ MASS SPECTROSCOPY, VFRMT14H

| | | | | |
|-------------------|-----------------------|---------------|-----------|--------|
| CLIENT | L-F-R | TOTAL VOLUME | 60.00 | LITERS |
| SAMPLE # | BAG # 1, LANDFILL GAS | MATRIX | TEDLAR GB | |
| COLLECTED | 8/15/00 | GC/ MS # | VOC30-36 | |
| ARTI # | 00-464 | CC ON COLUMN | 0.50 | CC |
| DATE RECEIVED | 8/15/00 | DATE ANALYZED | 8/15/00 | |
| INITIAL FLOW RATE | 1.00 L/ MIN | ANALYST | T.S. | |
| FINAL FLOW RATE | 1.00 L/ MIN | REPORT DATE | 9/30/00 | |
| SAMPLING TIME | 60 MIN | | | |

| TARGET COMPOUND | NANOGRAMS/ SAMPLE | | ANALYTE VALUE | |
|---------------------------|-------------------|---------|---------------|-----------|
| | FOUND | CORR(2) | MGS/M3 | PPBV (3) |
| FREON 12 | 0.55 | 0.55 | 1.09 | 220 |
| METHYLCHLORIDE | 1.38 | 1.38 | 2.76 | 1337 |
| FREON 114 | 0.55 | 0.55 | 1.09 | 156 |
| VINYLCHLORIDE | 2.15 | 2.15 | 4.29 | 1680 |
| METHYLBROMIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 26 |
| ETHYLCHLORIDE | 0.27 | 0.27 | 0.54 | 205 |
| FREON 11 | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 18 |
| VINYLDIENE CHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 25 |
| DICHLOROMETHANE | 1.19 | 1.19 | 2.37 | 682 |
| FREON 113 | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 13 |
| CARBON DISULFIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 32 |
| 1,1-DICHLOROETHANE | 1.02 | 1.02 | 2.04 | 504 |
| CIS-1,2-DICHLOROETHYLENE | 0.07 | 0.07 | 0.14 | 34 |
| CHLOROFORM | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 20 |
| 1,2-DICHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 25 |
| METHYL CHLOROFORM | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 18 |
| BENZENE | 0.43 | 0.43 | 0.85 | 267 |
| CARBON TETRACHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 16 |
| 1,2-DICHLOROPROPANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 22 |
| TRICHLOROETHYLENE | 0.94 | 0.94 | 1.88 | 349 |
| CIS-1,3-DICHLOROPROPENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 22 |
| TRANS-1,3-DICHLOROPROPENE | 0.05 | 0.05 | 0.10 | 22 |
| 1,1,2-TRICHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 18 |
| TOLUENE | 106 | 106 | 212 | 56300 |
| 1,2-DIBROMOETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 13 |
| TETRACHLOROETHYLENE | 1.88 | 1.88 | 3.75 | 488 |
| CHLOROBENZENE | 0.11 | 0.11 | 0.22 | 47 |
| ETHYLBENZENE | 8.09 | 8.09 | 16.18 | 3730 |
| 1,3-XYLENE + 1,4-XYLENE | 15.25 | 15.25 | 30.49 | 7030 |
| STYRENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 24 |
| 1,1,2,2-TETRACHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 15 |
| 1,2-XYLENE | 3.49 | 3.49 | 6.99 | 1610 |

| TARGET COMPOUND | ANALYTE VALUE | | ANALYTE VALUE | | |
|------------------------|---------------------|----------------|---------------|--------|-----|
| | NANOGRAMS/ FOUND | SAMPLE CORR | MGS/M3 | PPBV | |
| 4-ETHYLTOLUENE | 0.62 | 0.62 | 1.23 | | 251 |
| 1,3,5-TRIMETHYLBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2,4-TRIMETHYLBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| 1,3-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| BENZYLCHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| 1,4-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2,4-TRICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 13 |
| NAPHTHALENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| HEXACHLOROBUTADIENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 9 |

NON-TARGET COMPOUNDS FOUND

| | | | | | |
|----------------------------|------|------|------|--|------|
| TRANS-1,2-DICHLOROETHYLENE | 2.15 | 2.15 | 4.30 | | 1080 |
|----------------------------|------|------|------|--|------|

SURROGATE COMPOUND RESPONSE

| COMPOUND | % RECOVERY | ANALYTE VALUE | | |
|----------------------|------------|--------------------|------------------|------|
| | | NANOGRAMS FOUND | NLS (1) ADDED | |
| 4-BROMOFLUOROBENZENE | 94 | 5.66 | 6.00 | 0.79 |

(1) Milligrams/ M-3 = Milligrams per Cubic Meter Found in Sample. Corrected for Analyte Concentration in Trip Blank

(2) PPBV = Microliters/ cubic meter, Parts per Billion by Volume at 25.0 C Corrected for Blank Value.

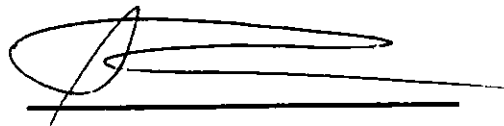
(3) N.D. = Non-Detectable, Less Than Concentration Indicated.

(4) A "J" Indicates an Estimated Value. This Flag is Used Either When Estimating a Concentration for a Tentatively Identified Compound or When a Given Compound is Suspected of Breakthrough and the Reported Value May be Biased Low.

(5) A "B" Indicates When the Analyte is Found in the Field Control as Well as the Sample. It Indicates Probable Contamination and Warns the Data User to Use This Result Cautiously.

TRIP BLANK CORRECTION VALUES

| TARGET COMPOUND | ANALYTE VALUE MILLIGRAMS/ M-3 FOUND |
|---------------------------|---|
| FREON 12 | 0.00 |
| METHYLCHLORIDE | 0.00 |
| FREON 114 | 0.00 |
| VINYLCHLORIDE | 0.00 |
| METHYLBROMIDE | 0.00 |
| ETHYLCHLORIDE | 0.00 |
| FREON 11 | 0.00 |
| ETHYL ETHER | 0.00 |
| VINYLDIENE CHLORIDE | 0.00 |
| DICHLOROMETHANE | 0.00 |
| FREON 113 | 0.00 |
| CARBONDISULFIDE | 0.00 |
| 1,1-DICHLOROETHANE | 0.00 |
| CIS-1,2-DICHLOROETHYLENE | 0.00 |
| CHLOROFORM | 0.00 |
| 1,2-DICHLOROETHANE | 0.00 |
| METHYL CHLOROFORM | 0.00 |
| BENZENE | 0.00 |
| CARBON TETRACHLORIDE | 0.00 |
| 1,2-DICHLOROPROPANE | 0.00 |
| TRICHLOROETHYLENE | 0.00 |
| CIS-1,3-DICHLOROPROPENE | 0.00 |
| TRANS-1,3-DICHLOROPROPENE | 0.00 |
| 1,1,2-TRICHLOROETHANE | 0.00 |
| TOLUENE | 0.00 |
| 1,2-DIBROMOETHANE | 0.00 |
| TETRACHLOROETHYLENE | 0.00 |
| CHLOROBENZENE | 0.00 |
| ETHYLBENZENE | 0.00 |
| 1,3-XYLENE + | |
| 1,4-XYLENE | 0.00 |
| STYRENE | 0.00 |
| 1,1,2,2-TETRACHLOROETHANE | 0.00 |
| 1,2-XYLENE | 0.00 |
| 4-ETHYLTOLUENE | 0.00 |
| 1,3,5-TRIMETHYLBENZENE | 0.00 |
| 1,2,4-TRIMETHYLBENZENE | 0.00 |
| 1,3-DICHLOROBENZENE | 0.00 |
| BENZYLCHLORIDE | 0.00 |
| 1,4-DICHLOROBENZENE | 0.00 |
| 1,2-DICHLOROBENZENE | 0.00 |
| 1,2,4-TRICHLOROBENZENE | 0.00 |
| NAPHTHALENE | 0.00 |
| HEXACHLOROBUTADIENE | 0.00 |



Thomas Sabatino, Chemist

ANALYSIS OF VOLATILE ORGANIC COMPOUNDS COLLECTED IN TEDLAR SAMPLING
 BAGS, USING USEPA ANALYTICAL METHOD TO-14, MODIFIED, CAPILLARY
 COLUMN GAS CHROMATOGRAPHY/ MASS SPECTROSCOPY, VFRMT14H

| | | | | | |
|-------------------|----------|--------------|---------------|-----------|--------|
| CLIENT | L-F-R | DUPLICATE | TOTAL VOLUME | 60.00 | LITERS |
| SAMPLE # | BAG # 1, | LANDFILL GAS | MATRIX | TEDLAR GB | |
| COLLECTED | | 8/15/00 | GC/ MS # | VOC30-37 | |
| ARTI # | | 00-464D | CC ON COLUMN | 0.50 | CC |
| DATE RECEIVED | | 8/15/00 | DATE ANALYZED | 8/15/00 | |
| INITIAL FLOW RATE | 1.00 | L/ MIN | ANALYST | T.S. | |
| FINAL FLOW RATE | 1.00 | L/ MIN | REPORT DATE | 9/30/00 | |
| SAMPLING TIME | 60 | MIN | | | |

| TARGET COMPOUND | NANOGRAMS/ SAMPLE | | ANALYTE VALUE | | |
|---------------------------|-------------------|---------|---------------|-----------|--|
| | FOUND | CORR(2) | MGS/M3 | PPBV (3) | |
| FREON 12 | 0.57 | 0.57 | 1.14 | 230 | |
| METHYLCHLORIDE | 1.45 | 1.45 | 2.90 | 1400 | |
| FREON 114 | 0.57 | 0.57 | 1.14 | 163 | |
| VINYLCHLORIDE | 2.03 | 2.03 | 4.05 | 1590 | |
| METHYLBROMIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 26 | |
| ETHYLCHLORIDE | 0.22 | 0.22 | 0.43 | 165 | |
| FREON 11 | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 18 | |
| VINYLDIENE CHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 25 | |
| DICHLOROMETHANE | 1.12 | 1.12 | 2.23 | 642 | |
| FREON 113 | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 13 | |
| CARBON DISULFIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 32 | |
| 1,1-DICHLOROETHANE | 0.98 | 0.98 | 1.95 | 482 | |
| CIS-1,2-DICHLOROETHYLENE | 0.06 | 0.06 | 0.12 | 31 | |
| CHLOROFORM | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 20 | |
| 1,2-DICHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 25 | |
| METHYL CHLOROFORM | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 18 | |
| BENZENE | 0.38 | 0.38 | 0.77 | 240 | |
| CARBON TETRACHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 16 | |
| 1,2-DICHLOROPROPANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 22 | |
| TRICHLOROETHYLENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 19 | |
| CIS-1,3-DICHLOROPROPENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 22 | |
| TRANS-1,3-DICHLOROPROPENE | 0.05 | 0.05 | 0.10 | 21 | |
| 1,1,2-TRICHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 18 | |
| TOLUENE | 96.53 | 96.53 | 193 | 51300 | |
| 1,2-DIBROMOETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 13 | |
| TETRACHLOROETHYLENE | 1.72 | 1.72 | 3.44 | 448 | |
| CHLOROBENZENE | 0.10 | 0.10 | 0.20 | 44 | |
| ETHYLBENZENE | 8.56 | 8.56 | 17.13 | 3950 | |
| 1,3-XYLENE + 1,4-XYLENE | 15.78 | 15.78 | 31.56 | 7280 | |
| STYRENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 24 | |
| 1,1,2,2-TETRACHLOROETHANE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < 15 | |
| 1,2-XYLENE | 3.68 | 3.68 | 7.36 | 1700 | |

| TARGET COMPOUND | ANALYTE VALUE | | ANALYTE VALUE | | |
|------------------------|---------------------|----------------|---------------|--------|-----|
| | NANOGRAMS/ FOUND | SAMPLE CORR | MGS/M3 | PPBV | |
| 4-ETHYLTOLUENE | 0.82 | 0.82 | 1.64 | | 335 |
| 1,3,5-TRIMETHYLBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2,4-TRIMETHYLBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| 1,3-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| BENZYLCHLORIDE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| 1,4-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2-DICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 17 |
| 1,2,4-TRICHLOROBENZENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 13 |
| NAPHTHALENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 19 |
| HEXACHLOROBUTADIENE | ND < 0.05 | 0.00 | ND < 0.05 | N.D. < | 9 |

NON-TARGET COMPOUNDS FOUND

| | | | | | |
|----------------------------|------|------|------|--|------|
| TRANS-1,2-DICHLOROETHYLENE | 2.02 | 2.02 | 4.04 | | 1020 |
|----------------------------|------|------|------|--|------|

SURROGATE COMPOUND RESPONSE

| COMPOUND | % RECOVERY | ANALYTE VALUE | | NLS (1) |
|----------------------|------------|--------------------|-------|---------|
| | | NANOGRAMS FOUND | ADDED | |
| 4-BROMOFLUOROBENZENE | 101 | 6.05 | 6.00 | 0.85 |

(1) Milligrams/ M-3 = Milligrams per Cubic Meter Found in Sample. Corrected for Analyte Concentration in Trip Blank

(2) PPBV = Microliters/ cubic meter, Parts per Billion by Volume at 25.0 C Corrected for Blank Value.

(3) N.D. = Non-Detectable, Less Than Concentration Indicated

(4) A "J" Indicates an Estimated Value. This Flag is Used Either When Estimating a Concentration for a Tentatively Identified Compound or When a Given Compound is Suspected of Breakthrough and the Reported Value May be Biased Low.

(5) A "B" Indicates When the Analyte is Found in the Field Control as Well as the Sample. It Indicates Probable Contamination and Warns the Data User to Use This Result Cautiously.

TRIP BLANK CORRECTION VALUES

| TARGET COMPOUND | ANALYTE VALUE MILLIGRAMS/ M-3 FOUND |
|---------------------------|---|
| FREON 12 | 0.00 |
| METHYLCHLORIDE | 0.00 |
| FREON 114 | 0.00 |
| VINYLCHLORIDE | 0.00 |
| METHYLBROMIDE | 0.00 |
| ETHYLCHLORIDE | 0.00 |
| FREON 11 | 0.00 |
| ETHYL ETHER | 0.00 |
| VINYLDIENE CHLORIDE | 0.00 |
| DICHLOROMETHANE | 0.00 |
| FREON 113 | 0.00 |
| CARBONDISULFIDE | 0.00 |
| 1,1-DICHLOROETHANE | 0.00 |
| CIS-1,2-DICHLOROETHYLENE | 0.00 |
| CHLOROFORM | 0.00 |
| 1,2-DICHLOROETHANE | 0.00 |
| METHYL CHLOROFORM | 0.00 |
| BENZENE | 0.00 |
| CARBON TETRACHLORIDE | 0.00 |
| 1,2-DICHLOROPROPANE | 0.00 |
| TRICHLOROETHYLENE | 0.00 |
| CIS-1,3-DICHLOROPROPENE | 0.00 |
| TRANS-1,3-DICHLOROPROPENE | 0.00 |
| 1,1,2-TRICHLOROETHANE | 0.00 |
| TOLUENE | 0.00 |
| 1,2-DIBROMOETHANE | 0.00 |
| TETRACHLOROETHYLENE | 0.00 |
| CHLOROBENZENE | 0.00 |
| ETHYLBENZENE | 0.00 |
| 1,3-XYLENE + | |
| 1,4-XYLENE | 0.00 |
| STYRENE | 0.00 |
| 1,1,2,2-TETRACHLOROETHANE | 0.00 |
| 1,2-XYLENE | 0.00 |
| 4-ETHYLTOLUENE | 0.00 |
| 1,3,5-TRIMETHYLBENZENE | 0.00 |
| 1,2,4-TRIMETHYLBENZENE | 0.00 |
| 1,3-DICHLOROBENZENE | 0.00 |
| BENZYLCHLORIDE | 0.00 |
| 1,4-DICHLOROBENZENE | 0.00 |
| 1,2-DICHLOROBENZENE | 0.00 |
| 1,2,4-TRICHLOROBENZENE | 0.00 |
| NAPHTHALENE | 0.00 |
| HEXACHLOROBUTADIENE | 0.00 |



Thomas Sabatino, Chemist

SECTION 2.0
TARGET COMPOUND RETENTION TIME TABLES FOR THE ANALYSIS
OF TO-14 VOLATILE ORGANIC COMPOUNDS IN GAS SAMPLES

Quantitation Report Quanfile: VOC30-29
 Comment: TO-14 VOC STANDARD 1.5 NGS0 + 4-BFB
 Sorted via: Entry Number ↑

Quan Entries: 45

(S) = Standard

| Cal | Name of Compound | S | Scan# | R Time | Me | Calc Amt(A) | Units |
|-----|------------------------|---|-------|--------|----|-------------|-------|
| 1 | Freon 12 | E | 88 | 1:32 | BB | 1.403 | NANOS |
| 2 | Methyl Chloride | E | 94 | 1:39 | VV | 1.709 | NANOS |
| 3 | Freon 114 | E | 88 | 1:32 | BB | 1.403 | NANOS |
| 4 | Vinyl Chloride | E | 102 | 1:47 | BB | 1.394 | NANOS |
| 5 | Methyl Bromide | E | 117 | 2:03 | BB | 1.390 | NANOS |
| 6 | Ethyl Chloride | E | 123 | 2:09 | BB | 1.392 | NANOS |
| 7 | Freon 11 | E | 151 | 2:39 | BB | 1.393 | NANOS |
| 8 | Ethyl Ether | E | 165 | 2:53 | BB | 1.397 | NANOS |
| 9 | 1,1-Dichloroethene | E | 178 | 3:07 | BB | 1.388 | NANOS |
| 10 | Methylene Chloride | E | 182 | 3:11 | BV | 1.394 | NANOS |
| 11 | Freon 113 | E | 195 | 3:25 | BB | 1.395 | NANOS |
| 12 | Carbon Disulfide | E | 198 | 3:28 | BB | 1.399 | NANOS |
| 13 | Cis-1,2-Dichloroethene | E | 229 | 4:00 | BV | 1.396 | NANOS |
| 14 | 1,1-Dichloroethane | E | 240 | 4:12 | VB | 1.395 | NANOS |
| 15 | Trans-1,2-Dichloroethe | E | 287 | 5:01 | BB | 1.393 | NANOS |
| 16 | Chloroform | E | 305 | 5:20 | BB | 1.394 | NANOS |
| 17 | 1,2-Dichloroethane | E | 350 | 6:07 | BV | 1.397 | NANOS |
| 18 | 1,1,1-Trichloroethane | E | 366 | 6:24 | BB | 1.390 | NANOS |
| 19 | Benzene | E | 395 | 6:55 | BB | 1.393 | NANOS |
| 20 | Carbontetrachloride | E | 405 | 7:05 | BB | 1.392 | NANOS |
| 21 | 1,2-Dichloropropane | E | 450 | 7:52 | BB | 1.393 | NANOS |
| 22 | Trichloroethene | E | 466 | 8:09 | BB | 1.397 | NANOS |
| 23 | Cis-1,3-Dichloropropen | E | 527 | 9:13 | BB | 1.396 | NANOS |
| 24 | Trans-1,3-Dichloroprop | E | 566 | 9:54 | BB | 1.397 | NANOS |
| 25 | 1,1,2-Trichloroethane | E | 578 | 10:07 | BB | 1.390 | NANOS |
| 26 | Toluene | E | 599 | 10:29 | BB | 1.398 | NANOS |

Quantitation Report Quanfile: VOC30-29
 Comment: T0-14 VOC STANDARD 1.5 NGSQ + 4-BFB
 Sorted via: Entry Number ↑

Quan Entries: 45

(S) = Standard

| Cal | Name of Compound | S | Scan# | R Time | Me | Calc Amt(A) | Units |
|-----|------------------------|---|-------|--------|----|-------------|-------|
| 27 | 1,2-Dibromoethane | E | 649 | 11:21 | BB | 1.400 | NANOS |
| 28 | Tetrachloroethene | E | 687 | 12:01 | BB | 1.397 | NANOS |
| 29 | Chlorobenzene | E | 742 | 12:59 | BB | 1.406 | NANOS |
| 30 | Ethylbenzene | E | 775 | 13:34 | BB | 1.406 | NANOS |
| 31 | 1,3+1,4-Xylenes | E | 790 | 13:49 | BB | 2.828 | NANOS |
| 32 | Styrene | E | 822 | 14:23 | BV | 1.412 | NANOS |
| 33 | 1,1,2,2-Tetrachloroeth | E | 830 | 14:31 | BB | 1.406 | NANOS |
| 34 | 1,2-Xylene | E | 831 | 14:32 | BB | 1.406 | NANOS |
| 35 | 4-Bromofluorobenzene | E | 872 | 15:16 | BB | 1.412 | NANOS |
| 36 | Benzylchloride | E | 930 | 16:16 | BB | 1.405 | NANOS |
| 37 | 4-Ethyltoluene | E | 950 | 16:38 | BV | 1.451 | NANOS |
| 38 | 1,3,5-Trimethylbenzene | E | 959 | 16:47 | VV | 1.422 | NANOS |
| 39 | 1,2,4-Trimethylbenzene | E | 1001 | 17:31 | BB | 1.488 | NANOS |
| 40 | 1,3-Dichlorobenzene | E | 1015 | 17:46 | BV | 1.584 | NANOS |
| 41 | 1,4-Dichlorobenzene | E | 1023 | 17:54 | MM | 1.375 | NANOS |
| 42 | 1,2-Dichlorobenzene | E | 1060 | 18:33 | BB | 1.823 | NANOS |
| 43 | 1,2,4-Trichlorobenzene | E | 1276 | 22:20 | BB | 1.591 | NANOS |
| 44 | Naphthalene | E | 1289 | 22:33 | BB | 1.676 | NANOS |
| 45 | Hexachlorobutadiene | E | 1341 | 23:28 | BB | 1.545 | NANOS |

SECTION 3.0
GC/ MS OPERATING CONDITIONS FOR THE ANALYSIS OF TO-14 VOLATILE
ORGANIC COMPOUNDS IN GAS SAMPLES

Log Information

Comment: TO-14 VOC STANDARD
Comment: T. SABATINO

File: D:\VOC30-29

Date: Aug-15-2000 14:00:29

1.5 NGS0 + 4-BFB

VOLATILE ORGANIC COMPOUND ANALYSIS

Acquisition Method Table

| | |
|----------------------------|-----------------------|
| Instrument set points | |
| Multiplier set voltage | 1700 volts |
| Manifold set temperature | 220 °C |
| Emmision set current | 10 micro amps |
| A/M amplitude set voltage | 4.0 volts |
| User cancelled acquisition | no |
| End instrument set points | |
| | |
| Acquire segment # | 1 |
| Target value | 10000 |
| Low mass | 49 amu |
| High mass | 300 amu |
| Scan rate | 1050 milli seconds |
| Segment acquire time | 28 minutes |
| Threshold | 1 counts |
| Filament delay | 10 seconds |
| Mass defect | 50 milli mass/100 amu |
| Background mass | 48 amu |
| Calibration gas | no |
| Scan mode | EI |
| Ionization control | automatic |
| End Acquire segment # | 1 |

↑-Prev seg ↓-Next seg

Esc to continue

Log Information

Comment: TO-14 VOC STANDARD
Comment: T. SABATINO

File: D:\VOC30-29

Date: Aug-15-2000 14:00:29

1.5 NGS0 + 4-BFB

VOLATILE ORGANIC COMPOUND ANALYSIS

Title Saturn Log File

Analysis list file name VOC30
Acquisition method file name T014FG
GC method file name T014G
Autosampler method file name

Column Table

Column tables

| Start temp °C | End temp °C | rate °C/minute | time minutes |
|------------------|----------------|-------------------|-----------------|
| 40 | 40 | 0.0 | 4.00 |
| 40 | 180 | 7.0 | 20.00 |
| 180 | 180 | 0.0 | 4.00 |

End column tables

Esc to continue

F7 - Display GC transfer line table

F8 - Display GC valve parameters

F9 - Display GC error table

F10 - User log notes and comments editor

? - Help

Esc - Exit

Log Information File: D:\VOC30-29 Date: Aug-15-2000 14:00:29
Comment: T0-14 VOC STANDARD 1.5 NGS0 + 4-BFB
Comment: T. SABATINO VOLATILE ORGANIC COMPOUND ANALYSIS

Title Saturn Log File

Analysis list file name VOC30
Acquisition method file name T014FG
GC method file name T014G
Autosampler method file name

Injector Table

Injector tables

| Start temp | End temp | rate | time |
|------------|----------|-----------|---------|
| °C | °C | °C/minute | minutes |
| 150 | 150 | 0.0 | 28.00 |

End injector tables

Esc to continue

F6 - Display GC injector table
F7 - Display GC transfer line table
F8 - Display GC valve parameters
F9 - Display GC error table

F10 - User log notes and comments editor

? - Help
Esc - Exit

Log Information

Comment: T0-14 VOC STANDARD

Comment: T. SABATINO

File: D:\VOC30-29

1.5 NGS0 + 4-BFB

VOLATILE ORGANIC COMPOUND ANALYSIS

Date: Aug-15-2000 14:00:29

Title Saturn Log File

Analysis list file name VOC30
Acquisition method file name T014FG
GC method file name T014G
Autosampler method file name

Transfer Line Table

Transfer line tables

| Start temp | End temp | rate | time |
|------------|----------|-----------|---------|
| °C | °C | °C/minute | minutes |
| 200 | 200 | 0.0 | 4.00 |

End transfer line tables

Esc to continue

F6 - Display GC injector table
F7 - Display GC transfer line table
F8 - Display GC valve parameters
F9 - Display GC error table

F10 - User log notes and comments editor

? - Help
Esc - Exit

SECTION 4.0
GC/ MS RAW DATA REPORTS FOR THE ANALYSIS OF TO-14 VOLATILE
ORGANIC COMPOUNDS IN GAS SAMPLES

Quantitation Report Quanfile: UOC30-35
 Comment: LAB AIR BLANK 0.50 CC + 4-BFB
 Sorted via: Entry Number ↑

Quan Entries: 15

(S) = Standard

| Cal | Name of Compound | S | Scan# | R Time | Me | Calc Amt(A) | Units |
|-----|------------------------|---|-------|--------|----|-------------|-------|
| 2 | Methyl Chloride | E | 105 | 1:50 | BV | 0.000 | NANOS |
| 10 | Methylene Chloride | E | 182 | 3:11 | BB | 0.000 | NANOS |
| 11 | Freon 113 | E | 196 | 3:26 | BB | 0.000 | NANOS |
| 12 | Carbon Disulfide | E | 198 | 3:28 | BB | 0.000 | NANOS |
| 16 | Chloroform | E | 306 | 5:21 | BB | 0.000 | NANOS |
| 19 | Benzene | E | 394 | 6:54 | BB | 0.000 | NANOS |
| 26 | Toluene | E | 599 | 10:29 | BB | 0.000 | NANOS |
| 30 | Ethylbenzene | E | 778 | 13:37 | BB | 0.000 | NANOS |
| 31 | 1,3+1,4-Xylenes | E | 778 | 13:37 | BB | 0.000 | NANOS |
| 35 | 4-Bromofluorobenzene | E | 872 | 15:16 | BB | 5.902 | NANOS |
| 37 | 4-Ethyltoluene | E | 951 | 16:39 | BB | 0.000 | NANOS |
| 38 | 1,3,5-Trimethylbenzene | E | 951 | 16:39 | BB | 0.000 | NANOS |
| 39 | 1,2,4-Trimethylbenzene | E | 1001 | 17:31 | BB | 0.000 | NANOS |
| 43 | 1,2,4-Trichlorobenzene | E | 1278 | 22:22 | BB | 0.000 | NANOS |
| 44 | Naphthalene | E | 1291 | 22:35 | BB | 0.000 | NANOS |

Quantitation Report

Quanfile: UOC30-36

Quan Entries: 34

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB

Sorted via: Entry Number ↑

(S) = Standard

| Cal | Name of Compound | S | Scan# | R Time | Me | Calc Amt(A) | Units |
|-----|--------------------------|---|-------|--------|----|-------------|-------|
| 1 | Freon 12 | E | 87 | 1:31 | BV | 0.545 | NANOS |
| 2 | Methyl Chloride | E | 105 | 1:50 | VV | 1.381 | NANOS |
| 3 | Freon 114 | E | 87 | 1:31 | BV | 0.545 | NANOS |
| 4 | Vinyl Chloride | E | 101 | 1:46 | BB | 2.147 | NANOS |
| 6 | Ethyl Chloride | E | 125 | 2:11 | BB | 0.270 | NANOS |
| 8 | Ethyl Ether | E | 176 | 3:05 | BB | 3.361 | NANOS |
| 10 | Methylene Chloride | E | 182 | 3:11 | VB | 1.185 | NANOS |
| 11 | Freon 113 | E | 195 | 3:25 | BB | 0.000 | NANOS |
| 12 | Carbon Disulfide | E | 198 | 3:28 | BB | 0.000 | NANOS |
| 13 | Cis-1,2-Dichloroethene | E | 229 | 4:00 | BV | 0.068 | NANOS |
| 14 | 1,1-Dichloroethane | E | 240 | 4:12 | VB | 1.020 | NANOS |
| 15 | Trans-1,2-Dichloroethane | E | 287 | 5:01 | VV | 2.149 | NANOS |
| 18 | 1,1,1-Trichloroethane | E | 366 | 6:24 | BB | 0.000 | NANOS |
| 19 | Benzene | E | 395 | 6:55 | BB | 0.426 | NANOS |
| 21 | 1,2-Dichloropropane | E | 449 | 7:51 | VB | 0.000 | NANOS |
| 22 | Trichloroethene | E | 466 | 8:09 | BB | 0.938 | NANOS |
| 24 | Trans-1,3-Dichloroprop | E | 565 | 9:53 | BB | 0.050 | NANOS |
| 25 | 1,1,2-Trichloroethane | E | 574 | 10:03 | BB | 0.000 | NANOS |
| 26 | Toluene | E | 599 | 10:29 | BB | 106.101 | NANOS |
| 28 | Tetrachloroethene | E | 687 | 12:01 | BB | 1.876 | NANOS |
| 29 | Chlorobenzene | E | 742 | 12:59 | BB | 0.108 | NANOS |
| 30 | Ethylbenzene | E | 774 | 13:33 | BB | 0.089 | NANOS |
| 31 | 1,3+1,4-Xylenes | E | 789 | 13:48 | BB | 15.247 | NANOS |
| 33 | 1,1,2,2-Tetrachloroeth | E | 814 | 14:15 | VV | 0.000 | NANOS |
| 34 | 1,2-Xylene | E | 831 | 14:33 | BV | 3.494 | NANOS |
| 35 | 4-Bromofluorobenzene | E | 871 | 15:15 | BV | 5.659 | NANOS |

Quantitation Report

Quanfile: VOC30-36

Quan Entries: 34

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB

Sorted via: Entry Number ↑

(S) = Standard

| Cal | Name of Compound | S | Scan# | R Time | Me | Calc Amt(A) | Units |
|-----|------------------------|---|-------|--------|----|-------------|-------|
| 36 | Benzylchloride | E | 935 | 16:22 | VU | 0.000 | NANOS |
| 37 | 4-Ethyltoluene | E | 946 | 16:33 | VU | 0.617 | NANOS |
| 38 | 1,3,5-Trimethylbenzene | E | 959 | 16:47 | VB | 0.000 | NANOS |
| 39 | 1,2,4-Trimethylbenzene | E | 1001 | 17:31 | BB | 0.000 | NANOS |
| 40 | 1,3-Dichlorobenzene | E | 1023 | 17:54 | BB | 0.000 | NANOS |
| 41 | 1,4-Dichlorobenzene | E | 1023 | 17:54 | BB | 0.000 | NANOS |
| 43 | 1,2,4-Trichlorobenzene | E | 1275 | 22:19 | BB | 0.000 | NANOS |
| 44 | Naphthalene | E | 1289 | 22:33 | BB | 0.000 | NANOS |

Quantitation Report

Quanfile: UOC38-37

Quan Entries: 34

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB DUPLICATE

Sorted via: Entry Number ↑

(S) = Standard

| Cal | Name of Compound | S | Scan# | R Time | Me | Calc Amt(A) | Units |
|-----|--------------------------|---|-------|--------|----|-------------|-------|
| 1 | Freon 12 | E | 87 | 1:31 | BV | 0.570 | NANOS |
| 2 | Methyl Chloride | E | 105 | 1:50 | VV | 1.452 | NANOS |
| 3 | Freon 114 | E | 87 | 1:31 | BV | 0.570 | NANOS |
| 4 | Vinyl Chloride | E | 101 | 1:46 | BB | 2.026 | NANOS |
| 6 | Ethyl Chloride | E | 124 | 2:10 | BB | 0.217 | NANOS |
| 7 | Freon 11 | E | 151 | 2:39 | BB | 0.001 | NANOS |
| 8 | Ethyl Ether | E | 176 | 3:05 | BB | 3.151 | NANOS |
| 10 | Methylene Chloride | E | 182 | 3:11 | VB | 1.115 | NANOS |
| 11 | Freon 113 | E | 193 | 3:23 | BB | 0.000 | NANOS |
| 12 | Carbon Disulfide | E | 197 | 3:27 | BB | 0.000 | NANOS |
| 13 | Cis-1,2-Dichloroethene | E | 228 | 3:59 | BV | 0.062 | NANOS |
| 14 | 1,1-Dichloroethane | E | 239 | 4:11 | BB | 0.976 | NANOS |
| 15 | Trans-1,2-Dichloroethane | E | 287 | 5:01 | VV | 2.022 | NANOS |
| 18 | 1,1,1-Trichloroethane | E | 366 | 6:24 | BB | 0.000 | NANOS |
| 19 | Benzene | E | 395 | 6:55 | BB | 0.383 | NANOS |
| 21 | 1,2-Dichloropropane | E | 449 | 7:51 | BB | 0.000 | NANOS |
| 24 | Trans-1,3-Dichloroprop | E | 565 | 9:53 | BB | 0.048 | NANOS |
| 25 | 1,1,2-Trichloroethane | E | 574 | 10:03 | BB | 0.000 | NANOS |
| 26 | Toluene | E | 599 | 10:29 | BB | 96.529 | NANOS |
| 28 | Tetrachloroethene | E | 687 | 12:01 | BB | 1.722 | NANOS |
| 29 | Chlorobenzene | E | 742 | 12:59 | BB | 0.102 | NANOS |
| 30 | Ethylbenzene | E | 775 | 13:34 | BB | 8.563 | NANOS |
| 31 | 1,3+1,4-Xylenes | E | 790 | 13:49 | VV | 15.779 | NANOS |
| 33 | 1,1,2,2-Tetrachloroeth | E | 837 | 14:39 | BV | 0.000 | NANOS |
| 34 | 1,2-Xylene | E | 831 | 14:33 | BB | 3.679 | NANOS |
| 35 | 4-Bromofluorobenzene | E | 872 | 15:16 | BV | 6.054 | NANOS |

Quantitation Report

Quanfile: VOC30-37

Quan Entries: 34

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB DUPLICATE

Sorted via: Entry Number ↑

(S) = Standard

| Cal | Name of Compound | S | Scan# | R Time | Me | Calc Amt(A) | Units |
|-----|------------------------|---|-------|--------|----|-------------|-------|
| 36 | Benzylchloride | E | 935 | 16:22 | VB | 0.000 | NANOS |
| 37 | 4-Ethyltoluene | E | 947 | 16:34 | BV | 0.822 | NANOS |
| 38 | 1,3,5-Trimethylbenzene | E | 959 | 16:47 | VB | 0.000 | NANOS |
| 39 | 1,2,4-Trimethylbenzene | E | 1001 | 17:31 | VB | 0.000 | NANOS |
| 40 | 1,3-Dichlorobenzene | E | 1023 | 17:54 | BB | 0.000 | NANOS |
| 41 | 1,4-Dichlorobenzene | E | 1023 | 17:54 | BB | 0.000 | NANOS |
| 43 | 1,2,4-Trichlorobenzene | E | 1276 | 22:20 | BB | 0.000 | NANOS |
| 44 | Naphthalene | E | 1289 | 22:33 | BB | 0.000 | NANOS |

SECTION 5.0
GC/ MS ION CHROMATOGRAMS FOR THE ANALYSIS OF TO-14 VOLATILE
ORGANIC COMPOUNDS IN GAS SAMPLES

Chromatogram Plot

File: D:\VOC30-29

Date: Aug-15-2000 14:00:29

Comment: T0-14 VOC STANDARD 1.5 NGS0 + 4-BFB

Scan No: 1

Retention Time: 0:01

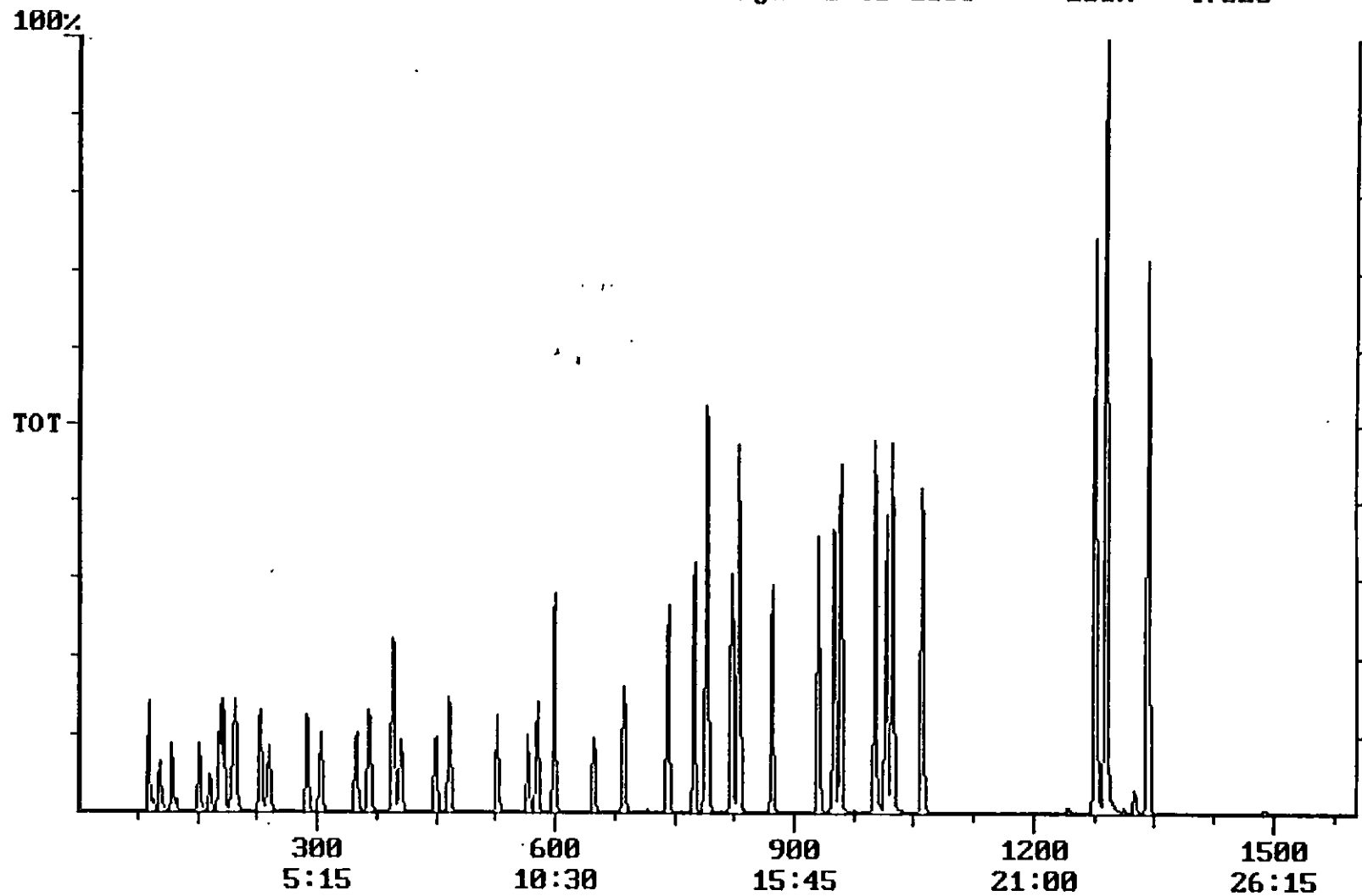
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1599

Range: 1 to 1599

100% = 47826



Chromatogram Plot

File: D:\VOC30-30

Date: Aug-15-2000 14:38:31

Comment: T0-14 VOC STANDARD 3.0 NGS0 + 4-BFB

Scan No: 1

Retention Time: 0:01

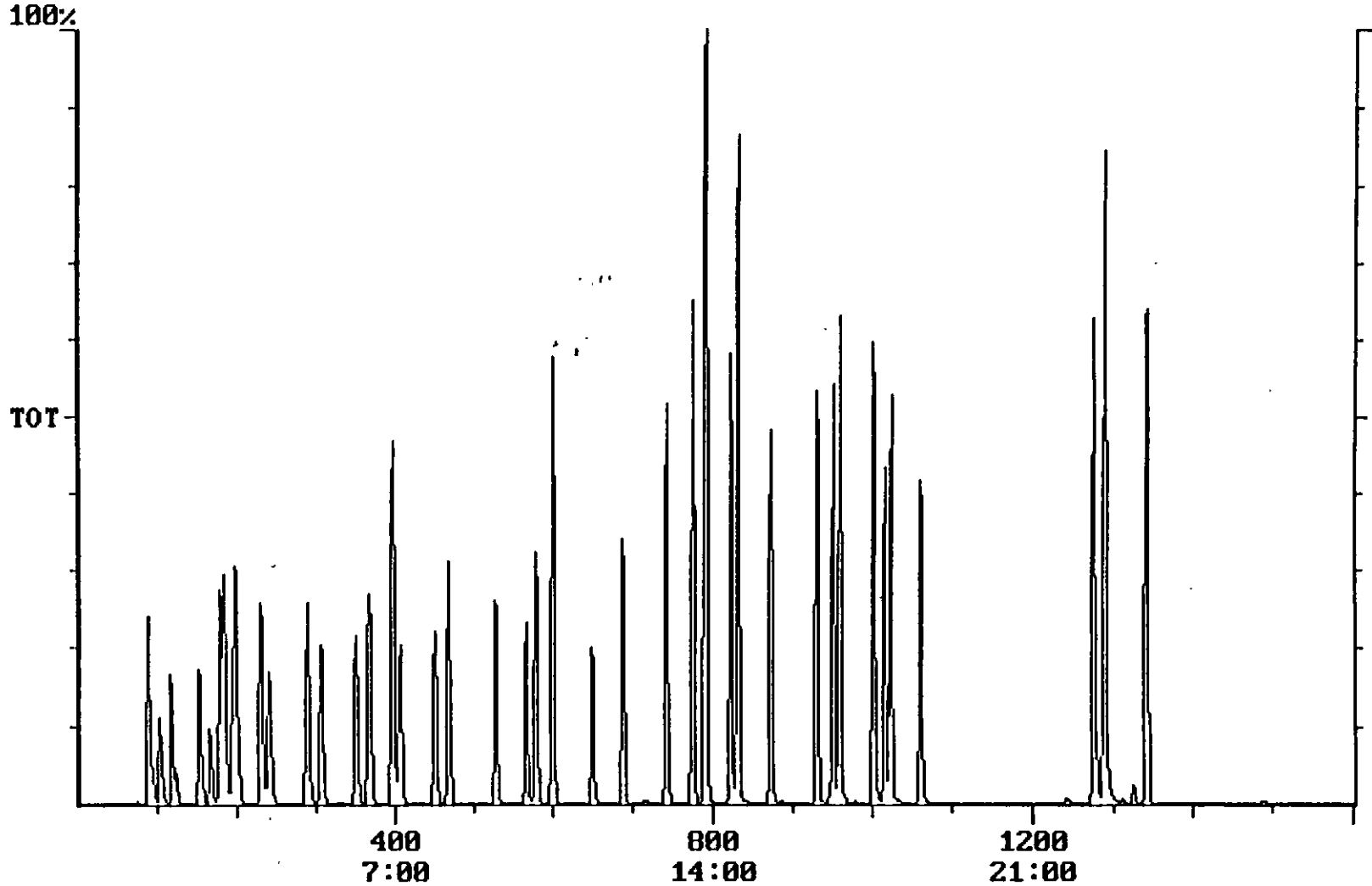
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1600

Range: 1 to 1600

100% = 50352



Chromatogram Plot

File: D:\VOC30-31

Date: Aug-15-2000 15:14:29

Comment: T0-14 VOC STANDARD 4.5 NGSE + 4-BFB

Scan No: 1

Retention Time: 0:01

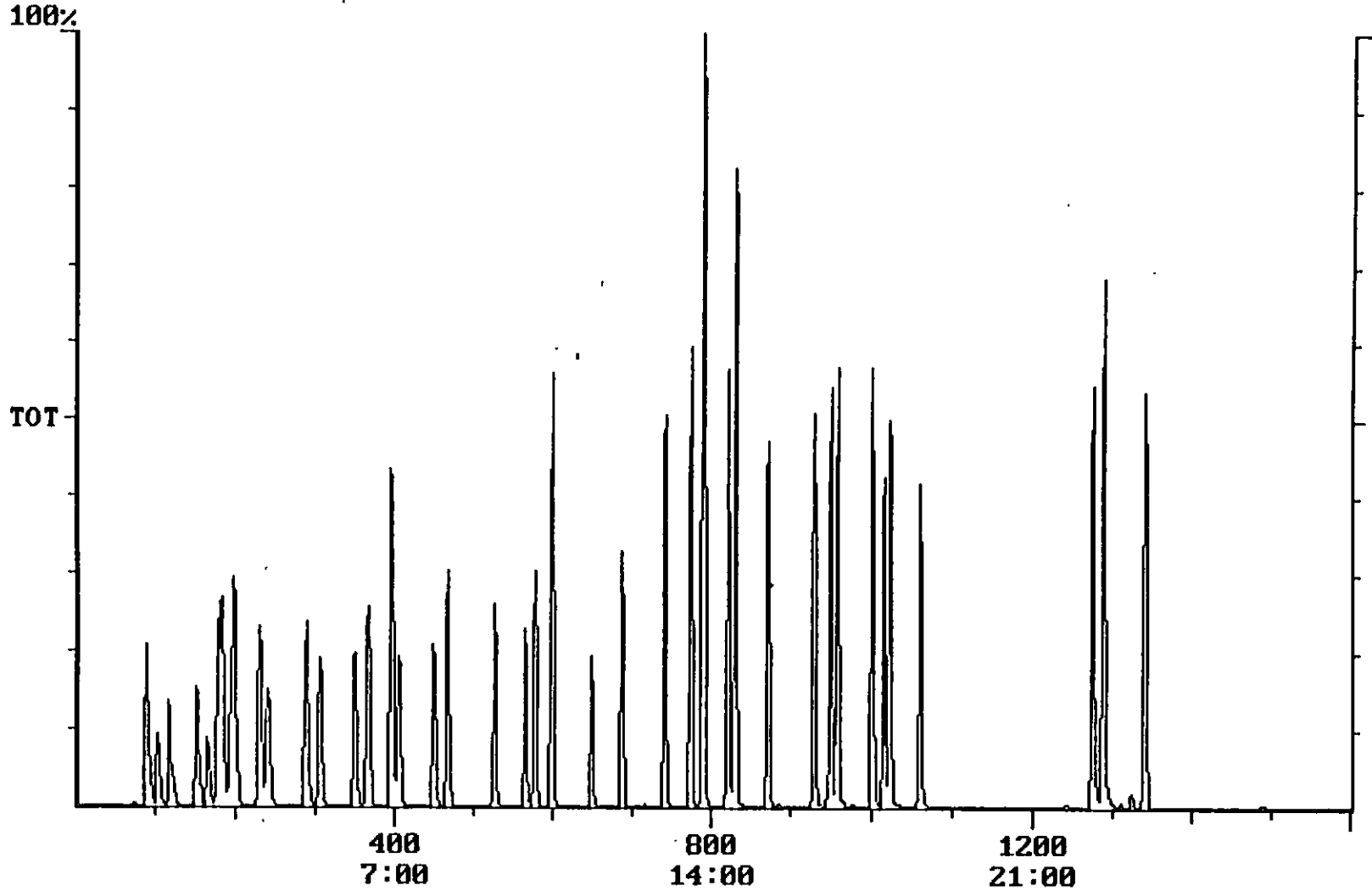
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1600

Range: 1 to 1600

100% = 65568



Chromatogram Plot

File: D:\VOC30-32

Date: Aug-15-2000 15:48:04

Comment: T0-14 VOC STANDARD 6.0 NGSO + 4-BFB

Scan No: 1

Retention Time: 0:01

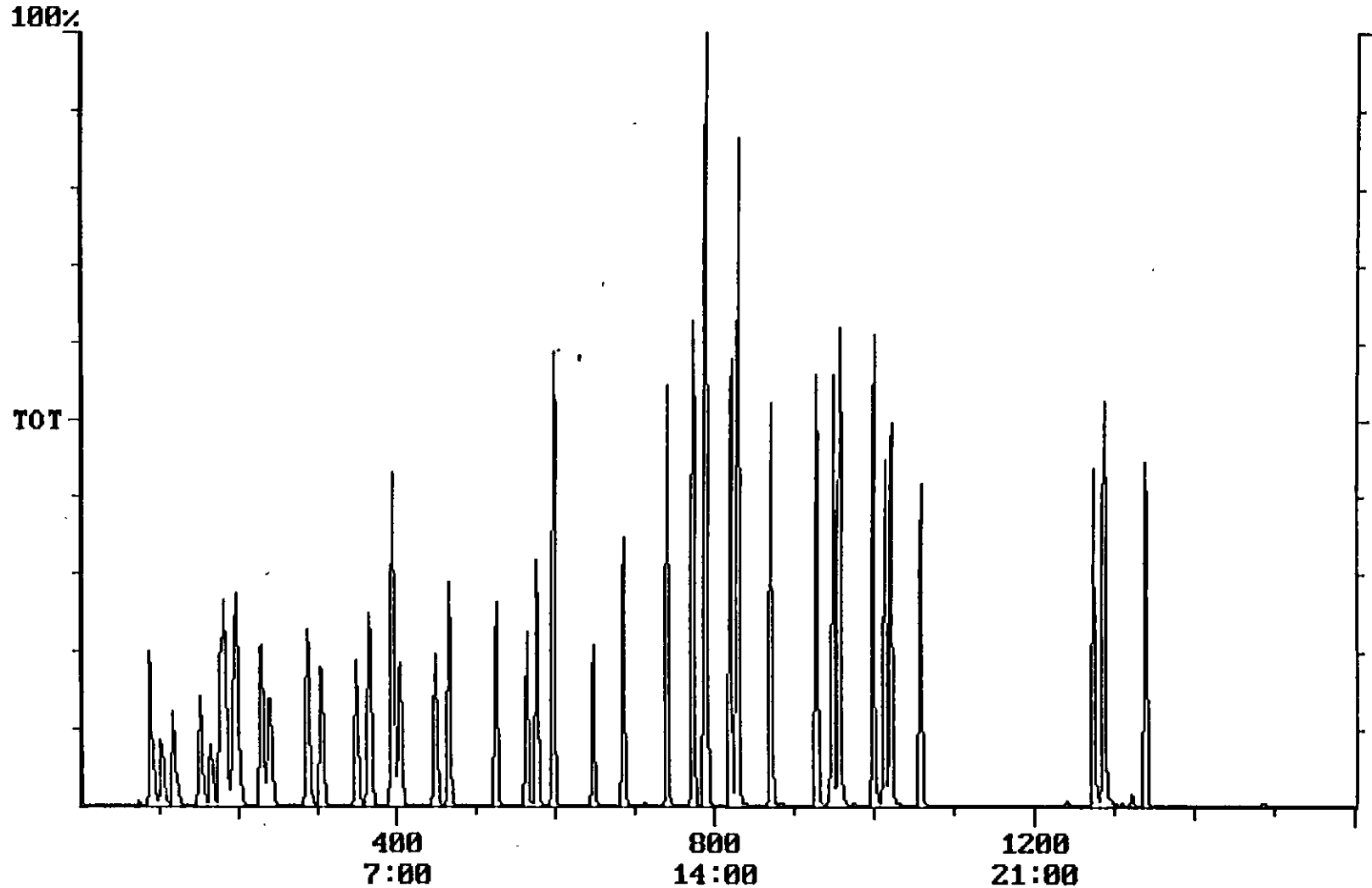
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1600

Range: 1 to 1600

100% = 84818



Chromatogram Plot

File: D:\UOC30-33

Date: Aug-15-2000 16:19:55

Comment: T0-14 VOC STANDARD 7.5 NGSQ + 4-BFB

Scan No: 1

Retention Time: 0:01

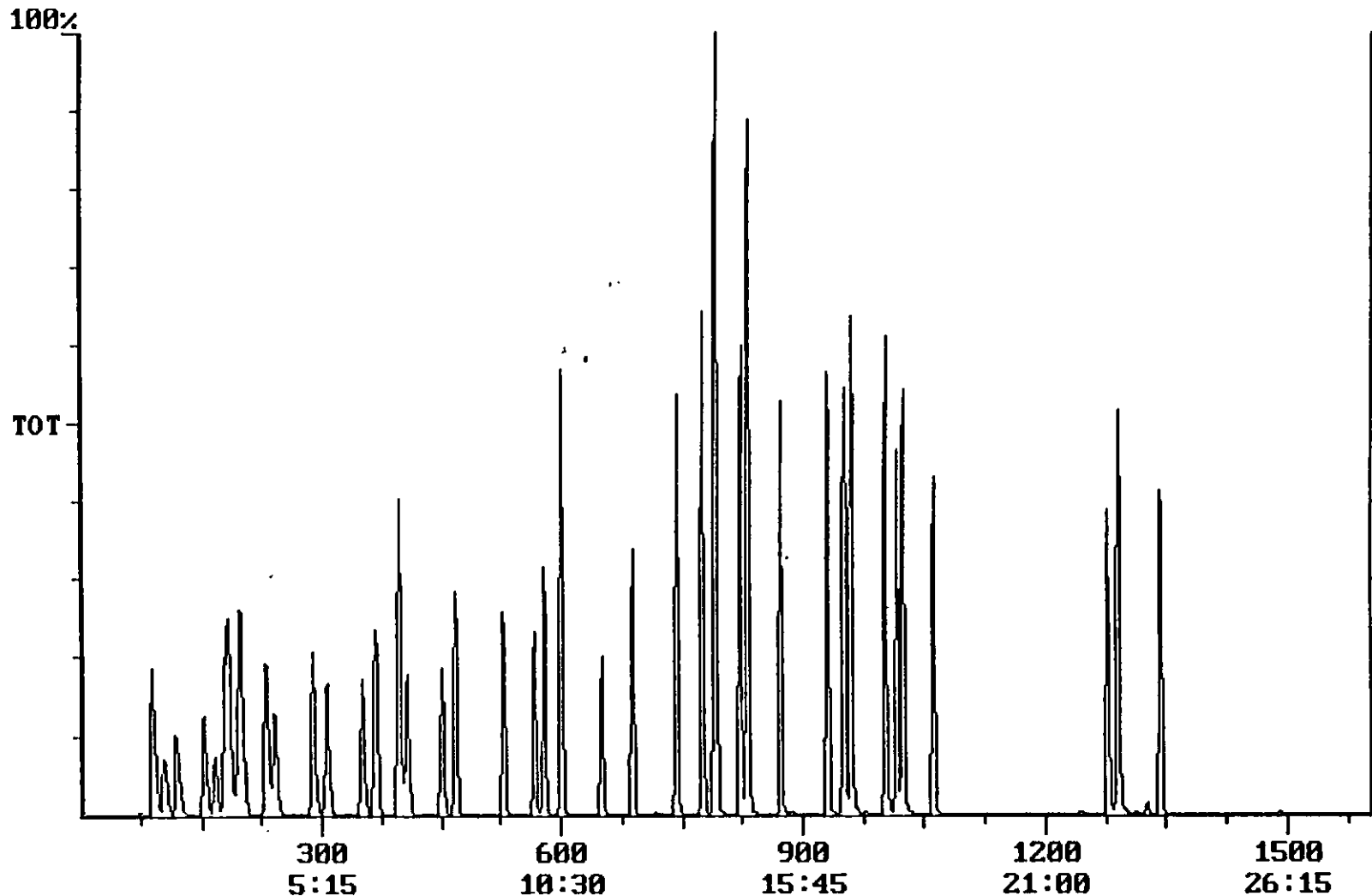
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1599

Range: 1 to 1599

100% = 99287



Chromatogram Plot

File: D:\NUOC30-34

Date: Aug-15-2000 16:55:33

Comment: TO-14 VOC STANDARD 9.0 NGS0 + 4-BFB

Scan No: 1

Retention Time: 0:01

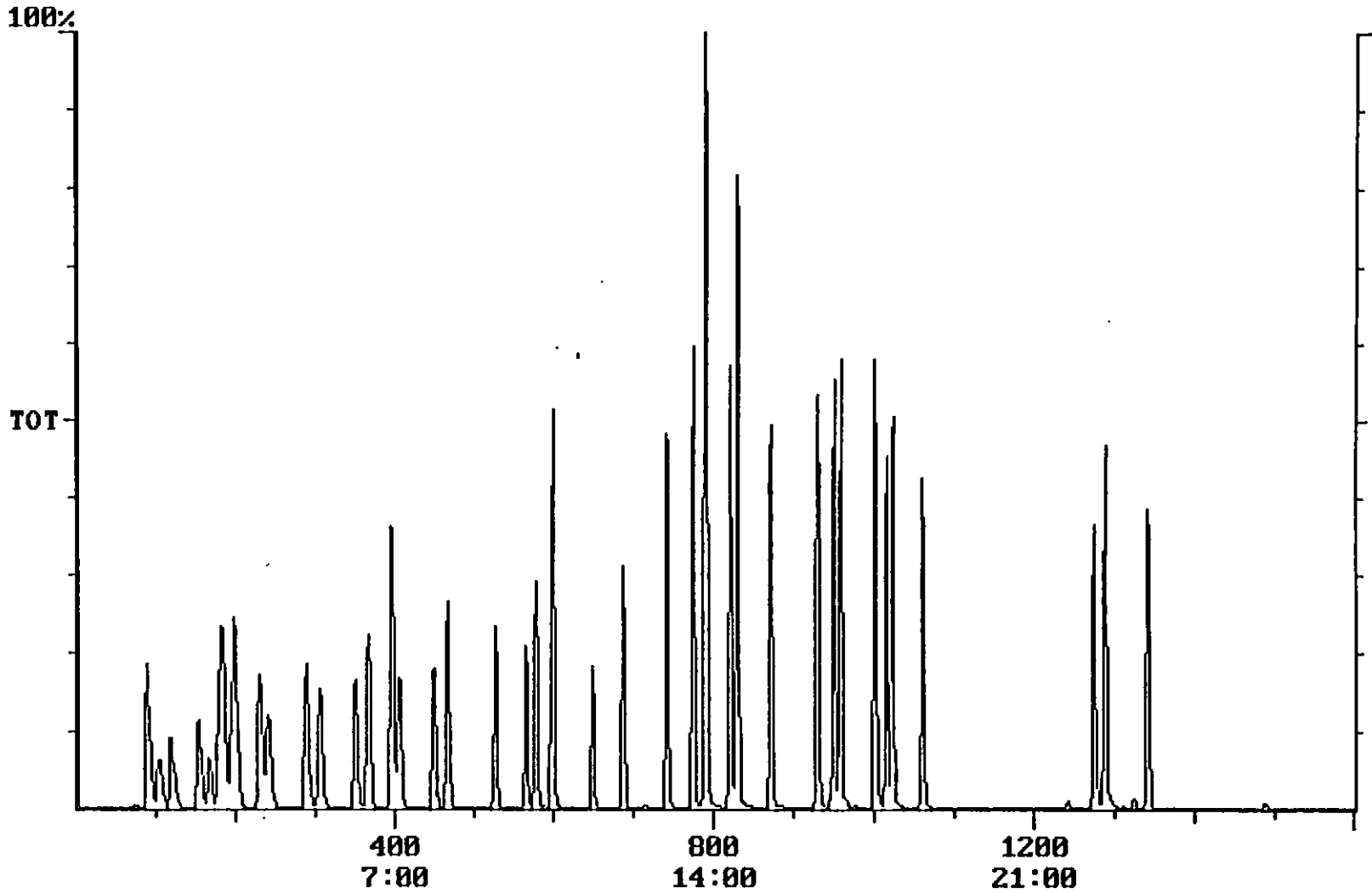
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1600

Range: 1 to 1600

100% = 121016



Chromatogram Plot

File: D:\VOC30-35

Date: Aug-15-2000 17:35:56

Comment: LAB AIR BLANK 0.50 CC + 4-BFB

Scan No: 1

Retention Time: 0:01

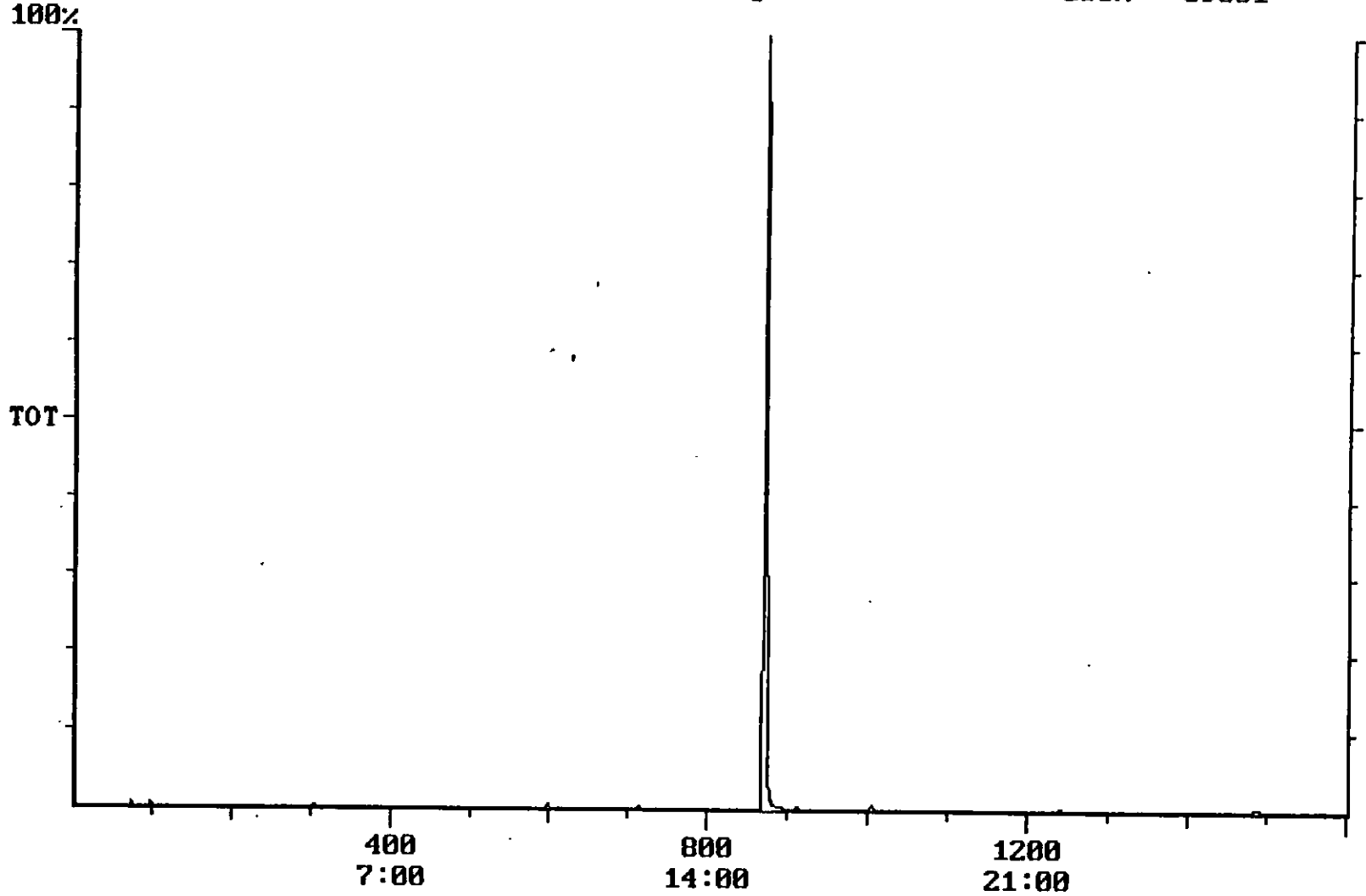
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1600

Range: 1 to 1600

100% = 39661



Chromatogram Plot

File: D:\U0030-36

Date: Aug-15-2000 18:11:27

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB

Scan No: 1

Retention Time: 0:01

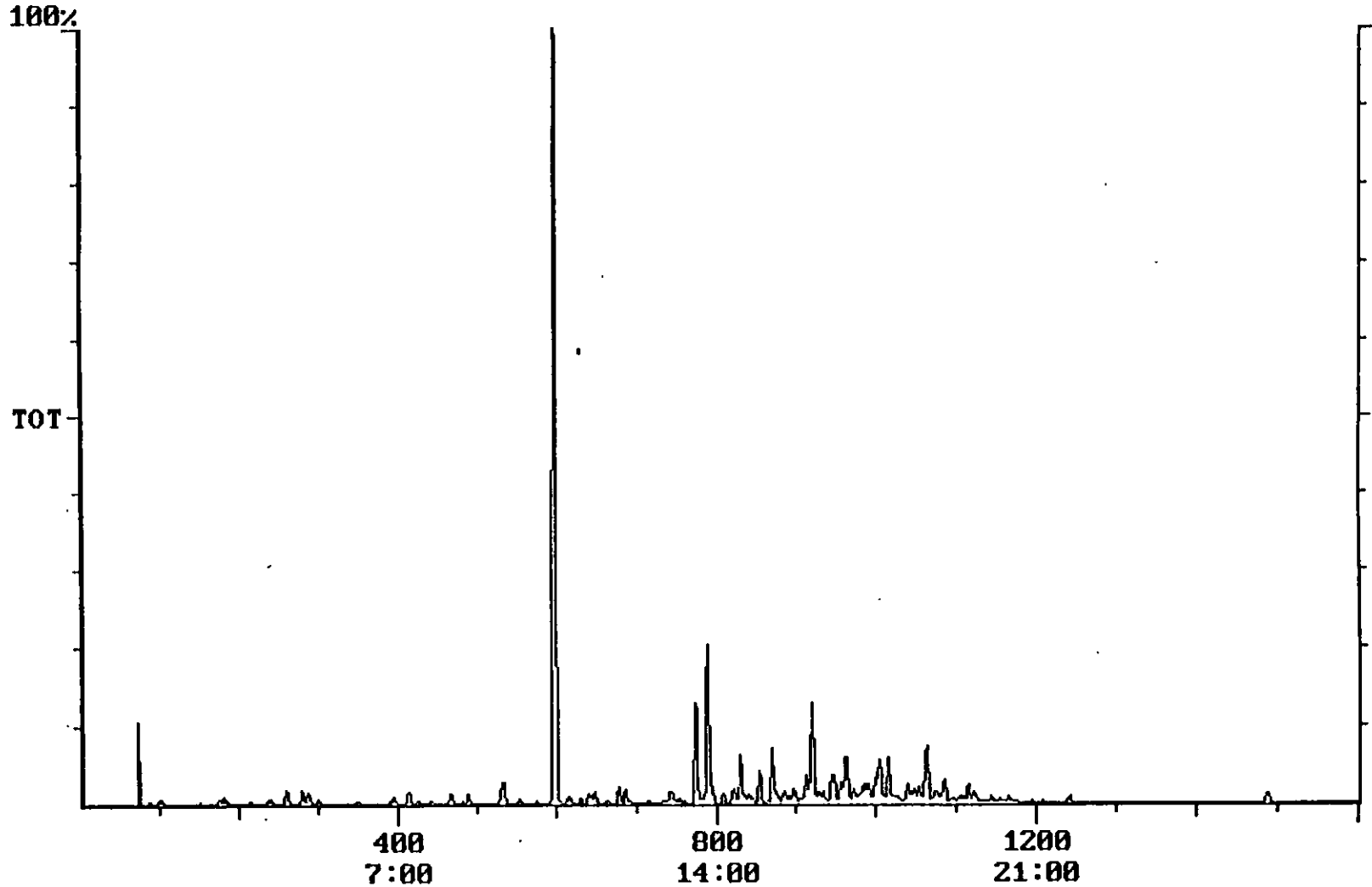
RIC: 0

Mass Range: 0 - 0

Plotted: 1 to 1600

Range: 1 to 1600

100% = 552634



Spectrum Plot

File: D:\VOC30-36

Date: Aug-15-2000 18:11:27

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB

Scan No: 599

Retention Time: 10:29

RIC: 546791

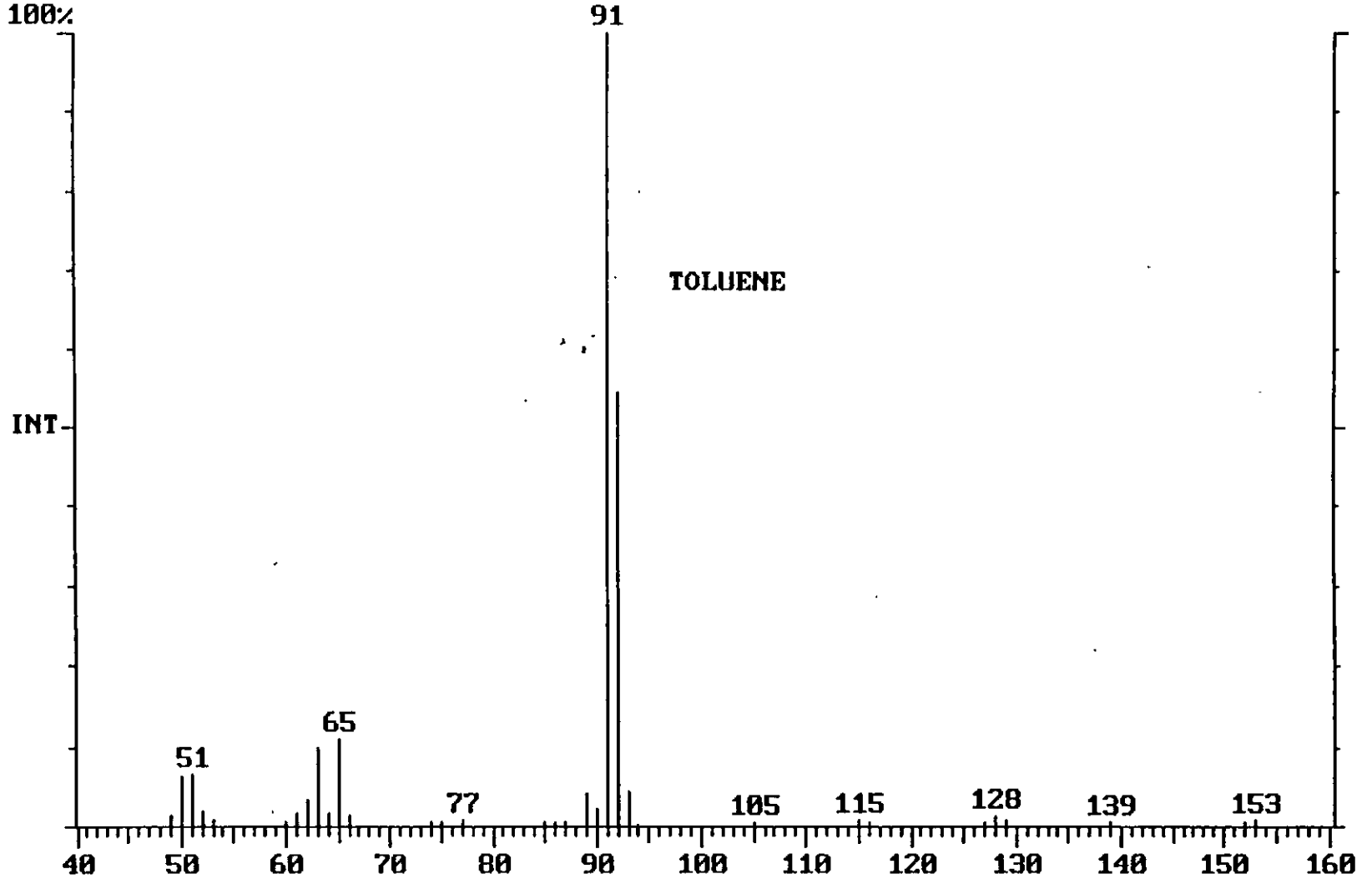
Mass Range: 49 - 153

Pks:33

Base Pk: 91

Int: 250420

100.00% = 250420



Spectrum Plot

File: D:\UOC30-36

Date: Aug-15-2000 18:11:27

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB

Scan No: 774

Retention Time: 13:33

RIC: 69881

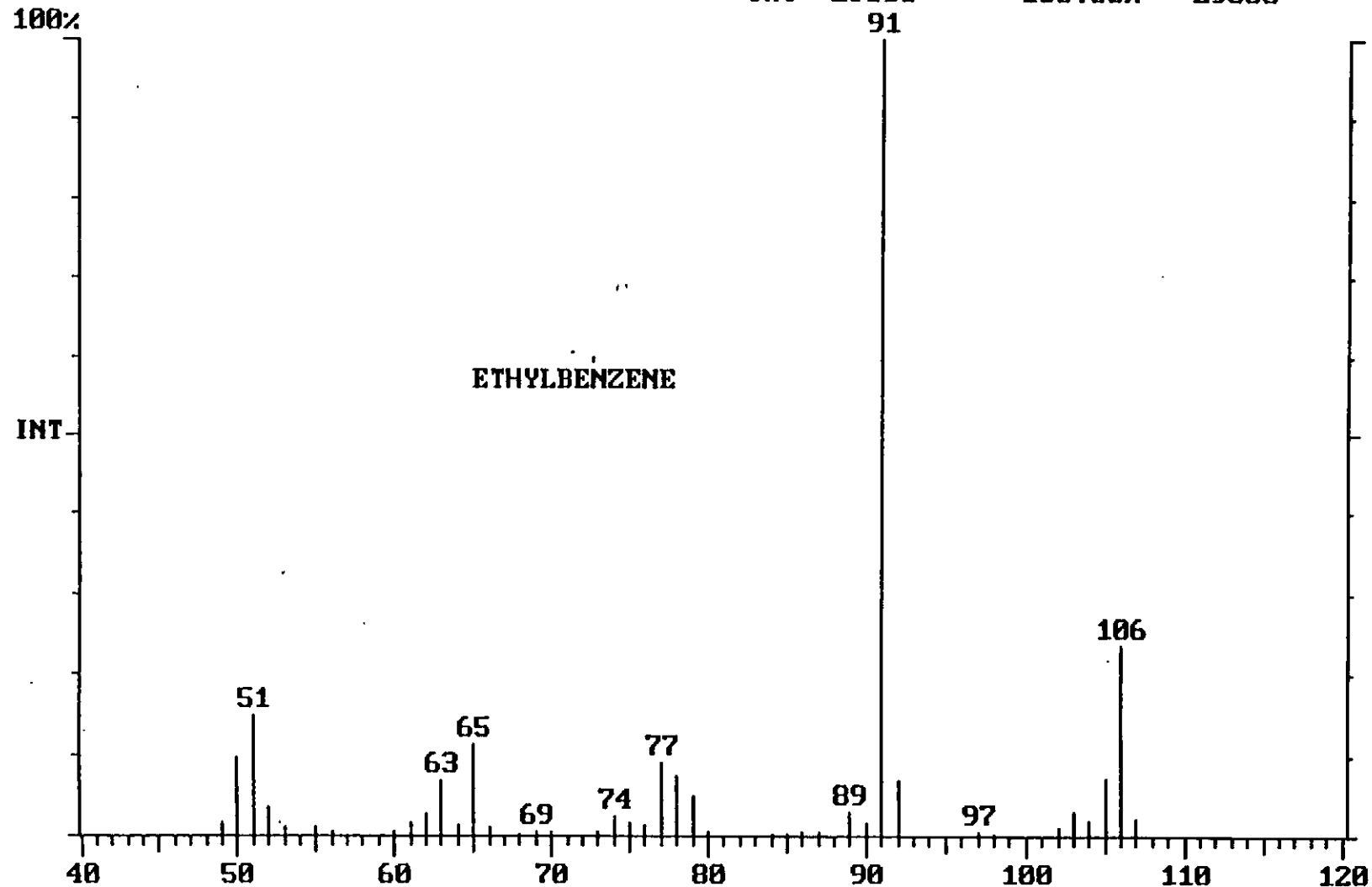
Mass Range: 49 - 107

Pks:41

Base Pk: 91

Int: 29565

100.00% = 29565



Spectrum Plot

File: D:\UOC30-36

Date: Aug-15-2000 18:11:27

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB

Scan No: 789

Retention Time: 13:48

RIC: 111895

Mass Range: 49 - 119

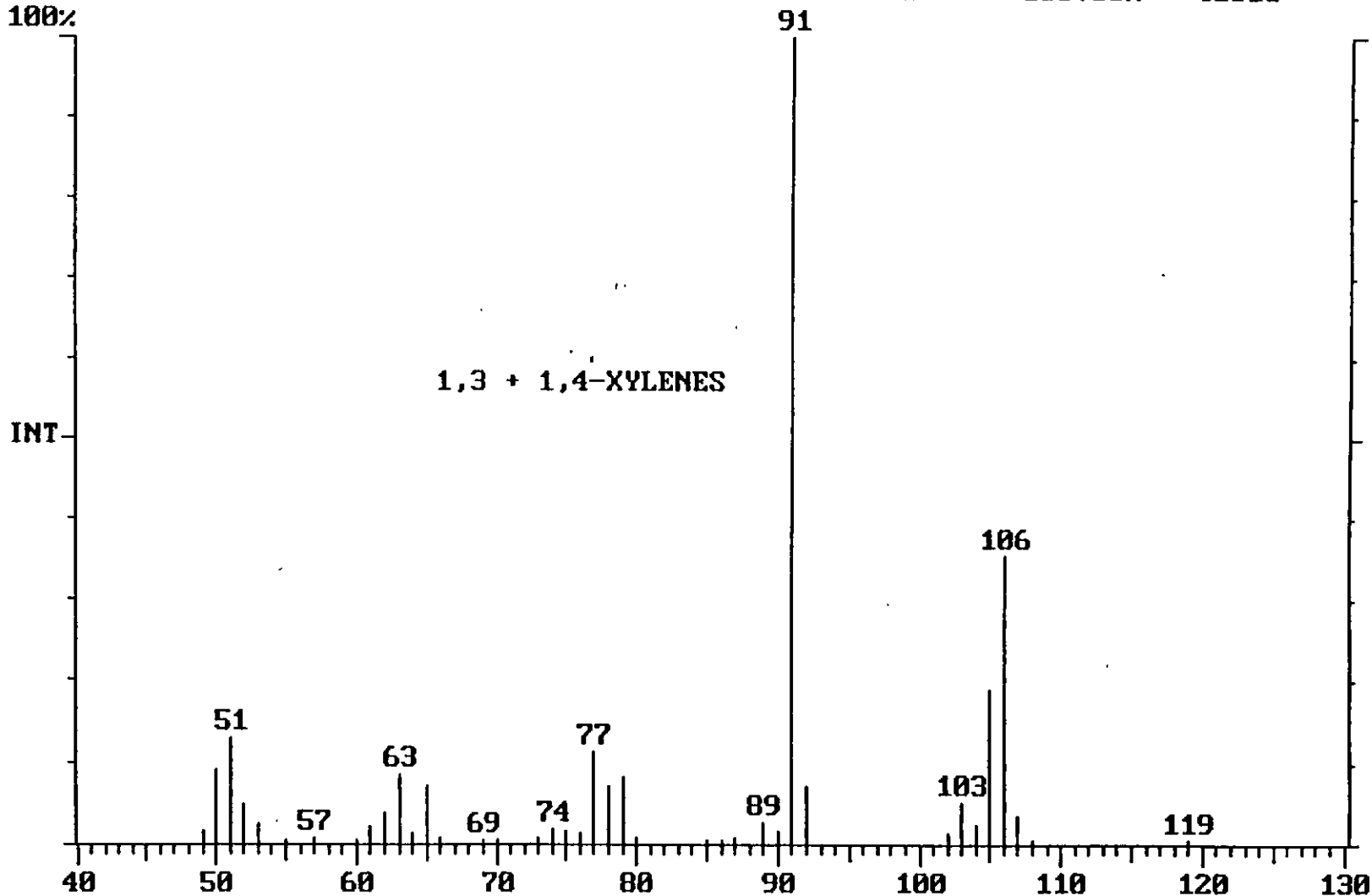
Pks:39

Base Pk: 91

Int: 41813

100.00% = 41813

100%



Spectrum Plot

File: D:\VOC30-36

Date: Aug-15-2000 18:11:27

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB

Scan No: 831

Retention Time: 14:33

RIC: 34360

Mass Range: 49 - 119

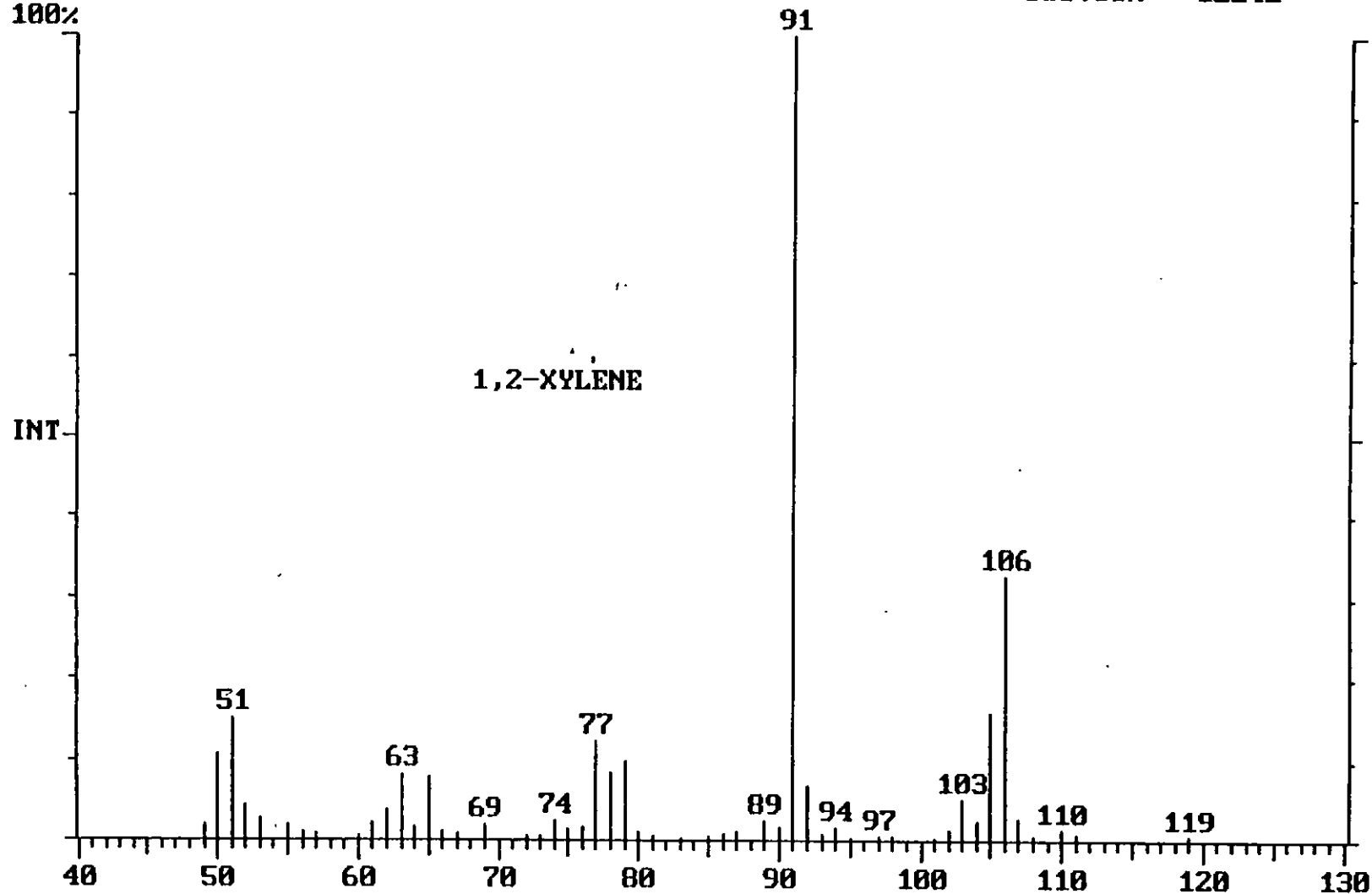
Pks: 51

Base Pk: 91

Int: 12242

100.00% = 12242

100%



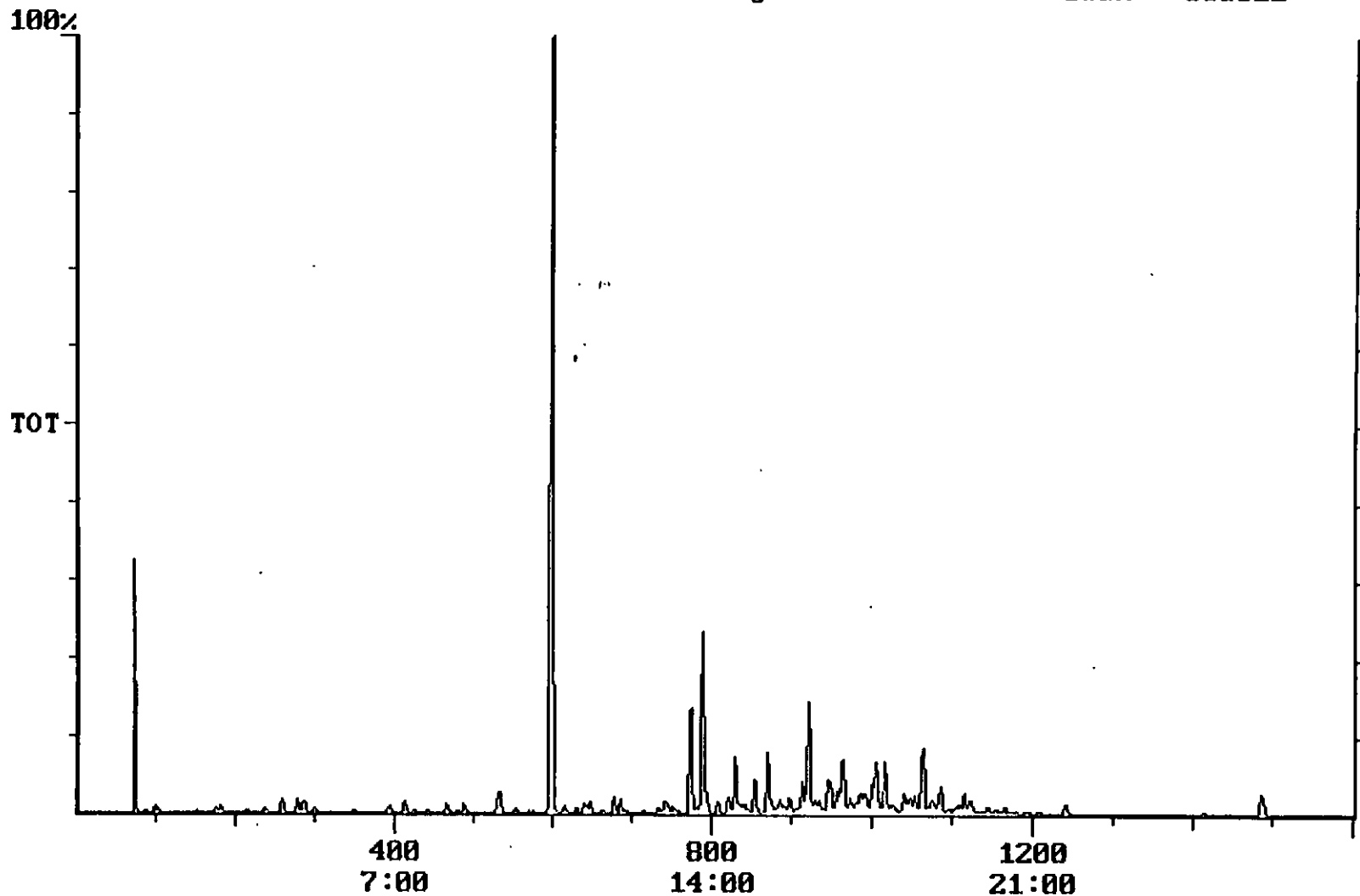
Chromatogram Plot

File: D:\VOC30-37 Date: Aug-15-2000 18:47:53

Comment: LFR GB 8/14/2000 DM LF 0.10 CC + 4-BFB DUPLICATE

Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0

Plotted: 1 to 1600 Range: 1 to 1600 100% = 508322



SECTION 6.0
4-BROMOFLUOROBENZENE TUNING CHARACTERISTICS AND RUN LOGS FOR
THE ANALYSIS OF TO-14 VOLATILE ORGANIC COMPOUNDS IN GAS SAMPLES

Spectrum Plot

Comment: T0-14 VOC STANDARD

Scan No: 872

Pks: 51

3.0 NGSQ + 4-BFB

Retention Time: 15:16

Base Pk: 176

File: D:\VOC30-30

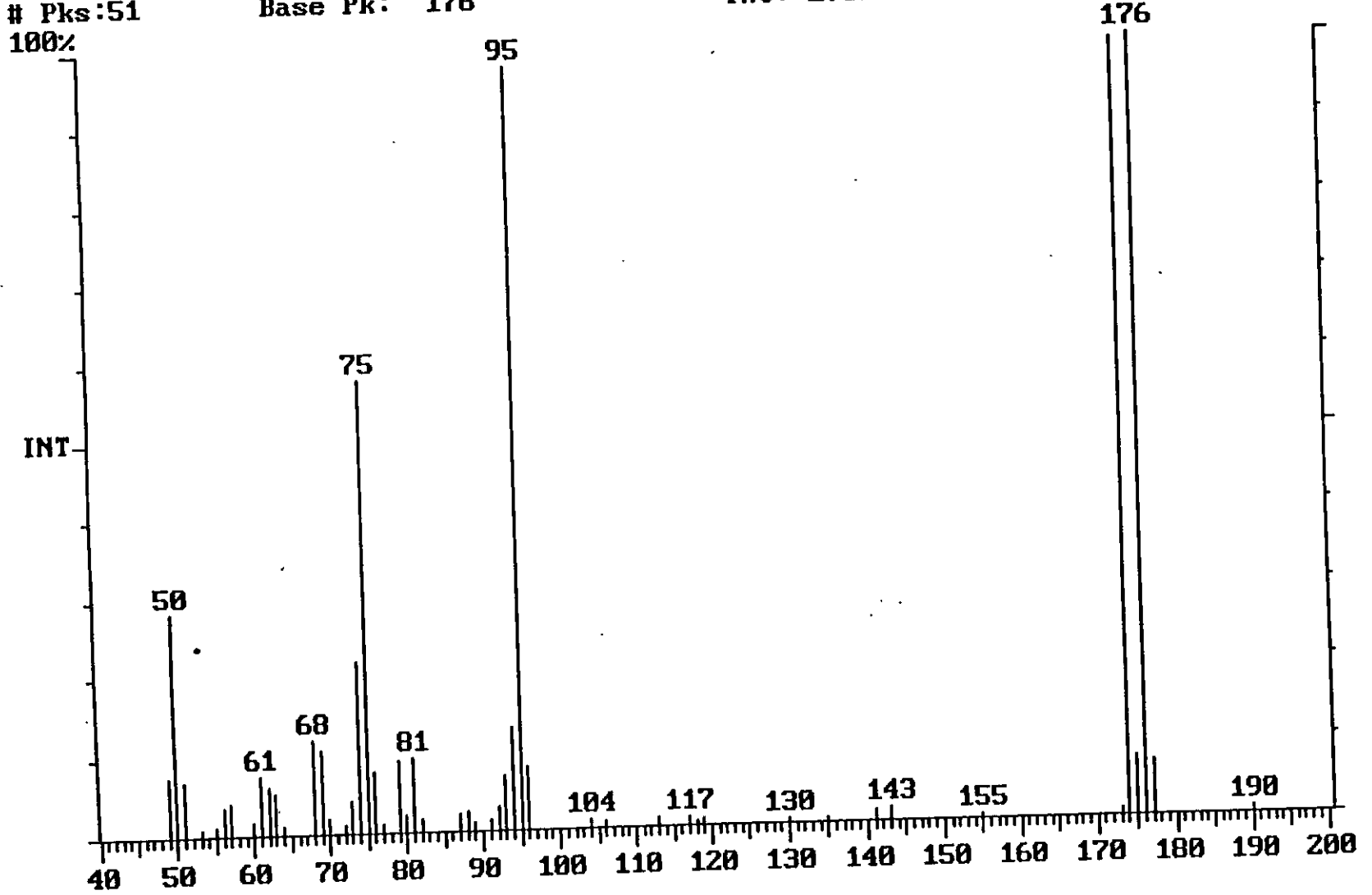
Date: Aug-15-2000 14:38:31

RIC: 15813

Int: 2782

Mass Range: 49 - 190

100.00% = 2782



Mass List

Comment: T0-14 VOC STANDARD
 Scan No: 872
 # Peaks: 51

3.0 NGSC + 4-BFB
 Retention Time: 15:16
 Mass of Base Peak: 176

File: D:\UOC30-30

Date: Aug-15-2000 14:38:31

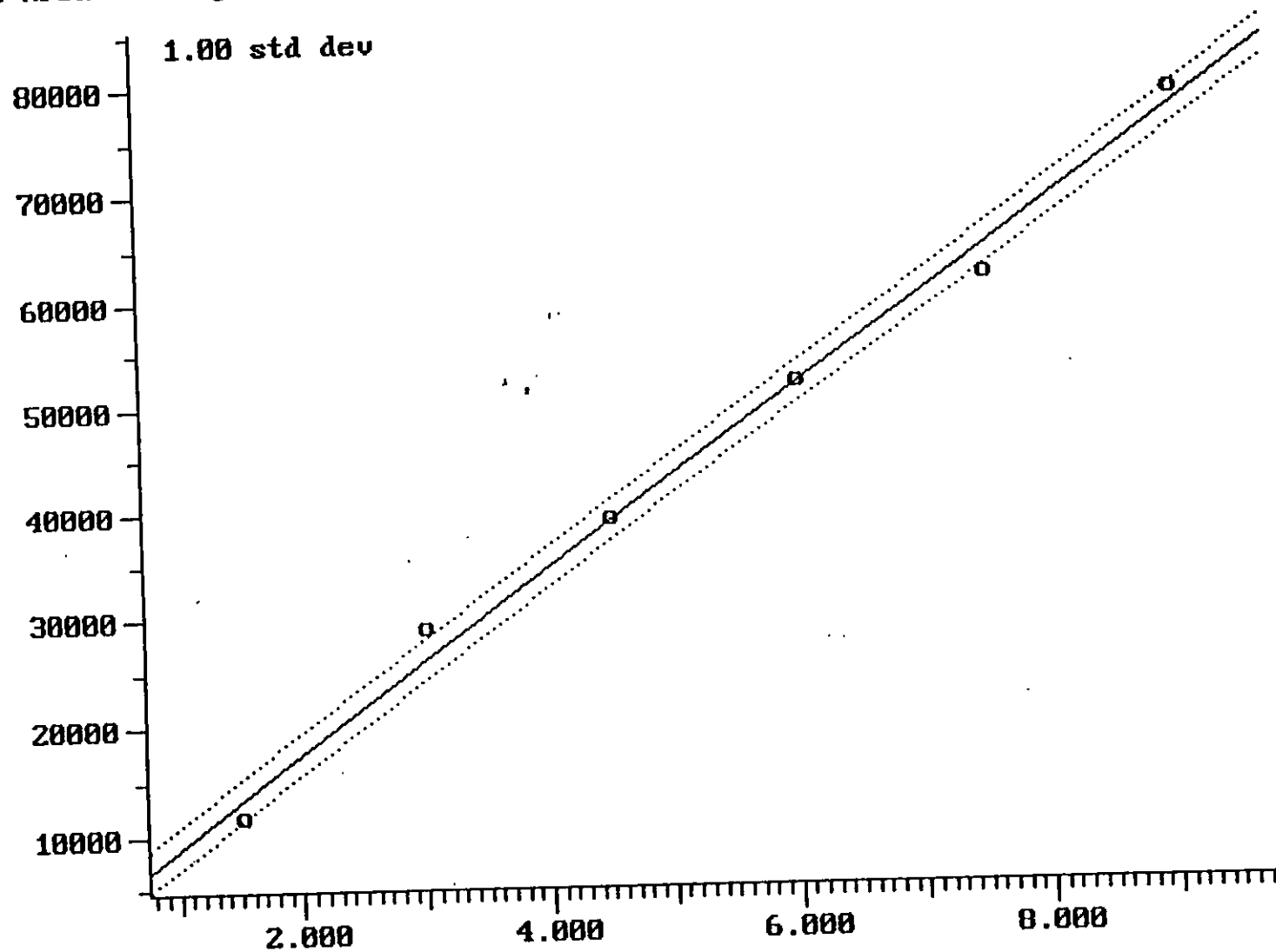
RIC: 15813
 Int: 2782

Mass Range: 49 - 190
 Threshold: 1.00%

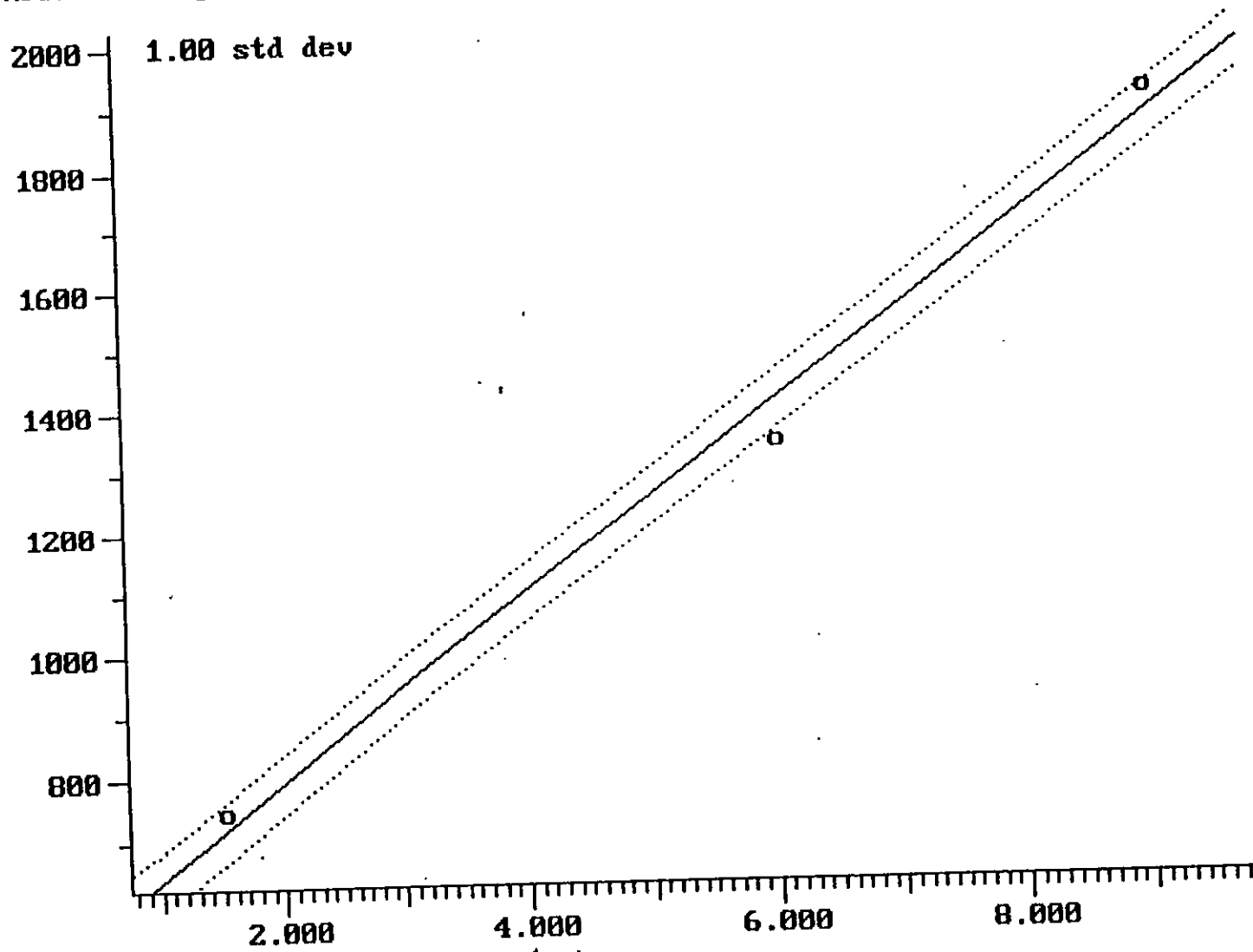
| Mass | Intensity | % Base | Mass | Intensity | % Base | Mass | Intensity | % Base |
|------|-----------|--------|------|-----------|--------|------|-----------|--------|
| 49 | 200 | 7.19 | 94 | 358 | 12.87 | | | |
| 50 | 782 | 28.11 | 95 | 2,716 | 97.63 | | | |
| 51 | 190 | 6.83 | 96 | 217 | 7.80 | | | |
| 56 | 95 | 3.41 | 104 | 30 | 1.08 | | | |
| 57 | 110 | 3.95 | 141 | 35 | 1.26 | | | |
| 60 | 42 | 1.51 | 143 | 45 | 1.62 | | | |
| 61 | 203 | 7.30 | 174 | 2,769 | 99.53 | | | |
| 62 | 167 | 6.00 | 175 | 200 | 7.19 | | | |
| 63 | 143 | 5.14 | 176 | 2,782 | 100.00 | | | |
| 68 | 328 | 11.79 | 177 | 185 | 6.65 | | | |
| 69 | 287 | 10.32 | | | | | | |
| 70 | 52 | 1.87 | | | | | | |
| 73 | 113 | 4.06 | | | | | | |
| 74 | 603 | 21.68 | | | | | | |
| 75 | 1,602 | 57.58 | | | | | | |
| 76 | 212 | 7.62 | | | | | | |
| 79 | 252 | 9.06 | | | | | | |
| 80 | 58 | 2.08 | | | | | | |
| 81 | 255 | 9.17 | | | | | | |
| 82 | 45 | 1.62 | | | | | | |
| 87 | 57 | 2.05 | | | | | | |
| 88 | 68 | 2.44 | | | | | | |
| 91 | 32 | 1.15 | | | | | | |
| 92 | 83 | 2.98 | | | | | | |
| 93 | 187 | 6.72 | | | | | | |

SECTION 7.0
INITIAL CALIBRATION CHARTS AND TABLES FOR THE ANALYSIS OF TO-14
VOLATILE ORGANIC COMPOUNDS IN GAS SAMPLES

Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.996
Freon 12 Compound: 1 of 45 Standard Deviation: 1.875
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin:Lin)

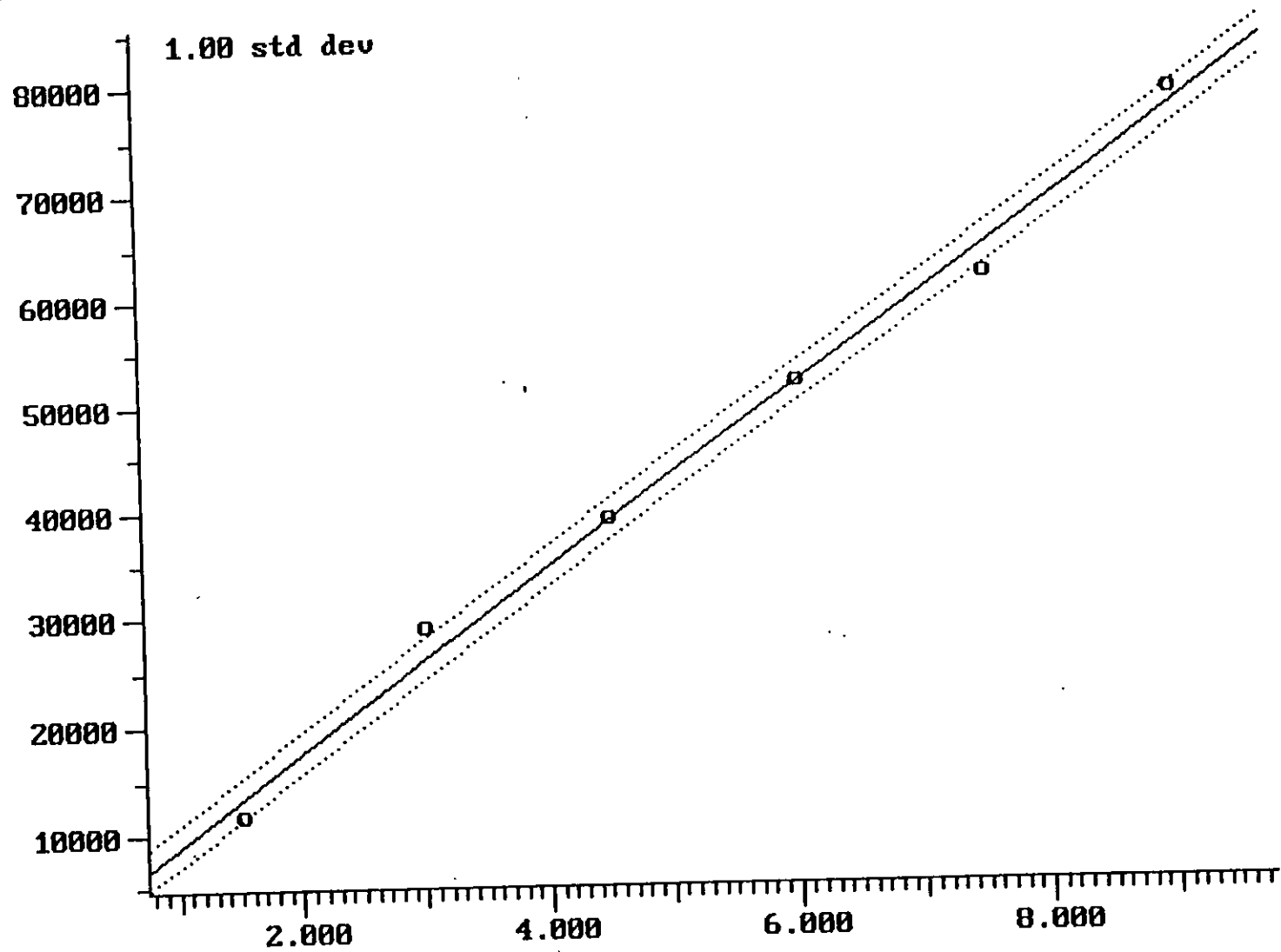


Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.995
Methyl Chloride Compound: 2 of 45 Standard Deviation: 0.047
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)

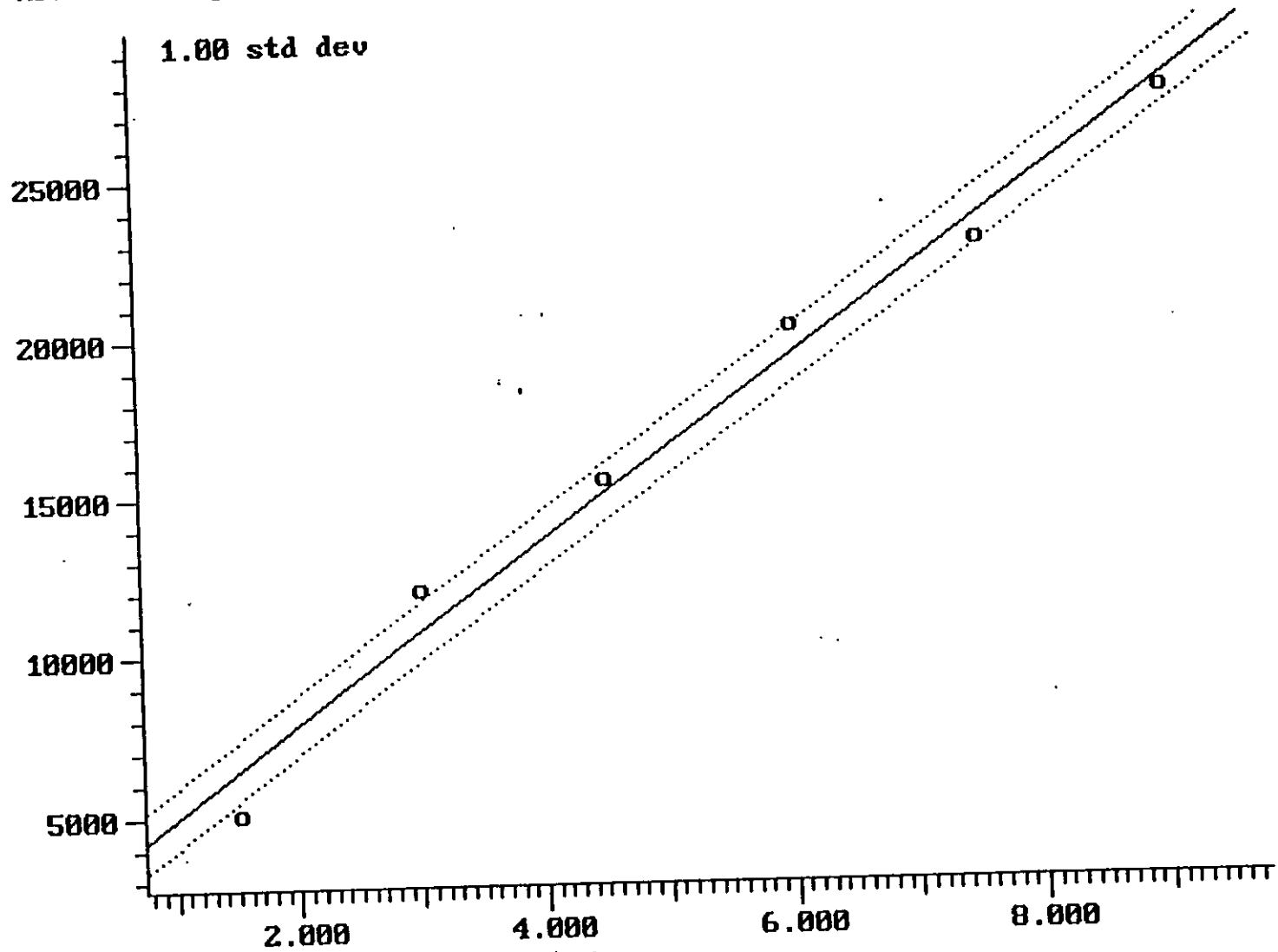


Calibration Plot (Ext Stds) Filename: T014GB
Freon 114 Compound: 3 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.996
Standard Deviation: 1.875
(Lin/Lin)



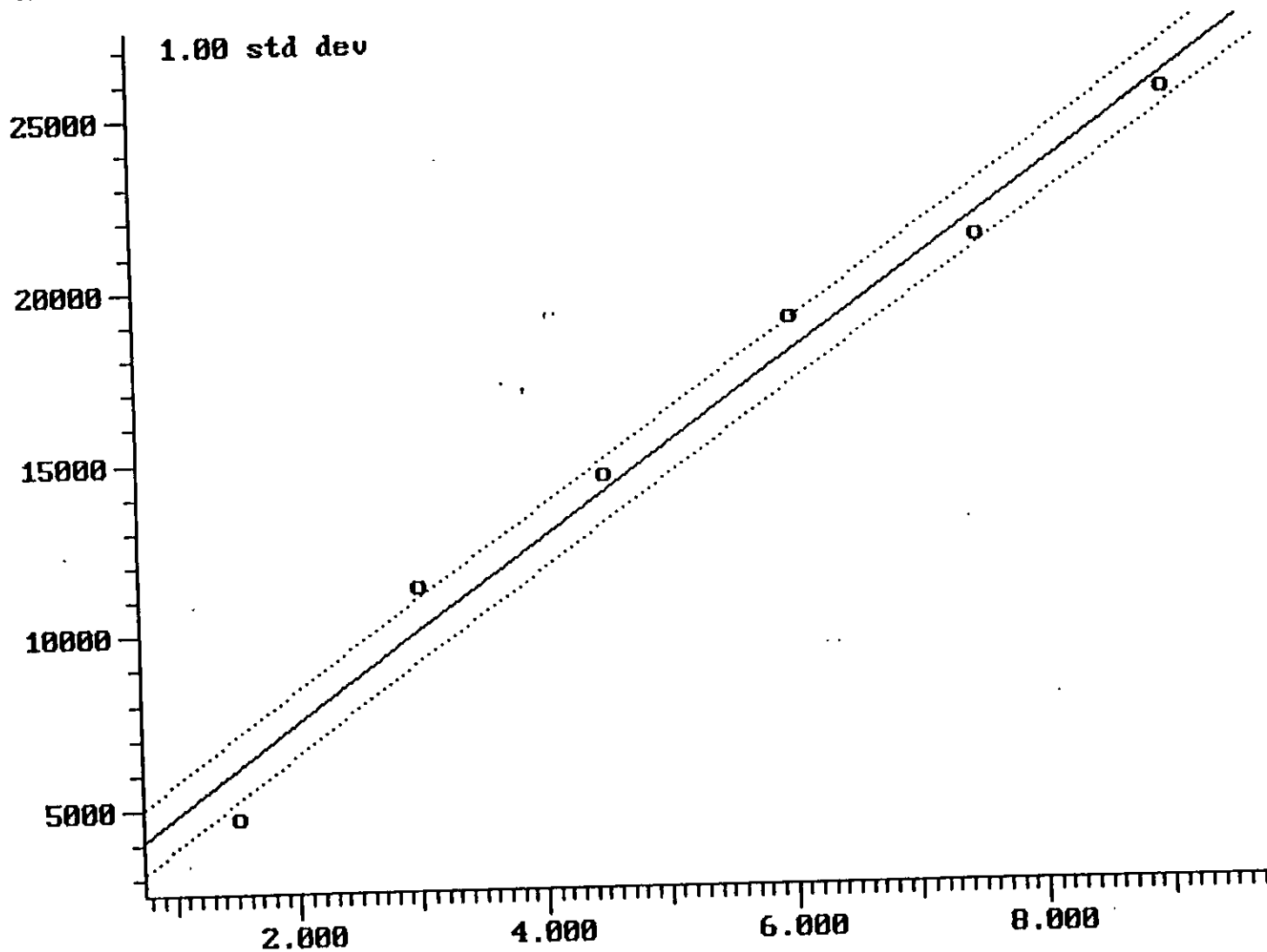
Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.992
Vinyl Chloride Compound: 4 of 45 Standard Deviation: 0.927
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)



Calibration Plot (Ext Stds)
Methyl Bromide
(Peak Area of Sample) vs (Amount of Sample Injected)

Filename: T014GB
Compound: 5 of 45

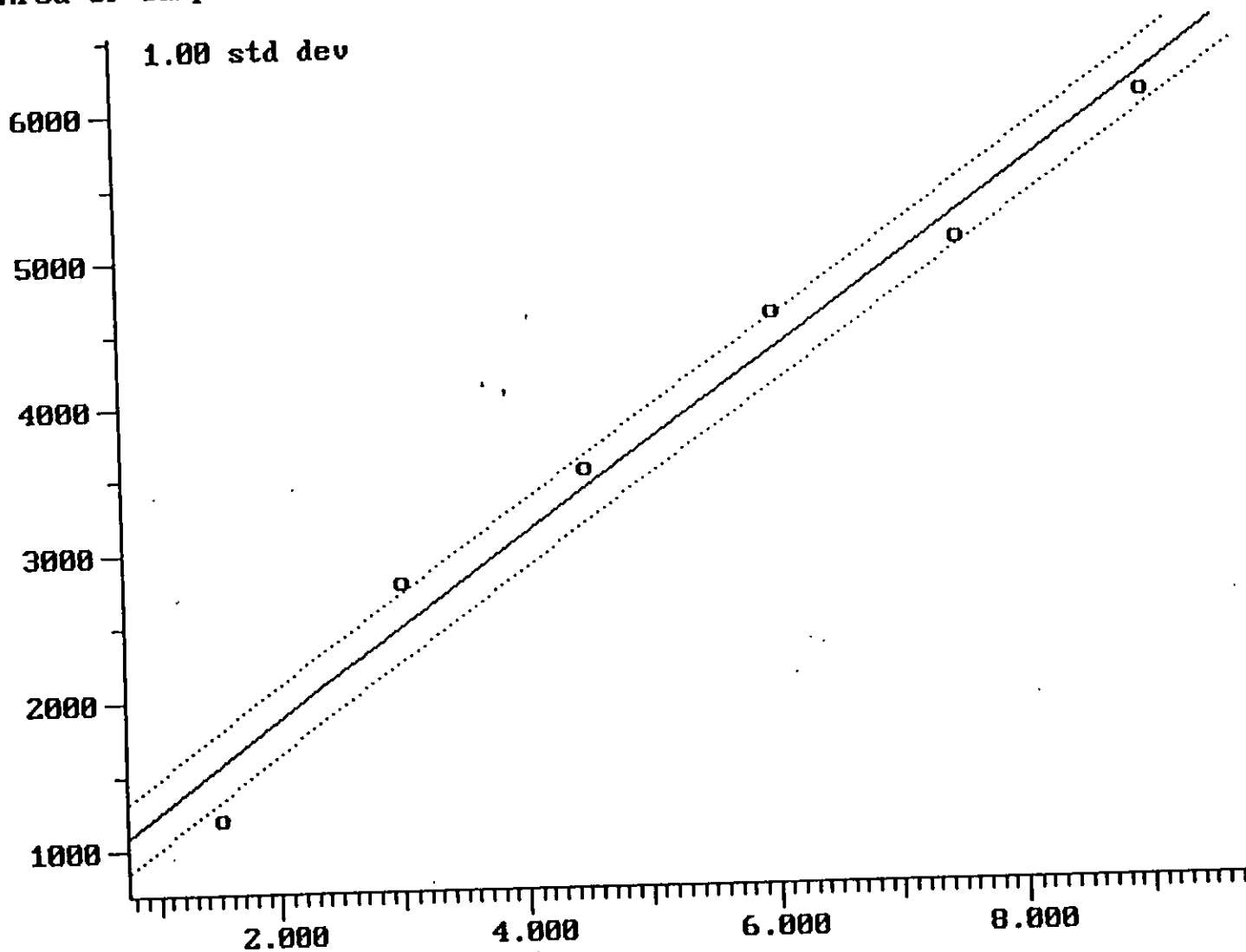
Correlation Coeff: 0.998
Standard Deviation: 0.943
(LinLin)



Calibration Plot (Ext Stds)
Ethyl Chloride
(Peak Area of Sample) vs (Amount of Sample Injected)

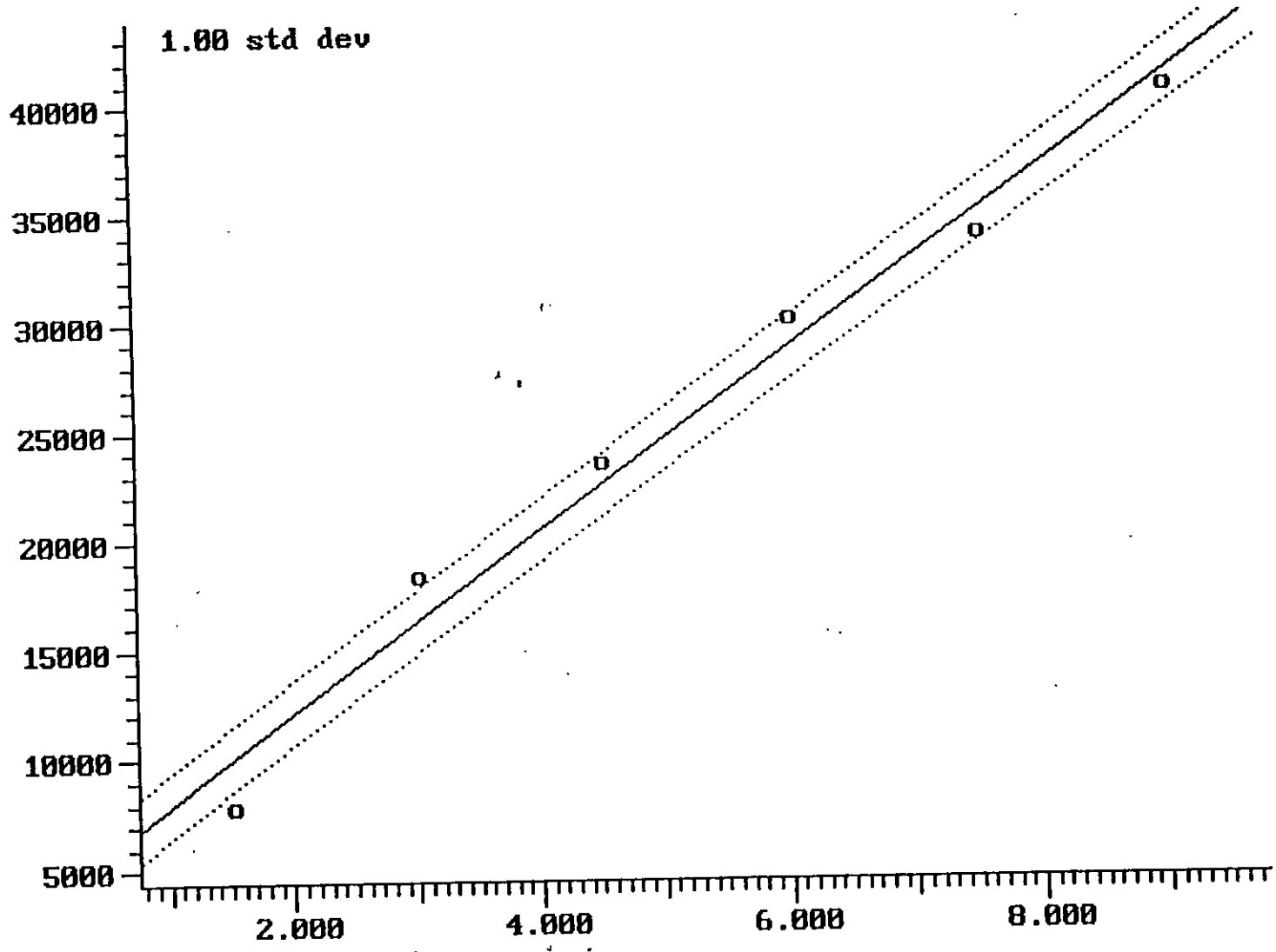
Filename: T014GB
Compound: 6 of 45

Correlation Coeff: 0.989
Standard Deviation: 0.237
(Lin/Lin)



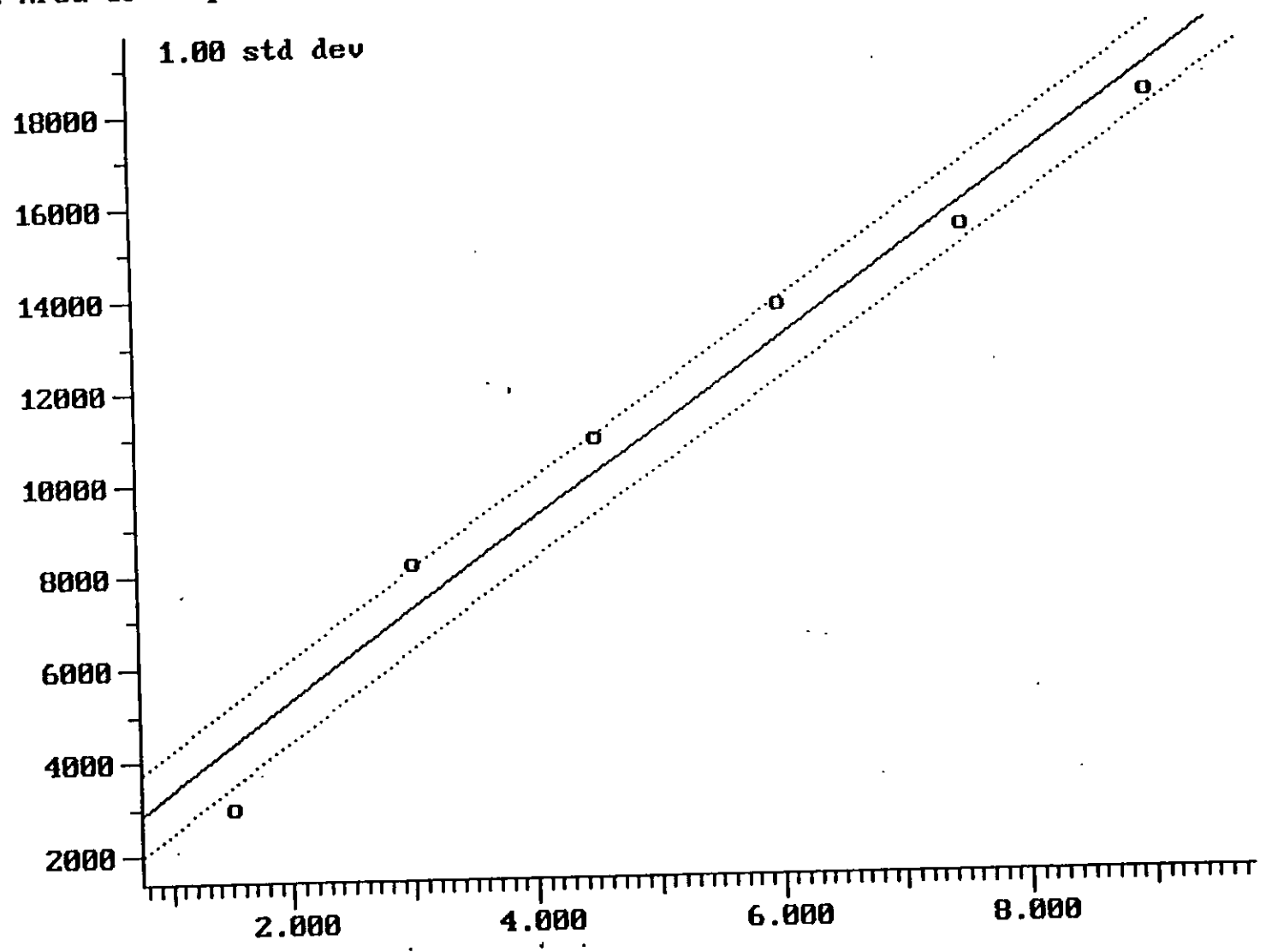
Calibration Plot (Ext Stds) Filename: T014GB
Freon 11 Compound: 7 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.991
Standard Deviation: 1.457
(Lin/Lin)



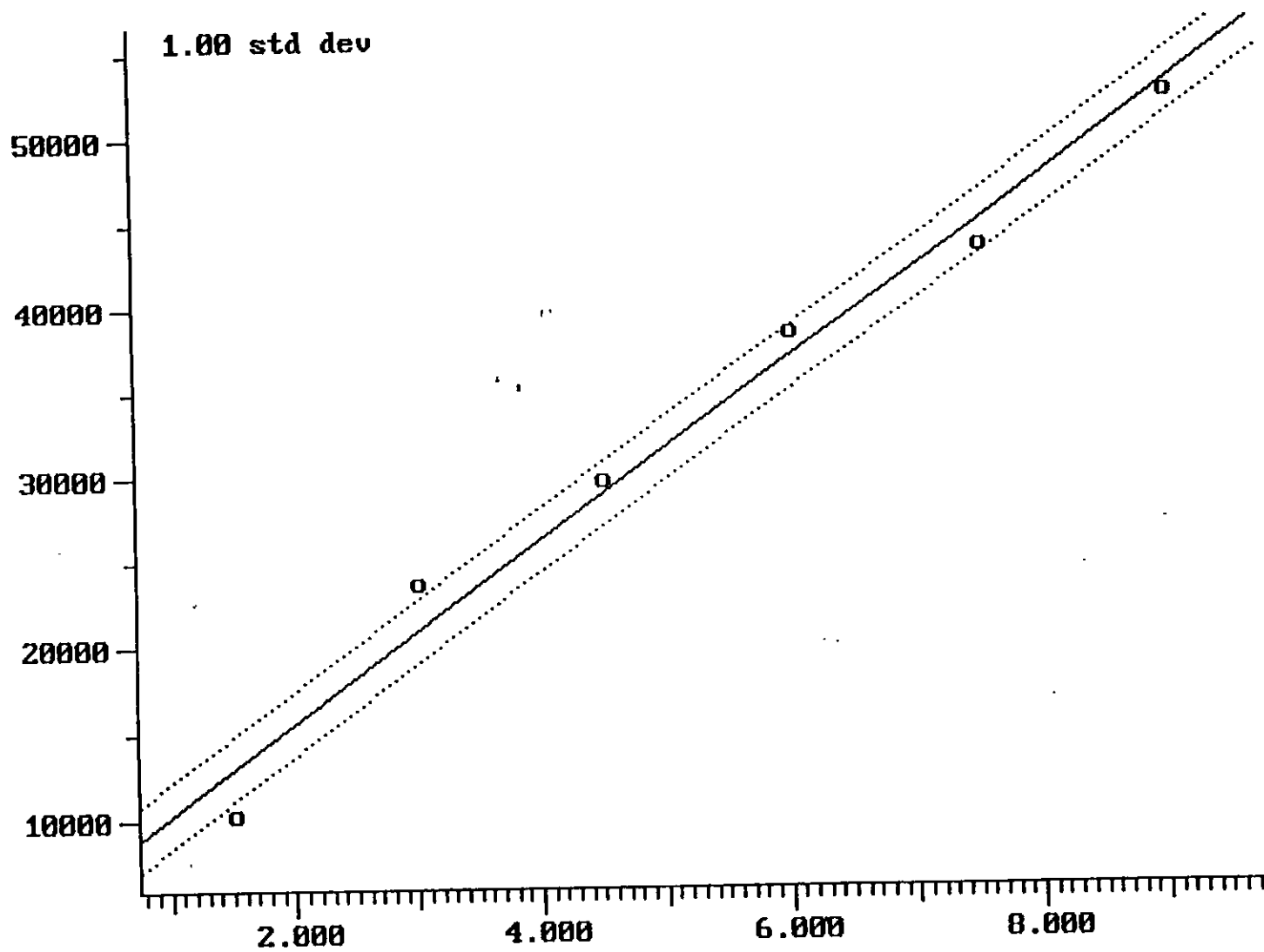
Calibration Plot (Ext Stds) Filename: T014GB
Ethyl Ether Compound: 8 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.985
Standard Deviation: 0.863
(LinLin)



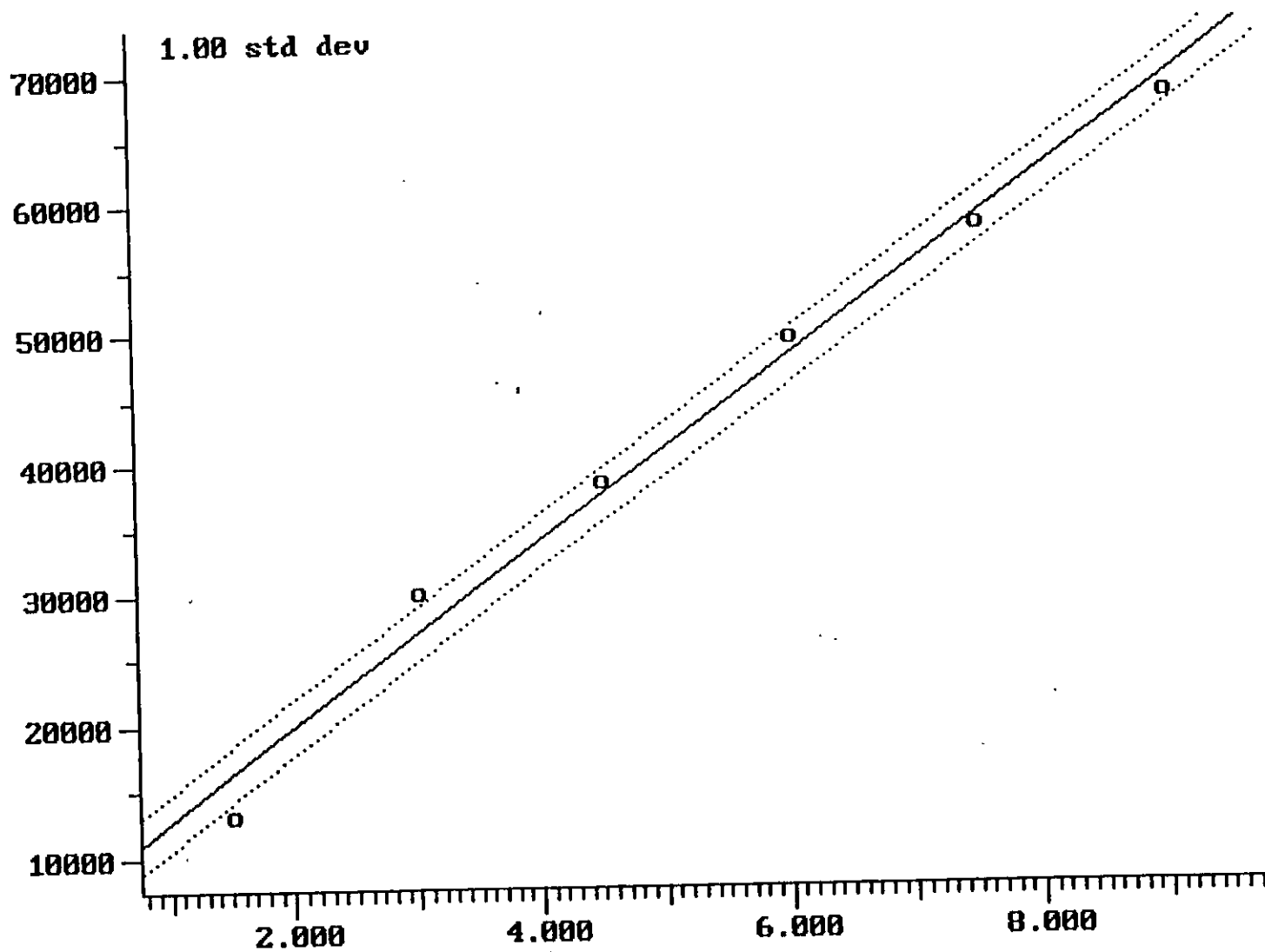
Calibration Plot (Ext Stds) Filename: T014GB
1,1-Dichloroethene Compound: 9 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.991
Standard Deviation: 1.823
(Lin/Lin)



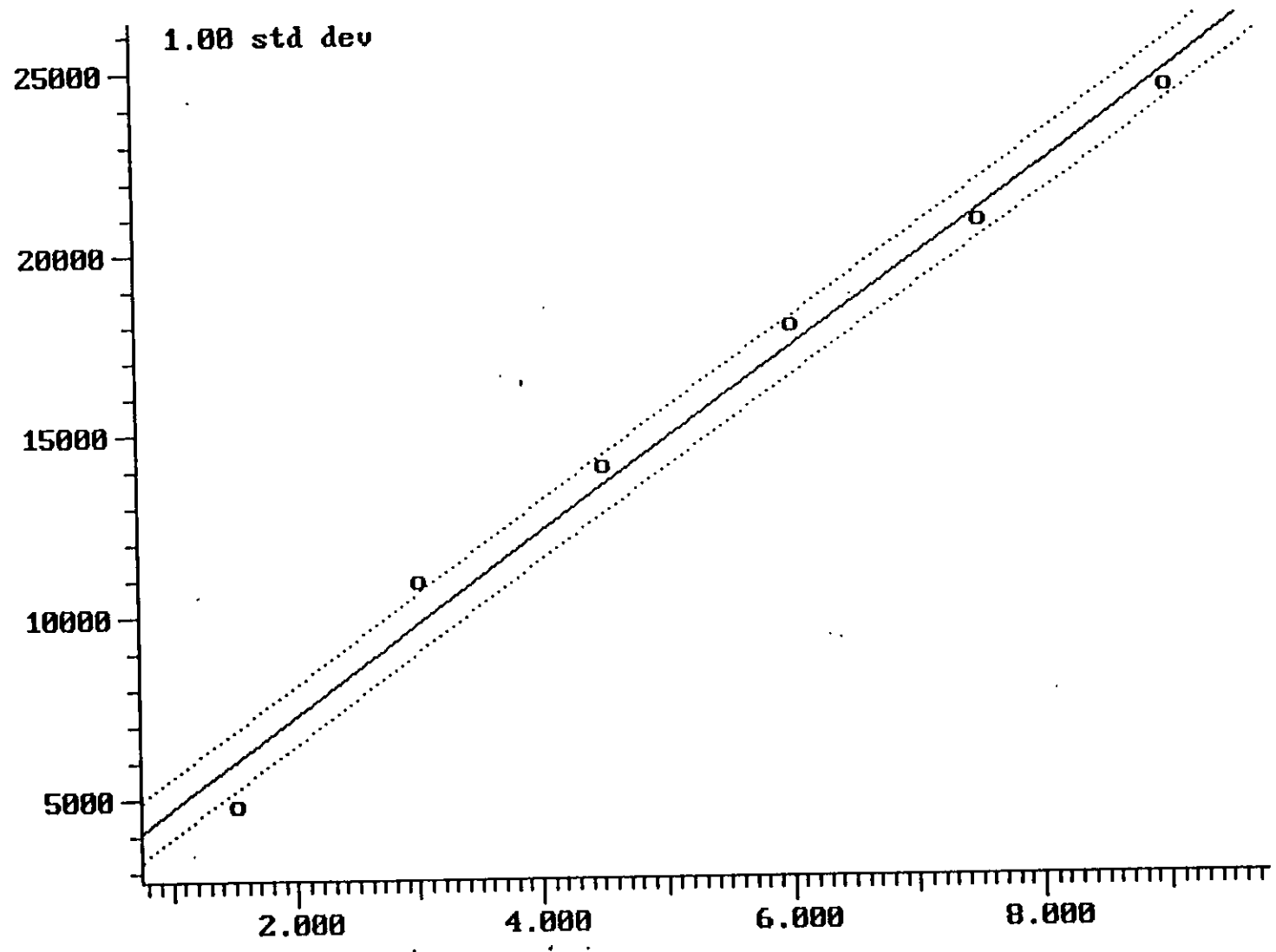
Calibration Plot (Ext Stds) Filename: T014GB
Methylene Chloride Compound: 10 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.994
Standard Deviation: 2.024
(Lin/Lin)



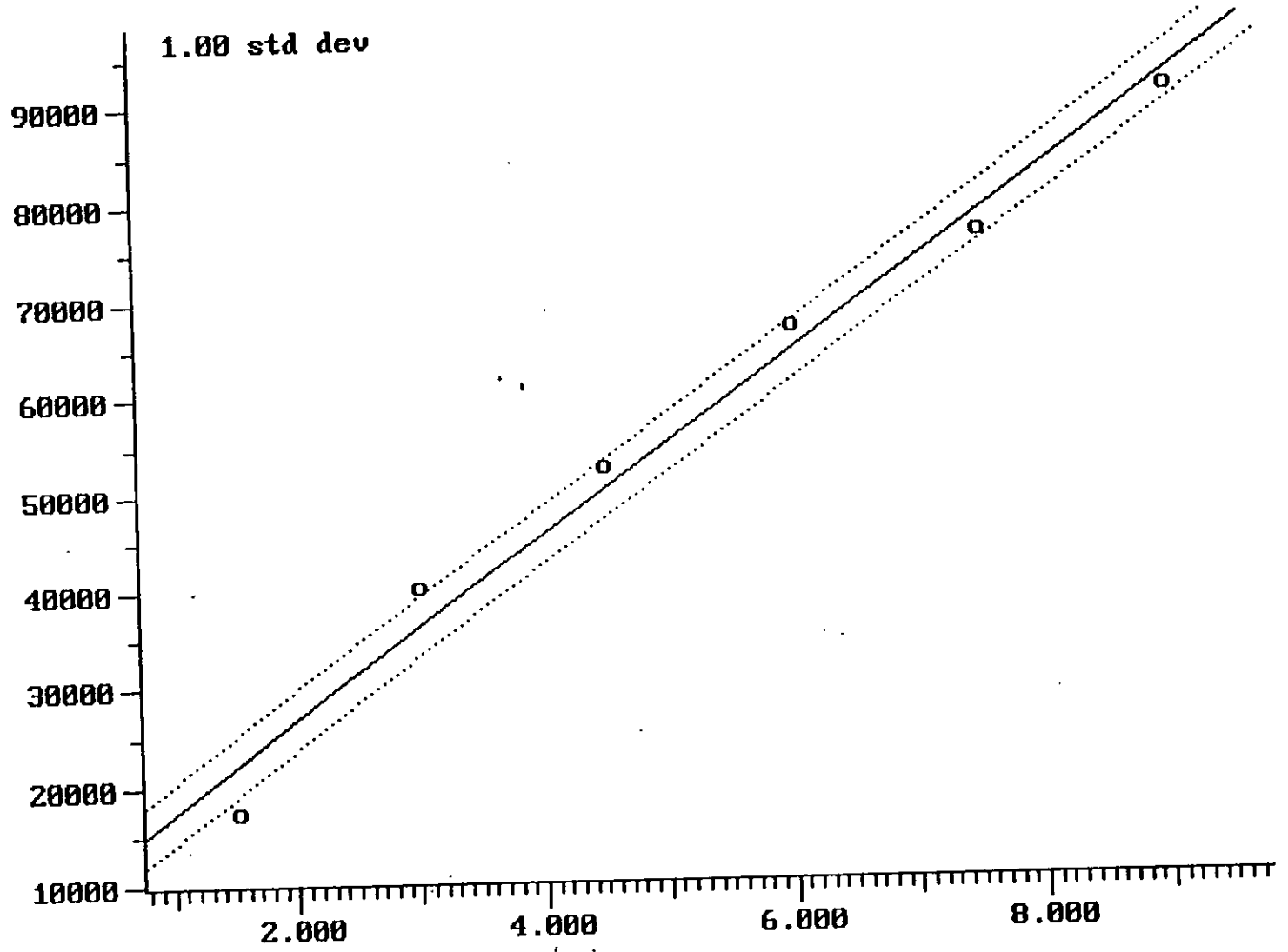
Calibration Plot (Ext Stds) Filename: T014GB
Freon 113 Compound: 11 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.993
Standard Deviation: 0.783
(Lin/Lin)



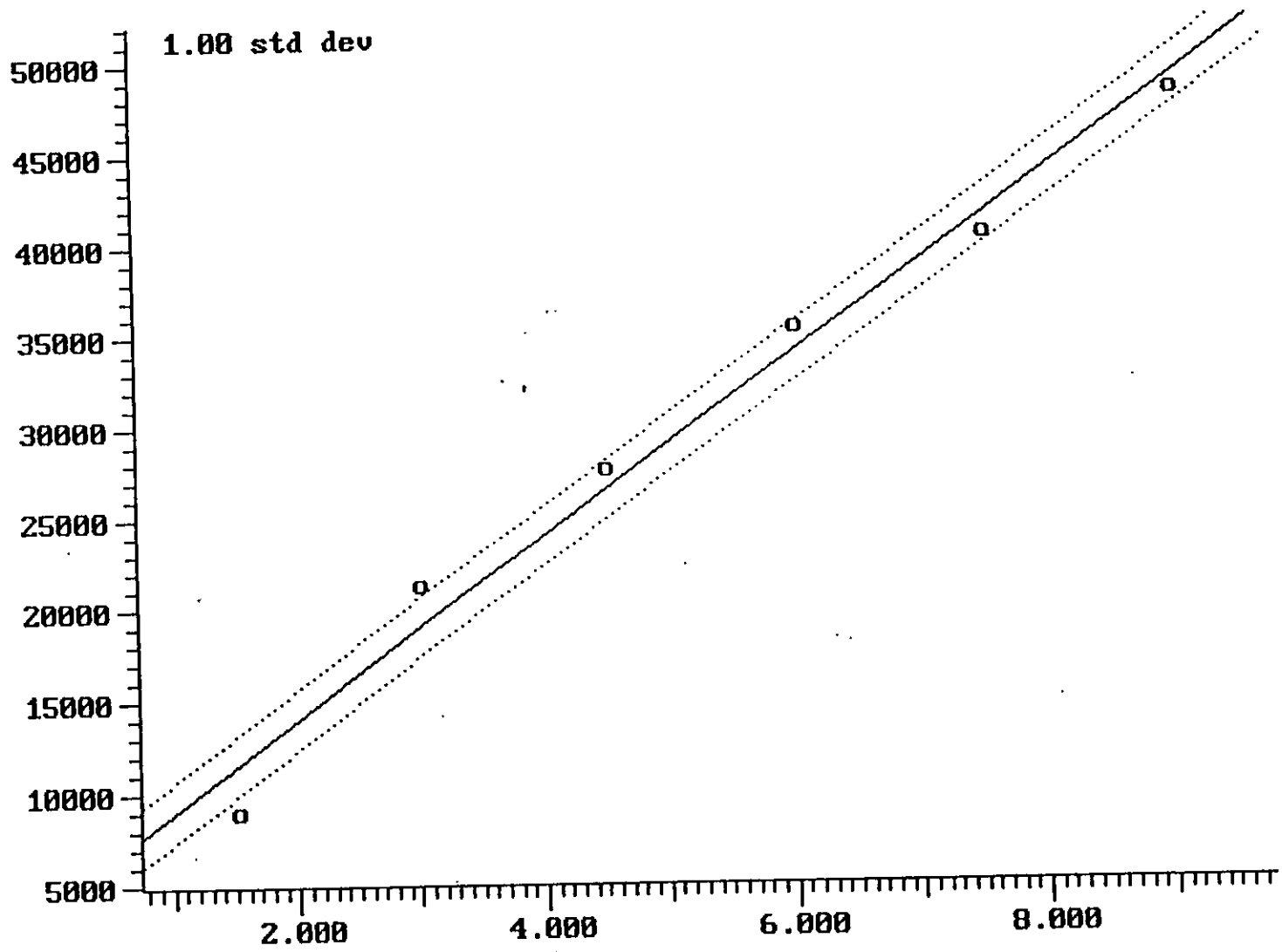
Calibration Plot (Ext Stds) Filename: T014GB
Carbon Disulfide Compound: 12 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.992
Standard Deviation: 3.042
(Lin/Lin)



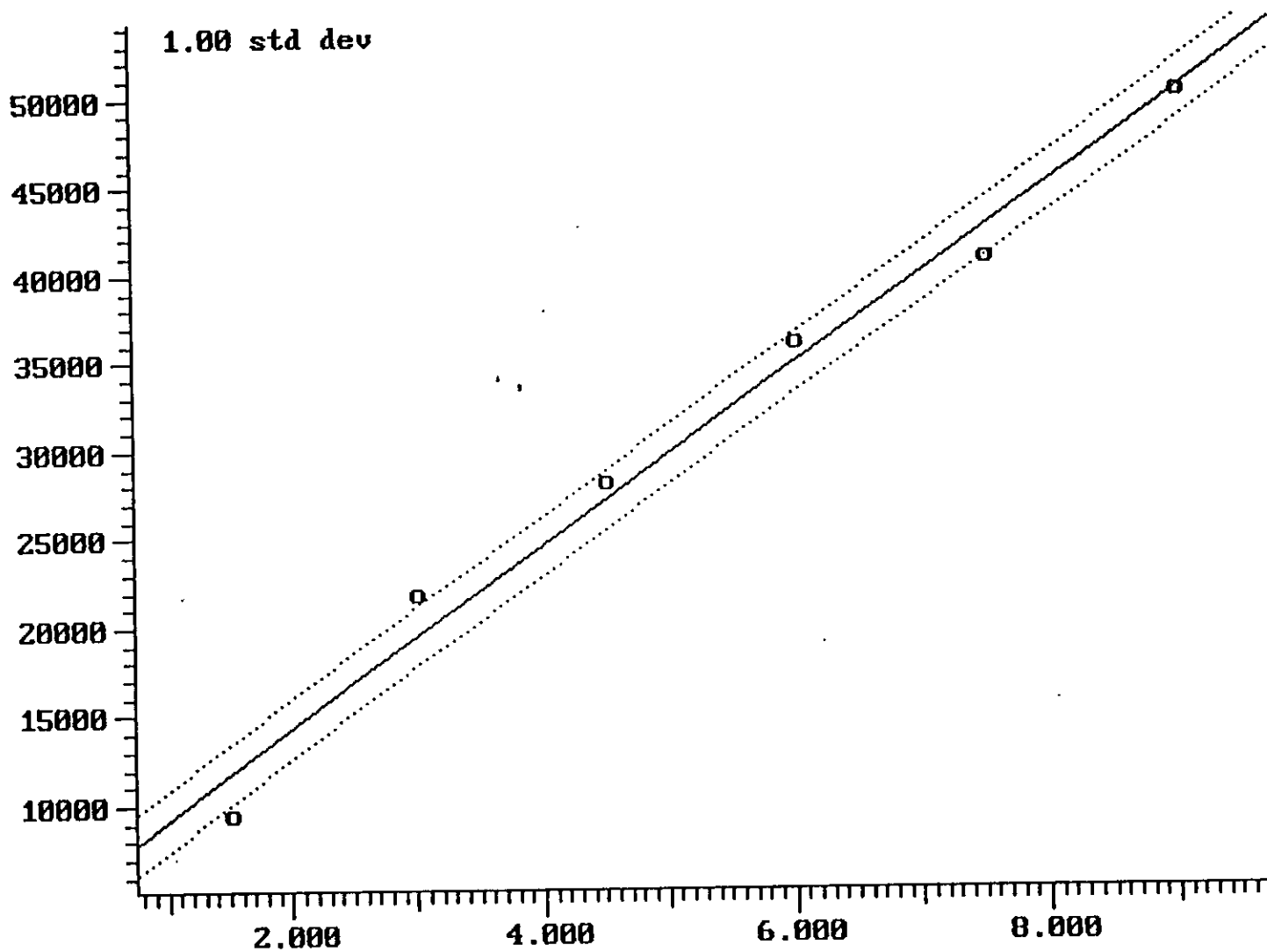
Calibration Plot (Ext Stds) Filename: T014GB
Cis-1,2-Dichloroethene Compound: 13 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.992
Standard Deviation: 1.621
(Lin/Lin)

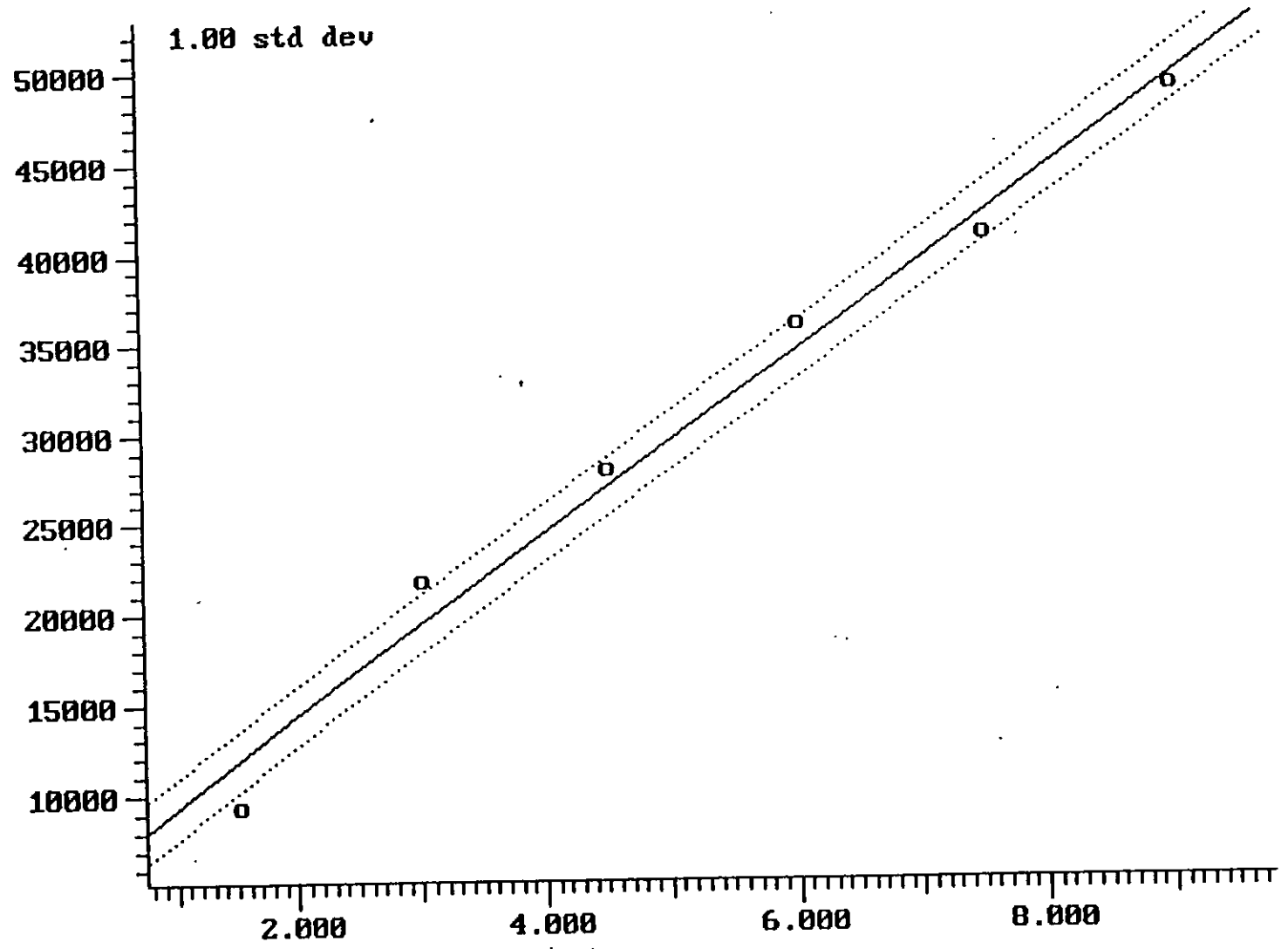


Calibration Plot (Ext Stds) Filename: T014GB
1,1-Dichloroethane Compound: 14 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.992
Standard Deviation: 1.622
(Lin/Lin)

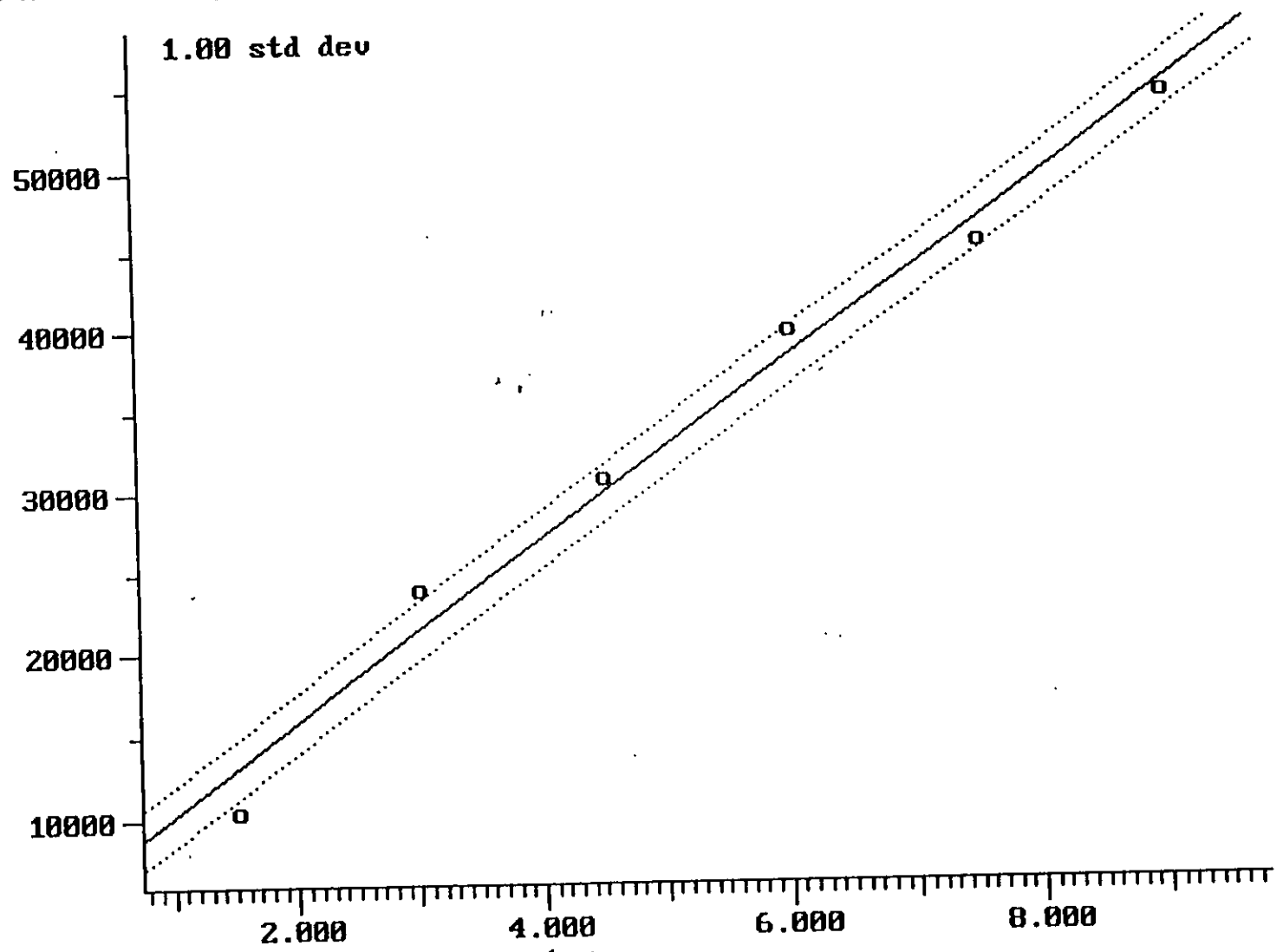


Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.992
Trans-1,2-Dichloroethene Compound: 15 of 45 Standard Deviation: 1.651
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)



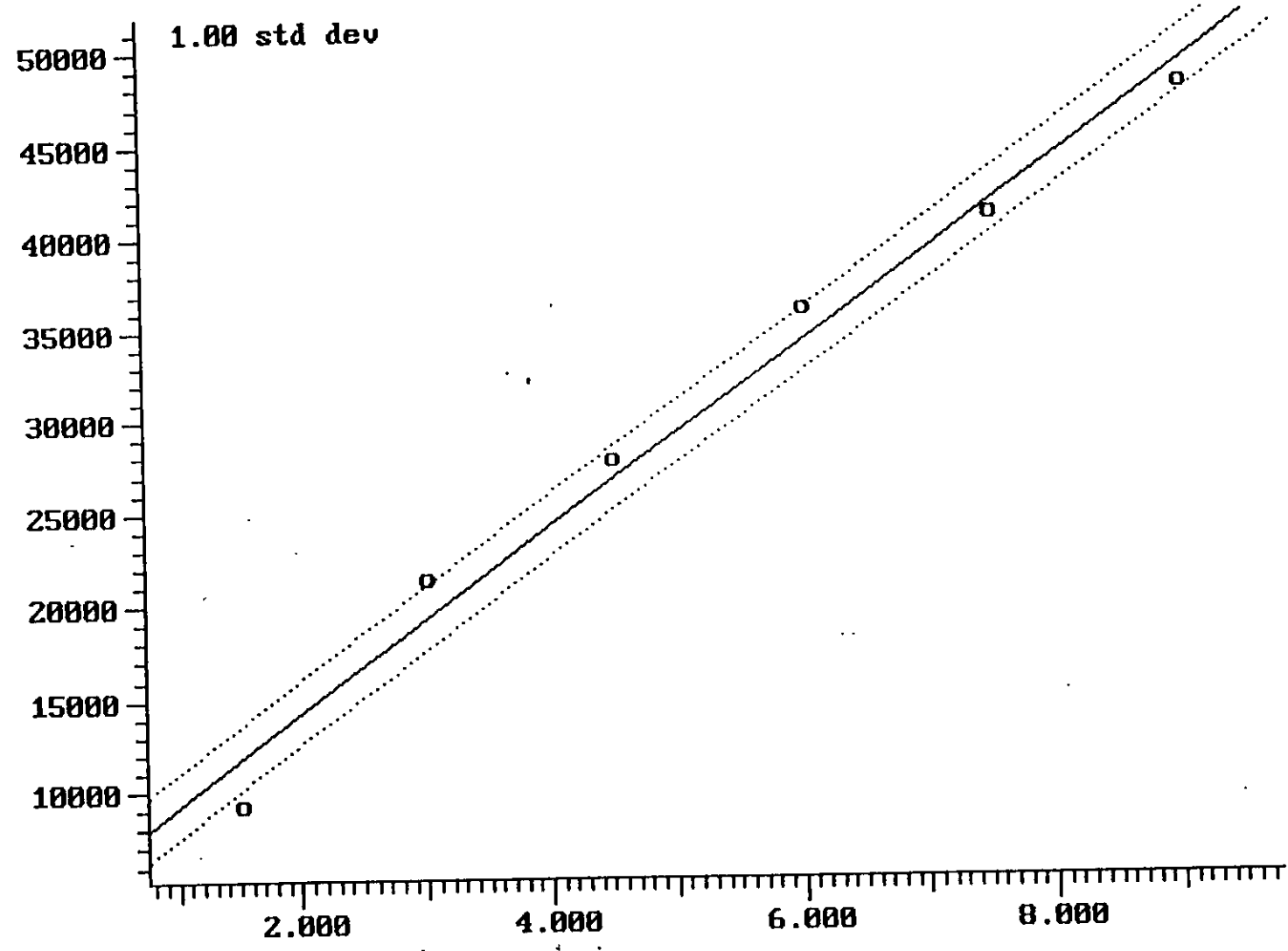
Calibration Plot (Ext Stds) Filename: T014GB
Chloroform Compound: 16 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.992
Standard Deviation: 1.762
(LinLin)



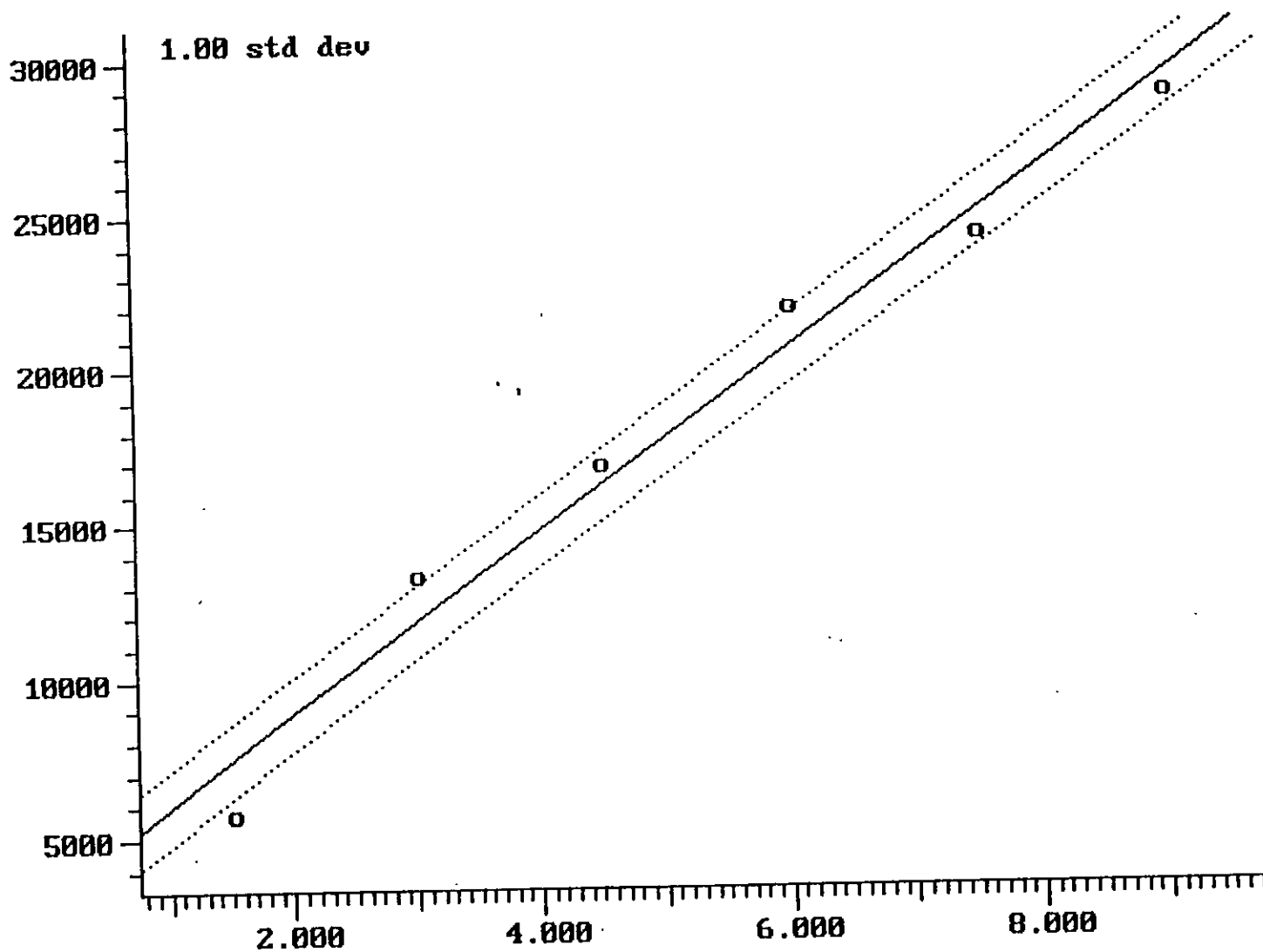
Calibration Plot (Ext Stds) Filename: T014GB
1,2-Dichloroethane Compound: 17 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.992
Standard Deviation: 1.675
(Lin:Lin)



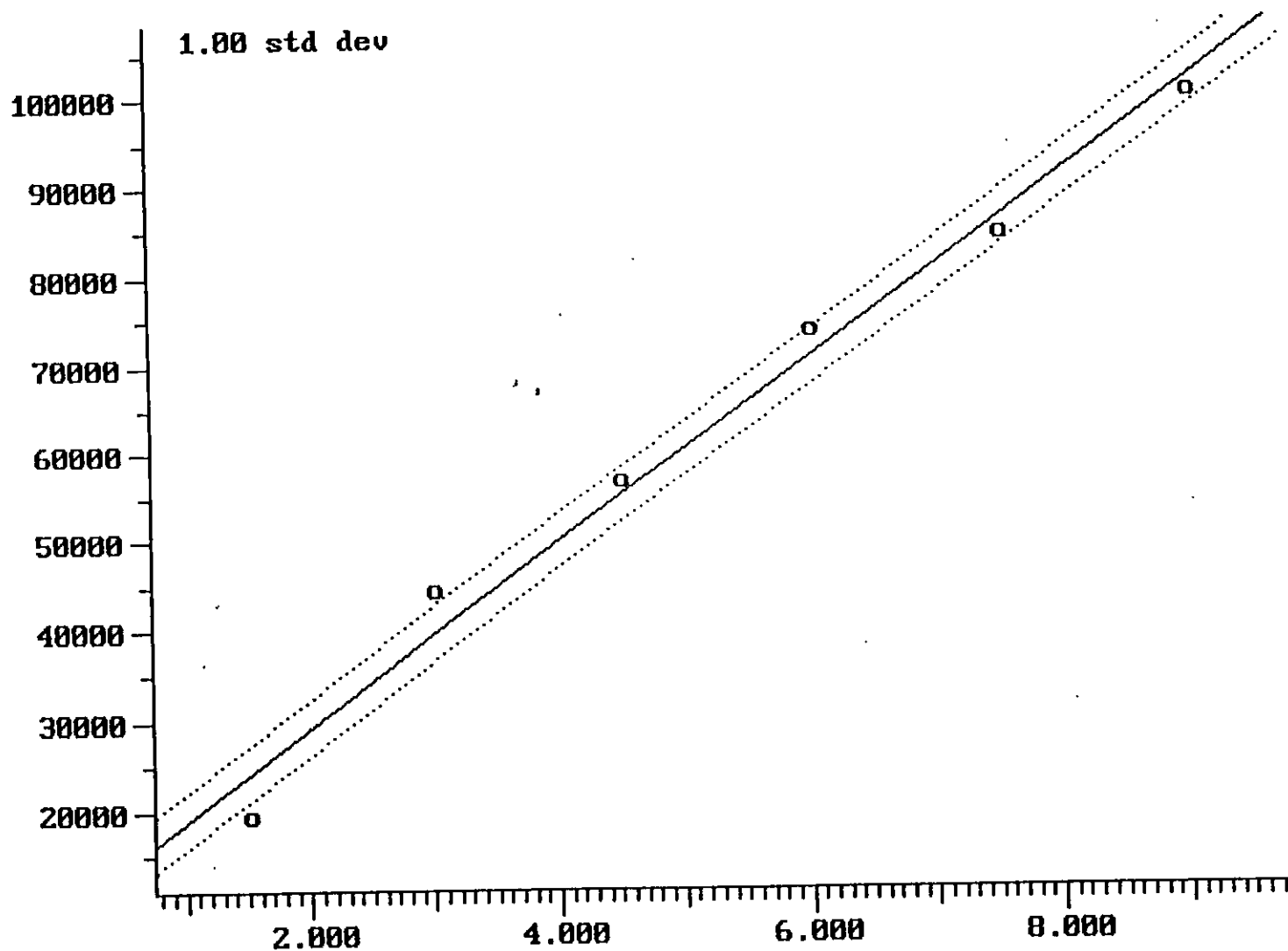
Calibration Plot (Ext Stds) Filename: T014GB
1,1,1-Trichloroethane Compound: 18 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.988
Standard Deviation: 1.163
(LinLin)



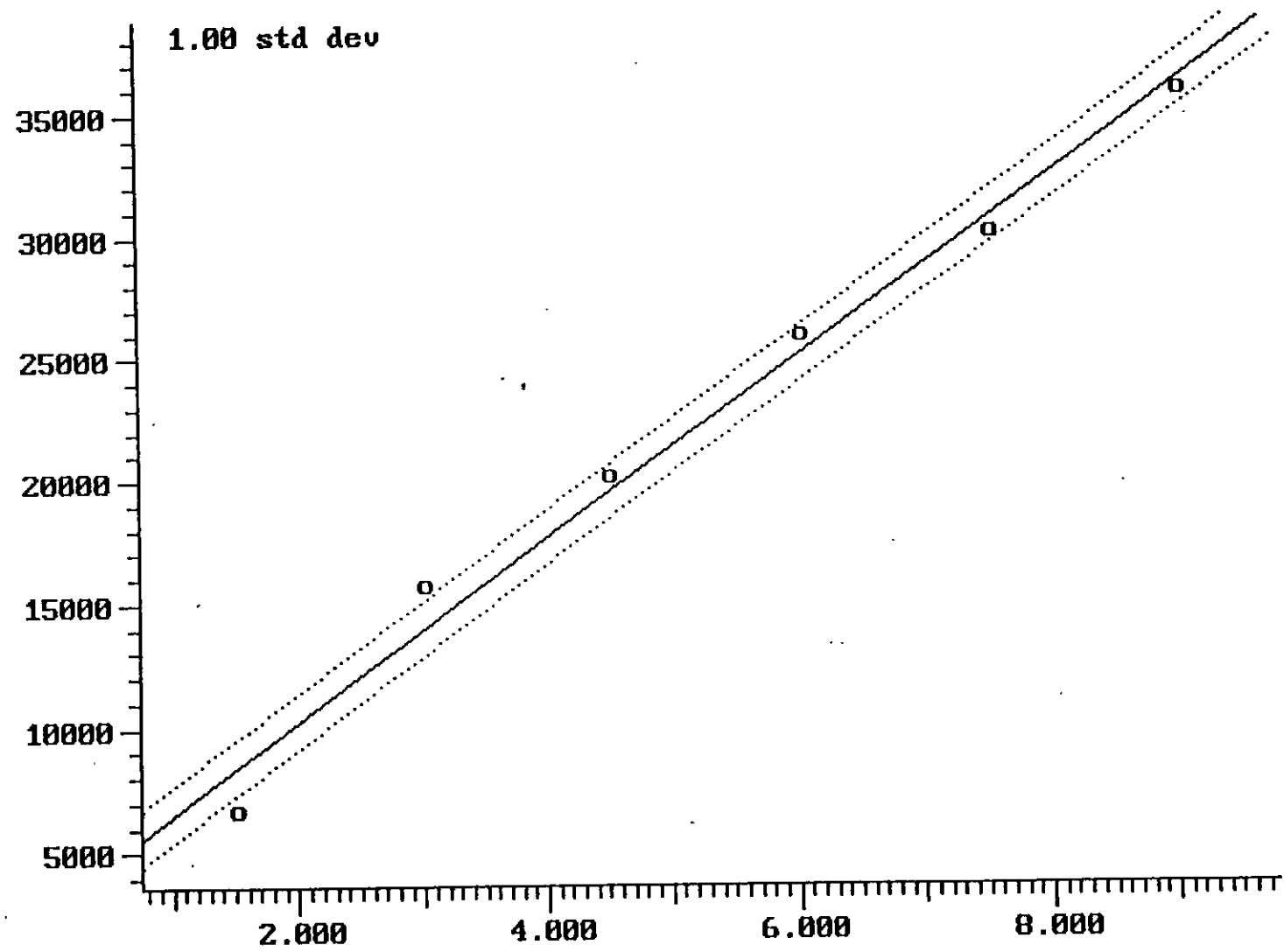
Calibration Plot (Ext Stds) Filename: T014GB
Benzene Compound: 19 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.993
Standard Deviation: 3.199
(Lin/Lin)



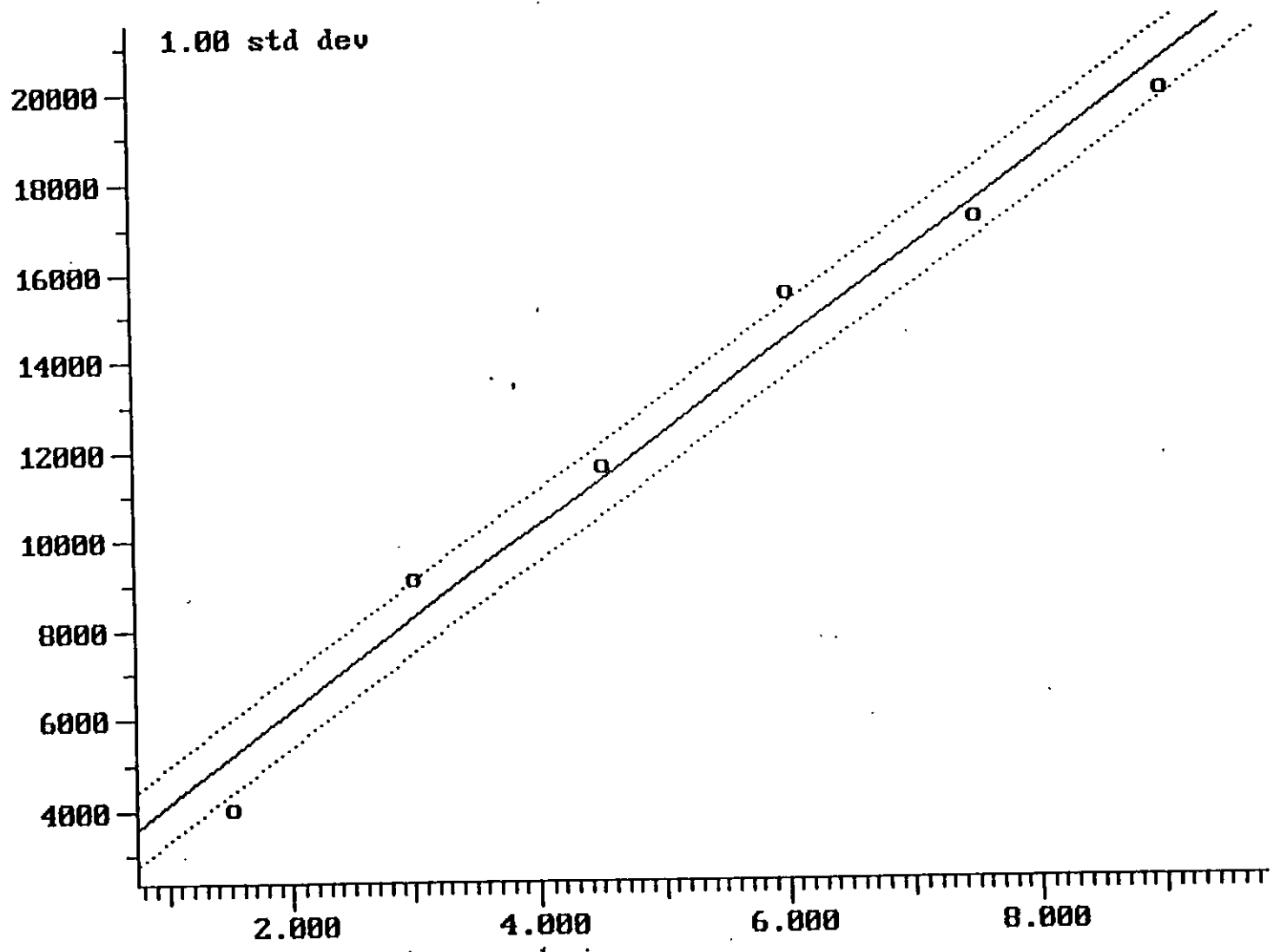
Calibration Plot (Ext Stds) Filename: T014GB
Carbontetrachloride Compound: 20 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.993
Standard Deviation: 1.122
(Lin:Lin)

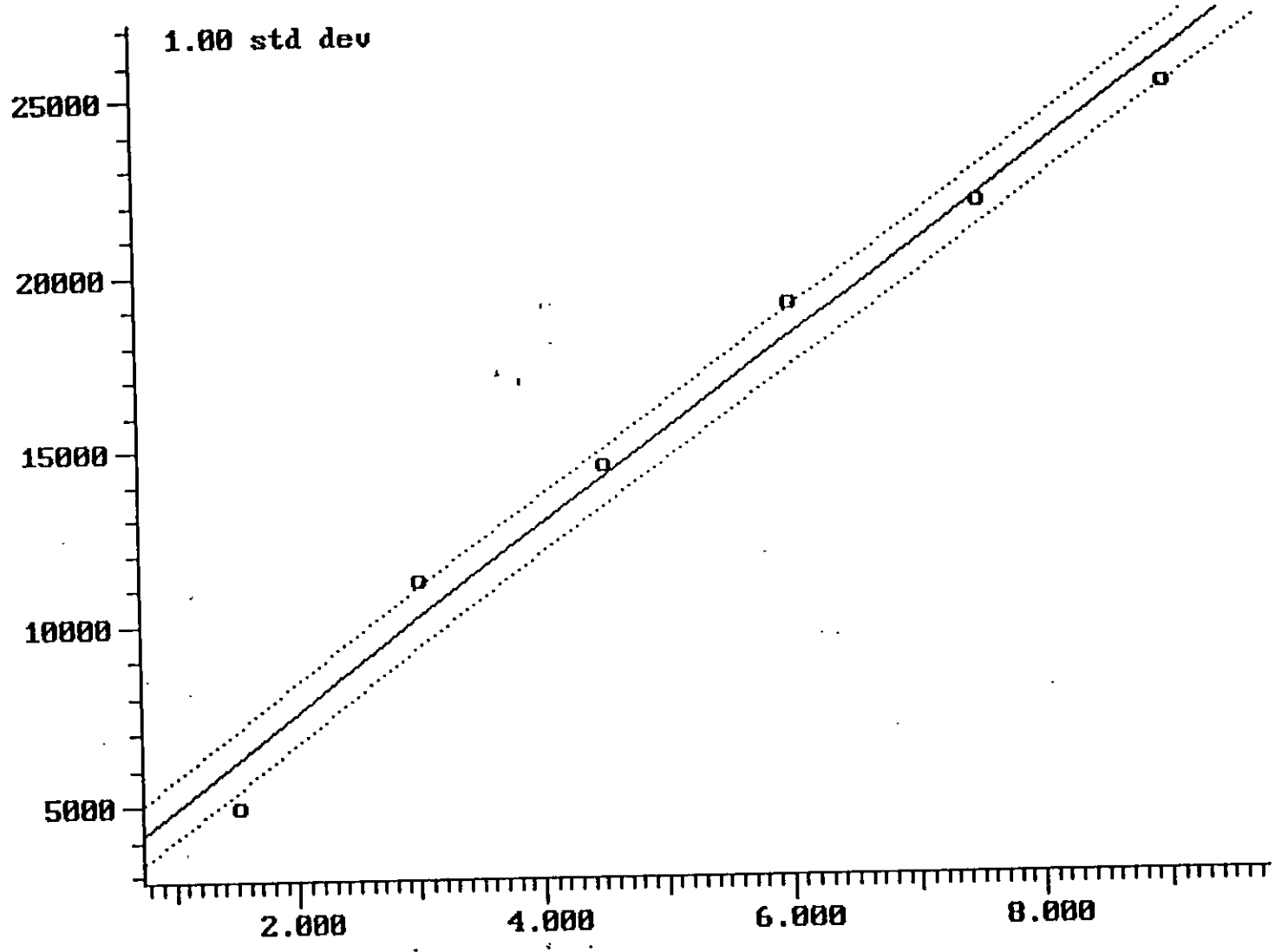


Calibration Plot (Ext Stds) Filename: T014GB
1,2-Dichloropropane Compound: 21 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

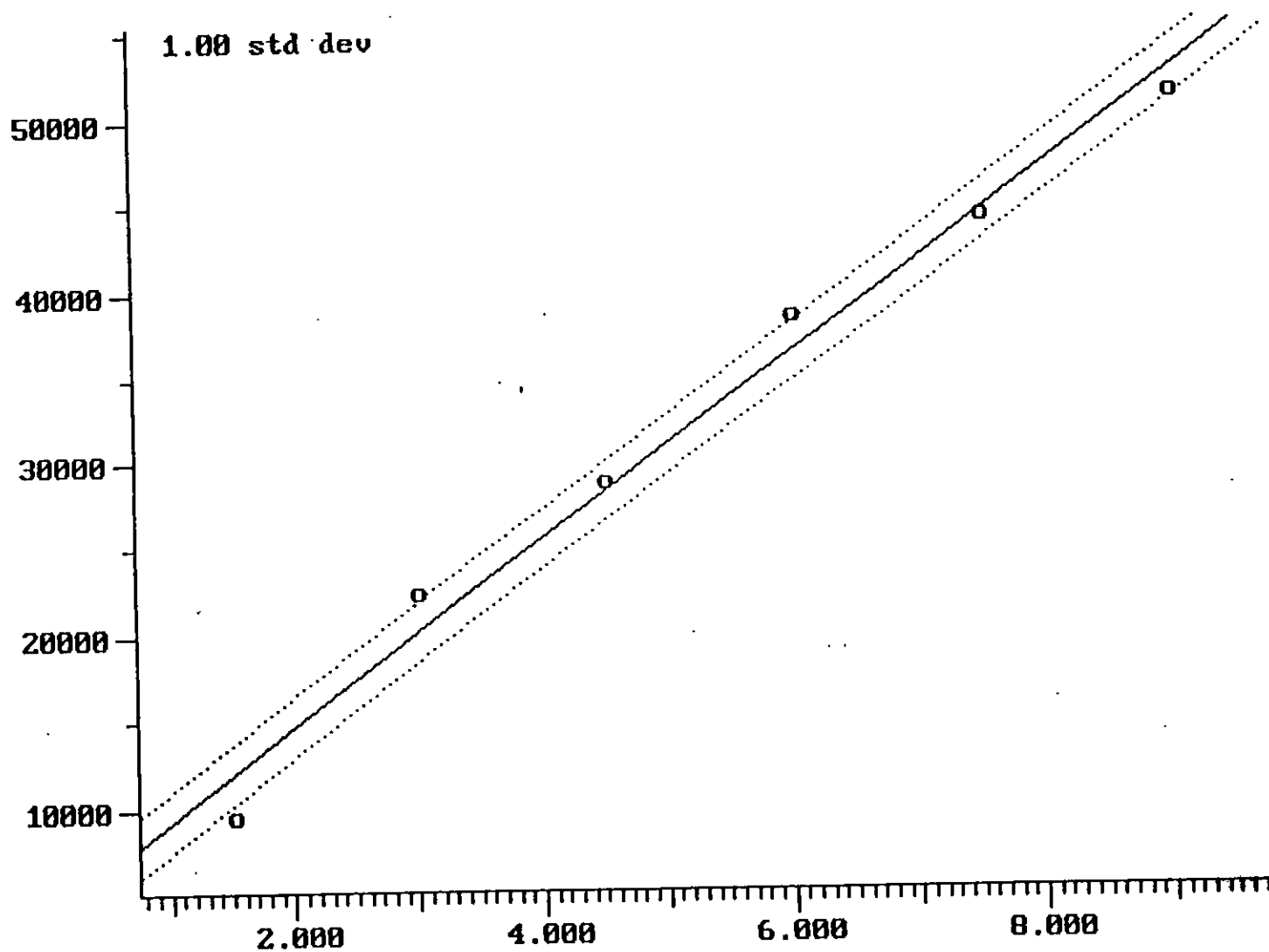
Correlation Coeff: 0.989
Standard Deviation: 0.798
(Lin/Lin)



Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.992
Trichloroethene Compound: 22 of 45 Standard Deviation: 0.863
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)

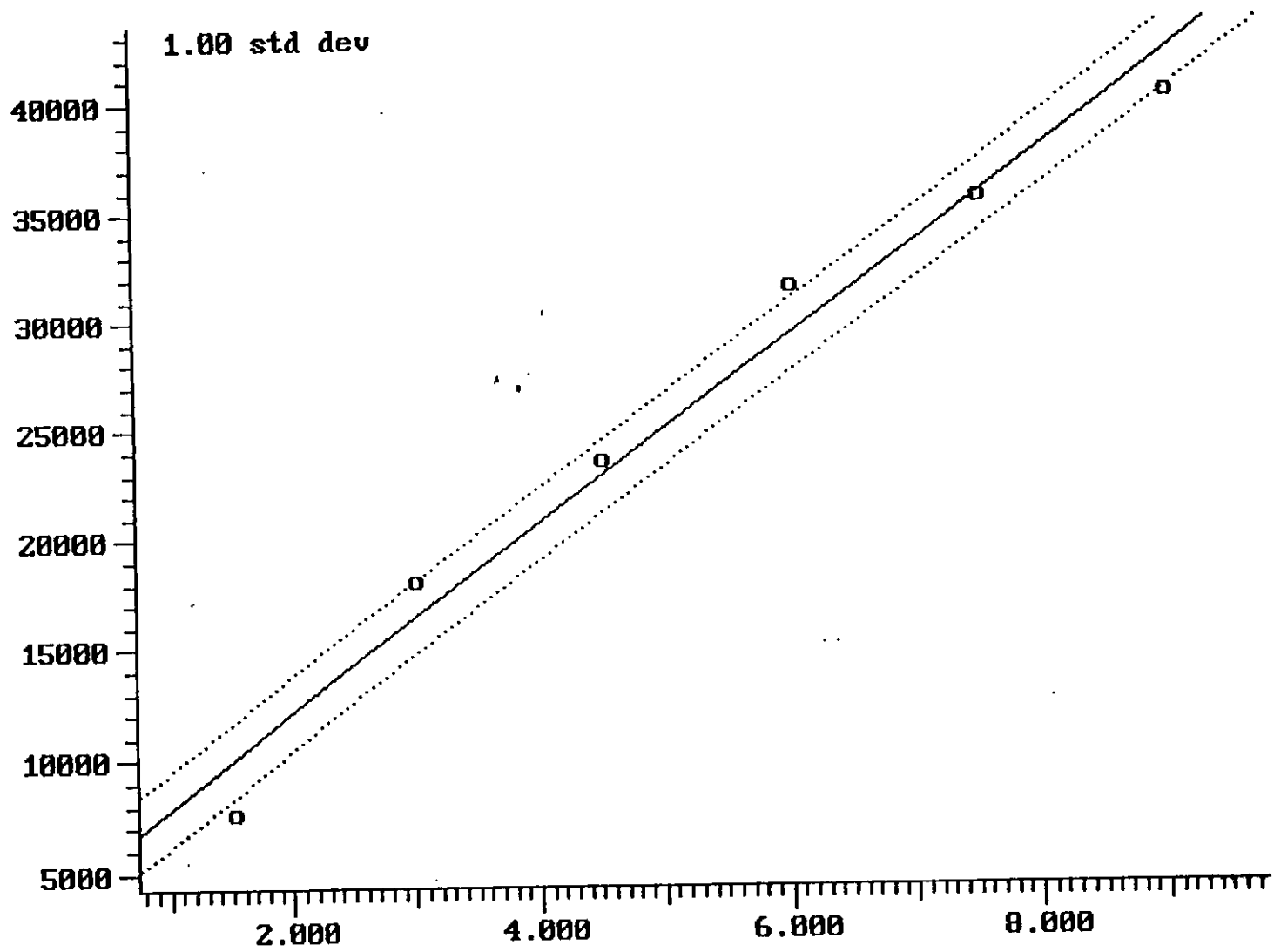


Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.993
Cis-1,3-Dichloropropene Compound: 23 of 45 Standard Deviation: 1.690
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)

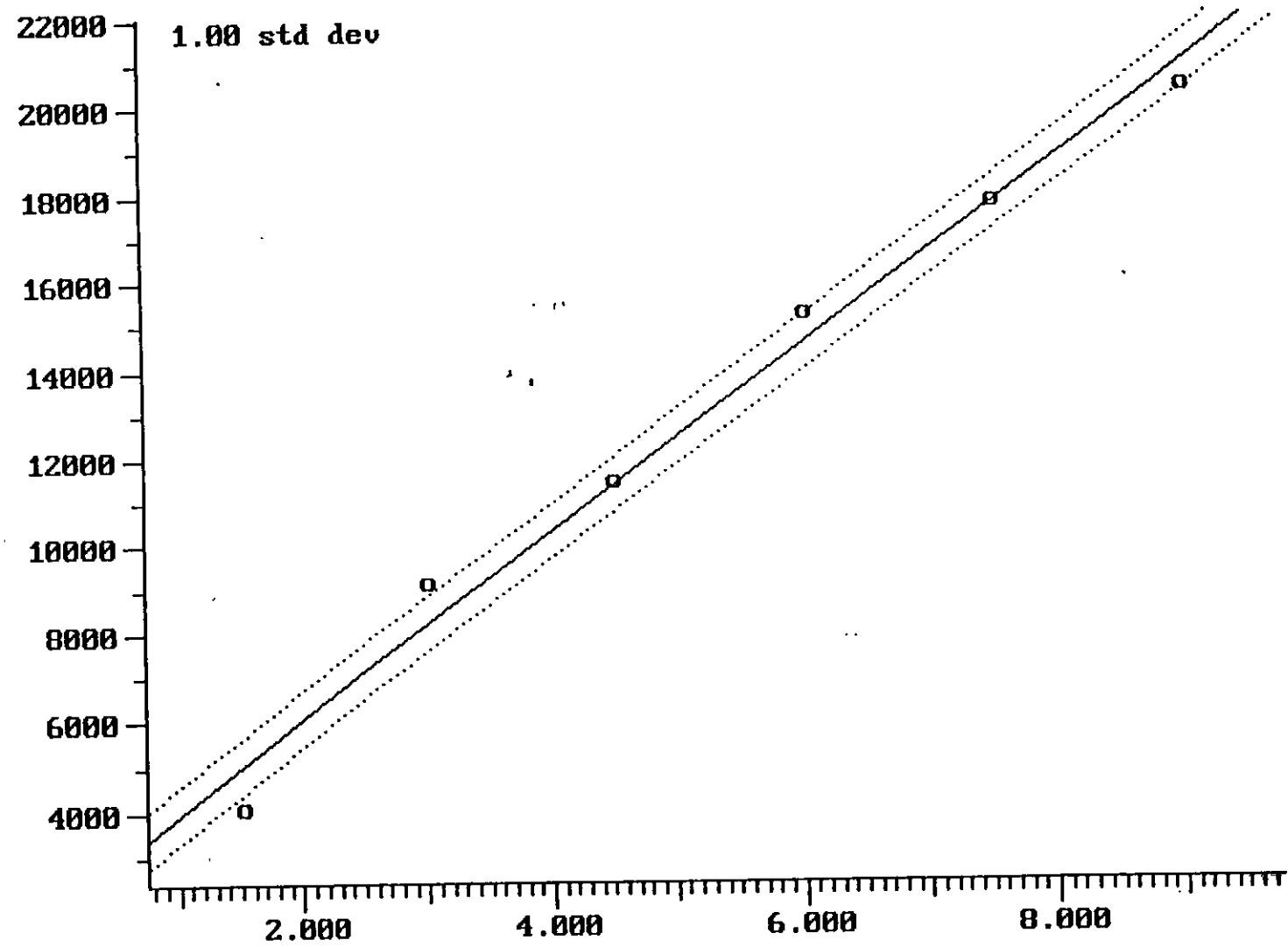


Calibration Plot (Ext Stds) Filename: T014GB
Trans-1,3-Dichloropropene Compound: 24 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.988
Standard Deviation: 1.680
(LinLin)

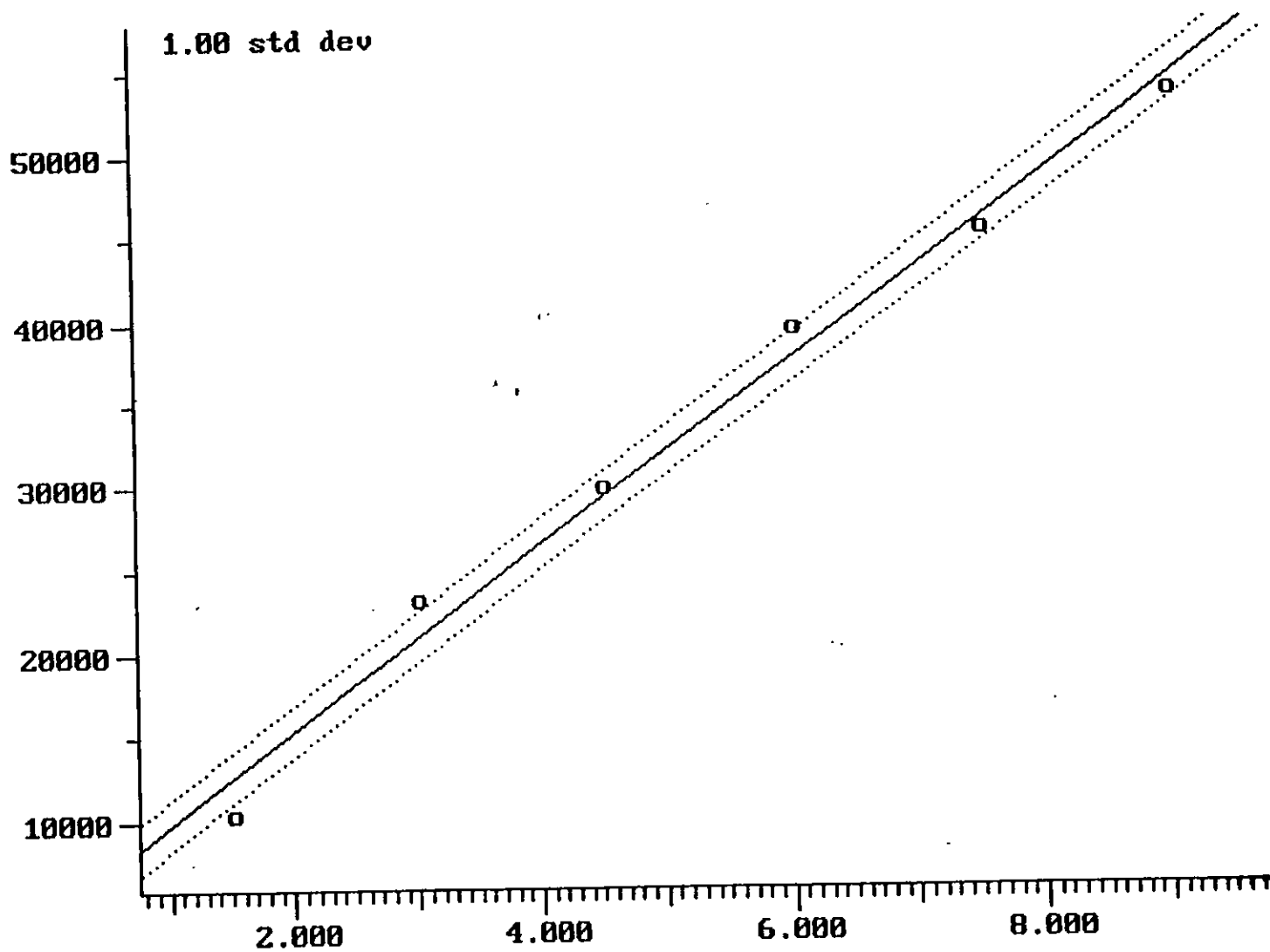


Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.993
1,1,2-Trichloroethane Compound: 25 of 45 Standard Deviation: 0.642
(Peak Area of Sample) vs (Amount of Sample Injected) (LinLin)



Calibration Plot (Ext Stds) Filename: T014GB
Toluene Compound: 26 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

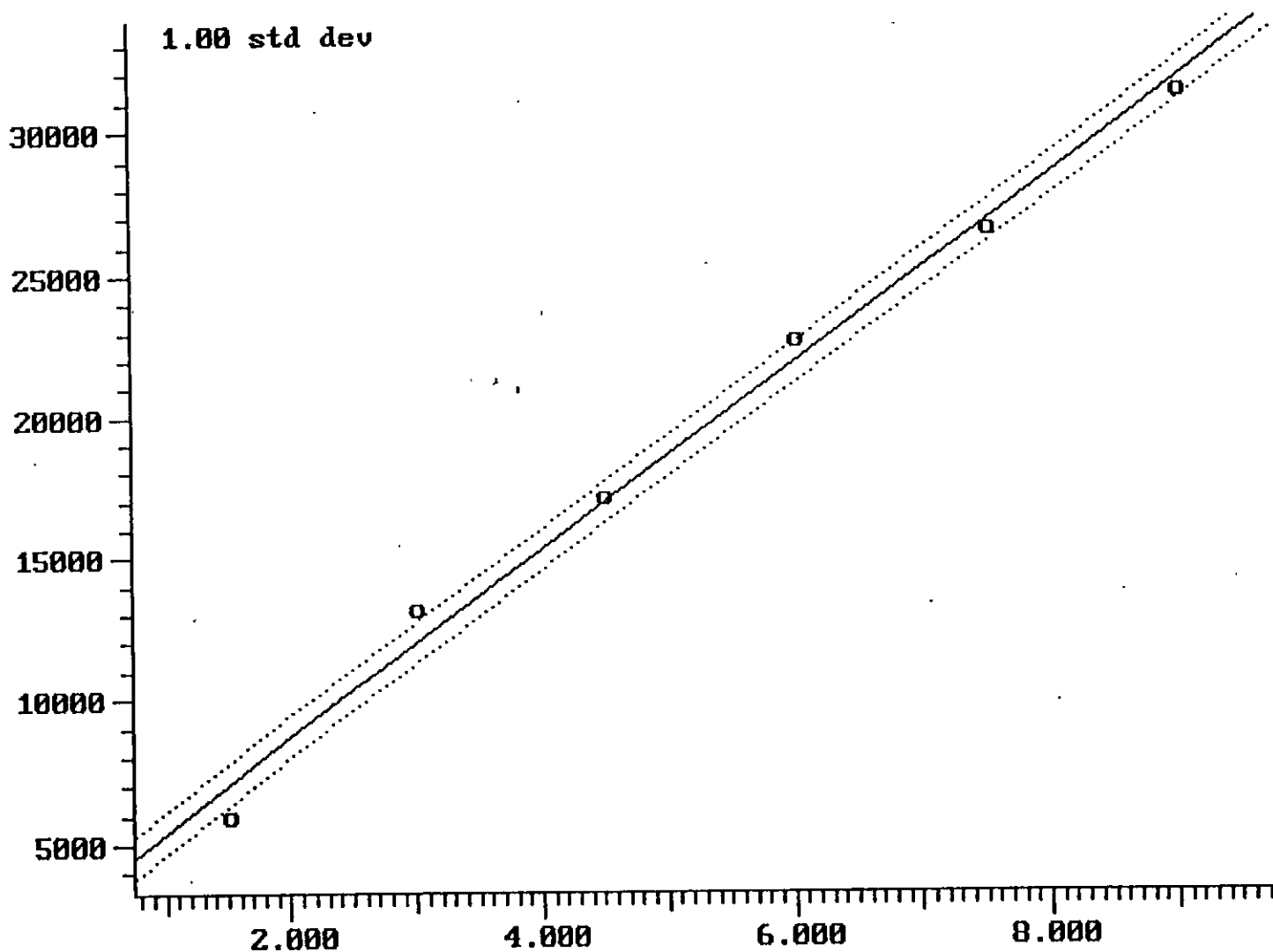
Correlation Coeff: 0.994
Standard Deviation: 1.547
(Lin:Lin)



Calibration Plot (Ext Stds)
1,2-Dibromoethane
(Peak Area of Sample) vs (Amount of Sample Injected)

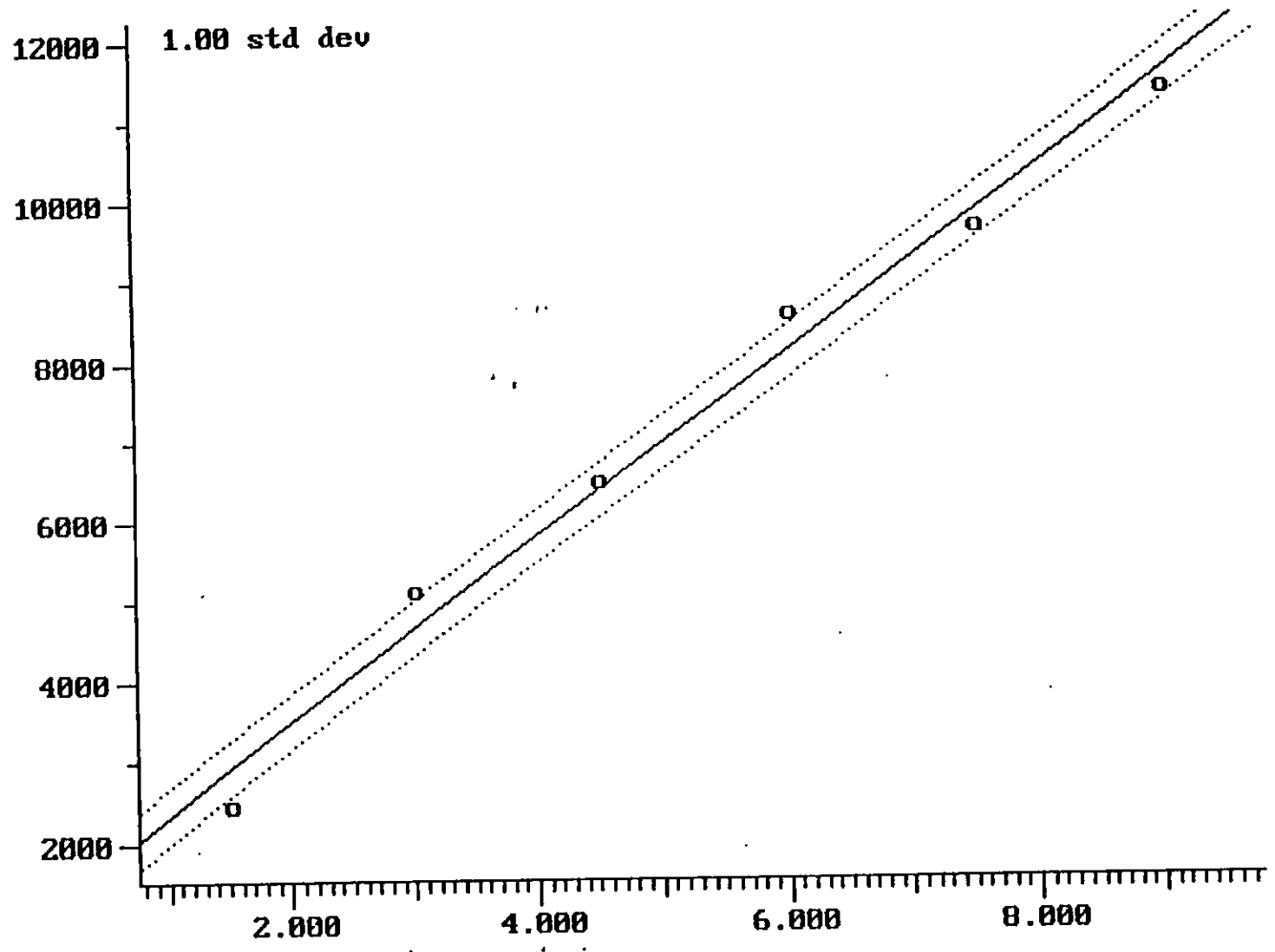
Filename: T014GB
Compound: 27 of 45

Correlation Coeff: 0.996
Standard Deviation: 0.768
(Lin/Lin)

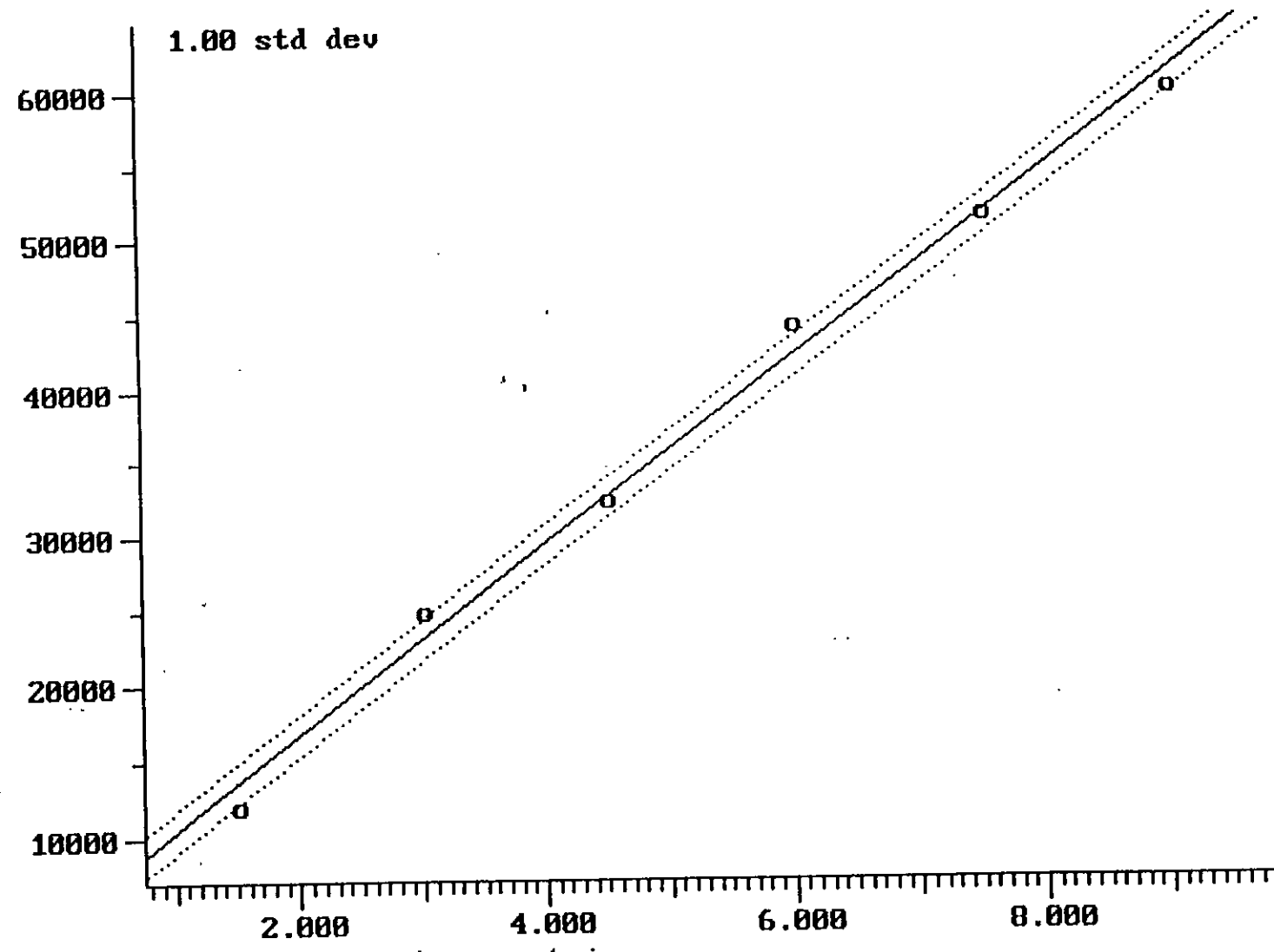


Calibration Plot (Ext Stds) Filename: T014GB
Tetrachloroethene Compound: 28 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

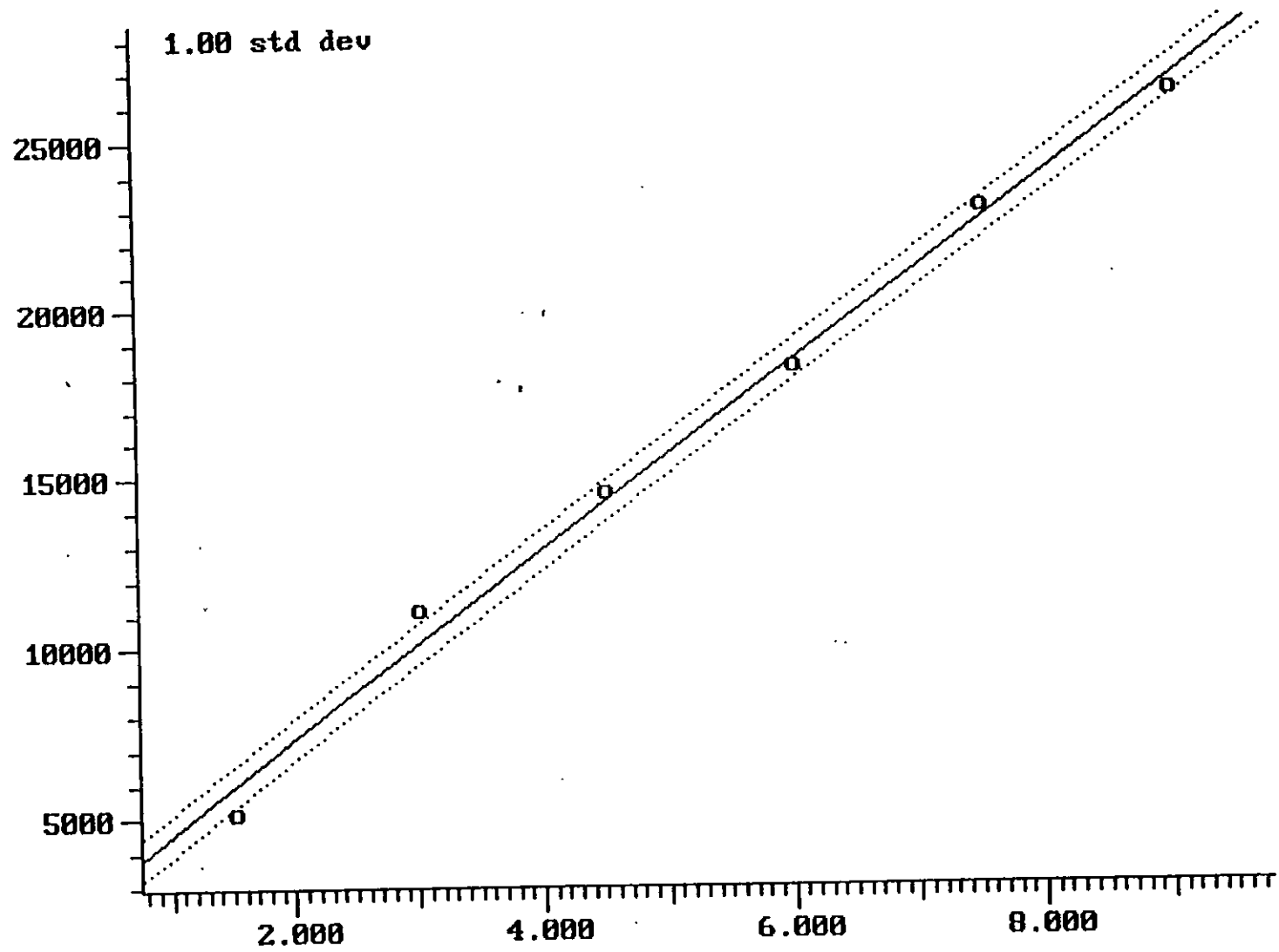
Correlation Coeff: 0.993
Standard Deviation: 0.342
(Lin/Lin)



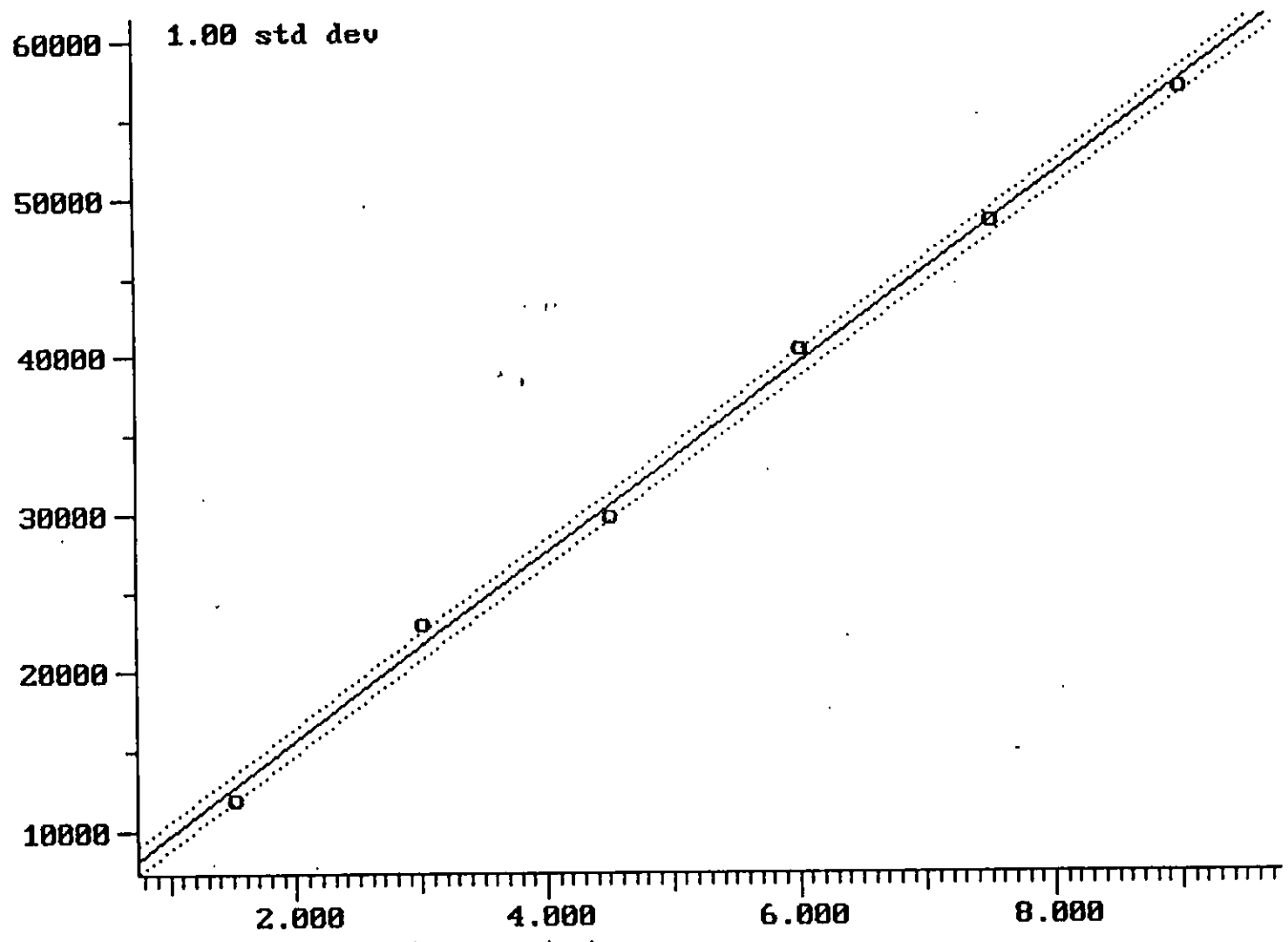
Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.996
Chlorobenzene Compound: 29 of 45 Standard Deviation: 1.380
(Peak Area of Sample) vs (Amount of Sample Injected) (LinLin)



Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.997
Ethylbenzene Compound: 30 of 45 Standard Deviation: 0.577
(Peak Area of Sample) vs (Amount of Sample Injected) (LinLin)

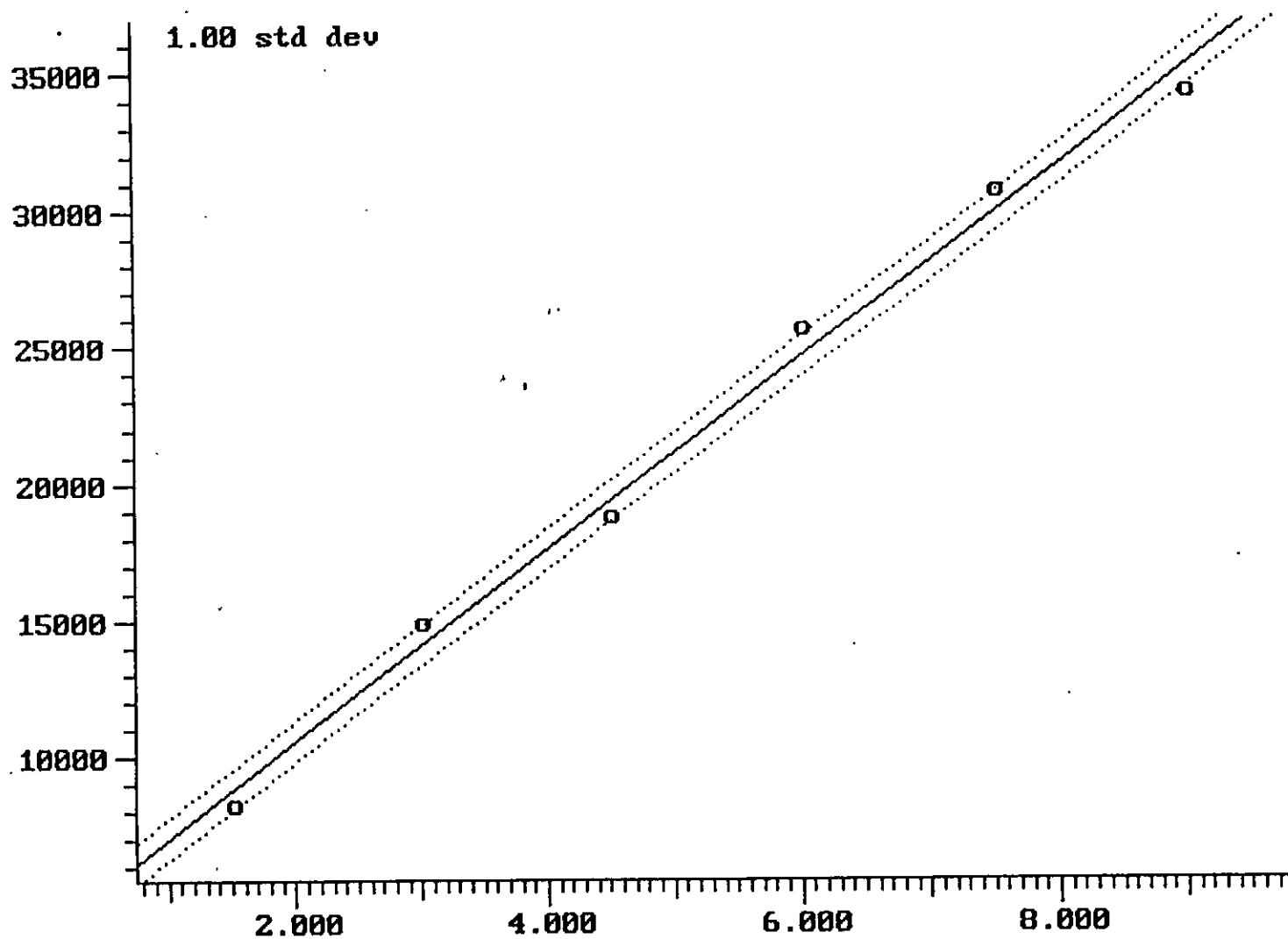


Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.999
Styrene Compound: 32 of 45 Standard Deviation: 0.809
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)

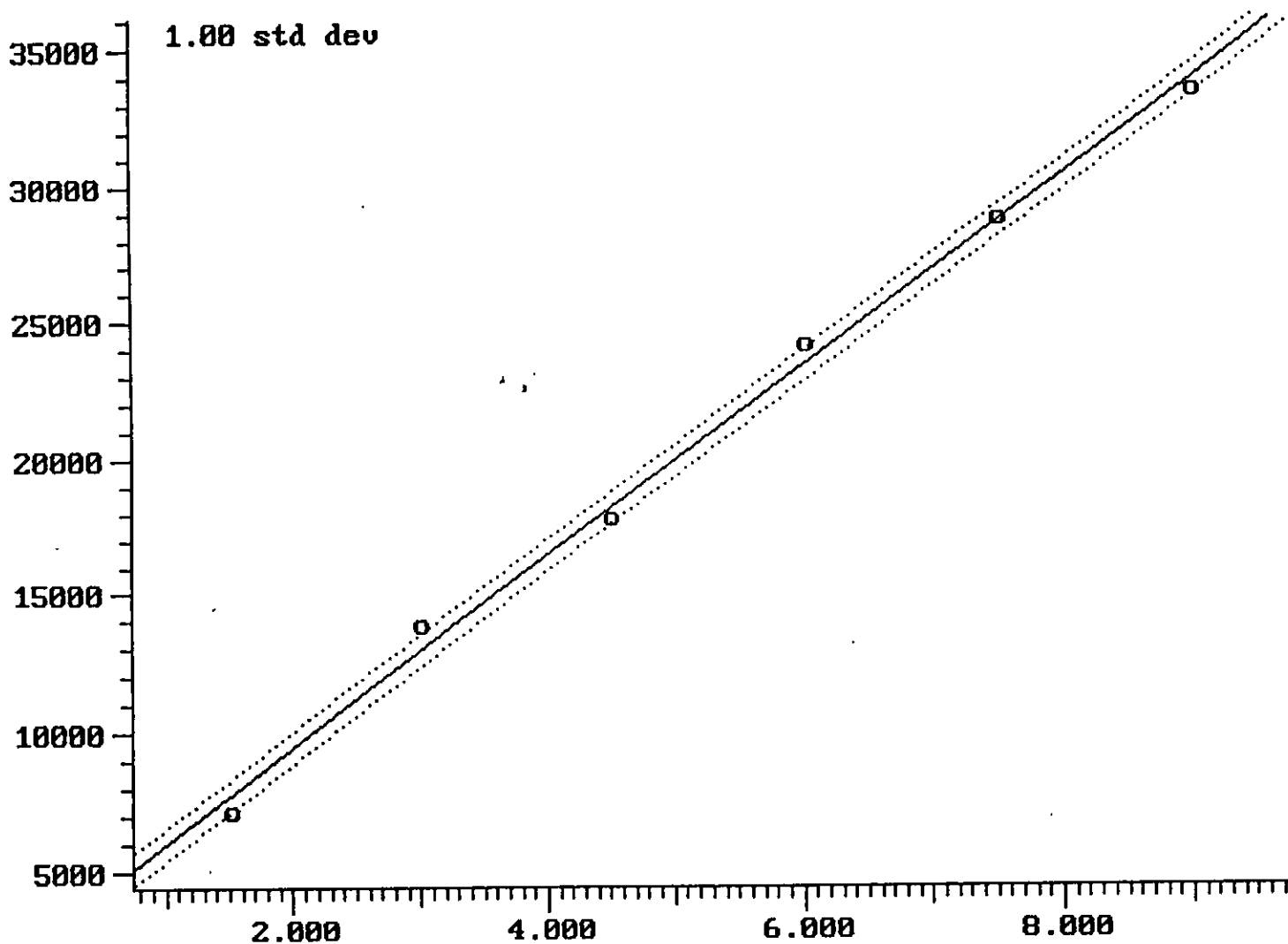


Calibration Plot (Ext Stds) Filename: T014GB
1,1,2,2-Tetrachloroethane Compound: 33 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

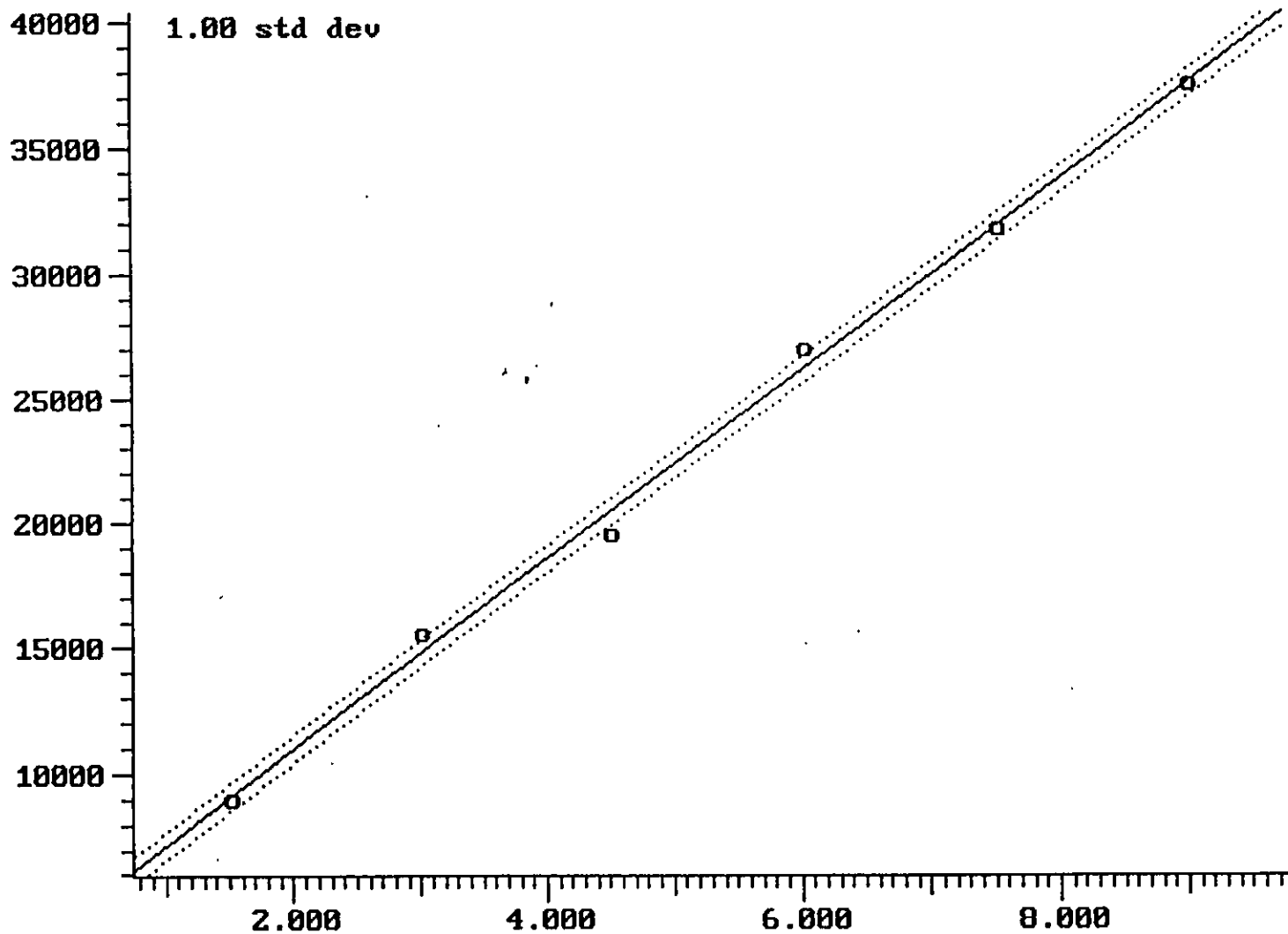
Correlation Coeff: 0.996
Standard Deviation: 0.786
(Lin/Lin)



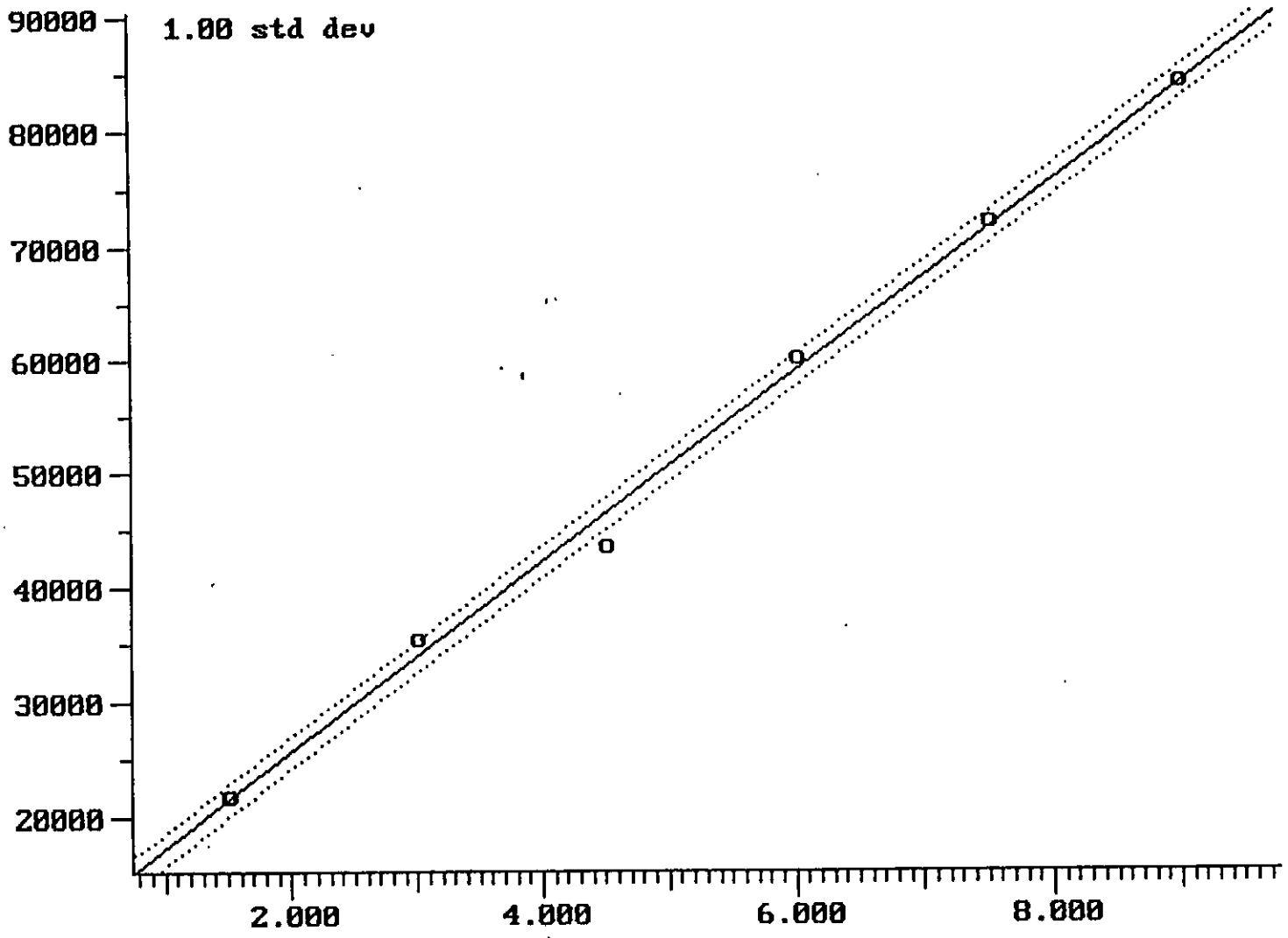
Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.998
1,2-Xylene Compound: 34 of 45 Standard Deviation: 0.579
(Peak Area of Sample) vs (Amount of Sample Injected) (LinLin)



Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.998
4-Bromofluorobenzene Compound: 35 of 45 Standard Deviation: 0.585
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)

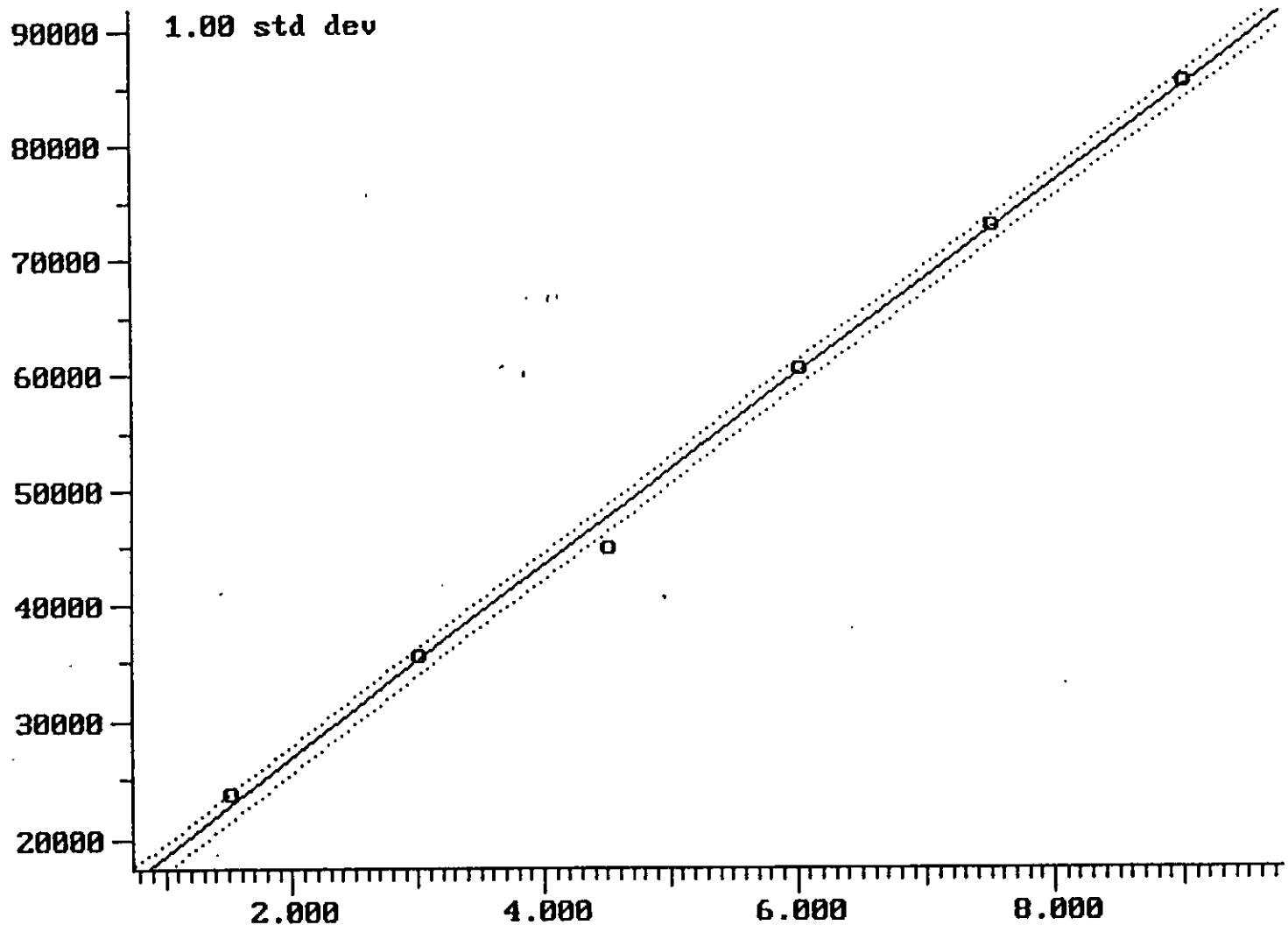


Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.998
Benzylchloride Compound: 36 of 45 Standard Deviation: 1.452
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)



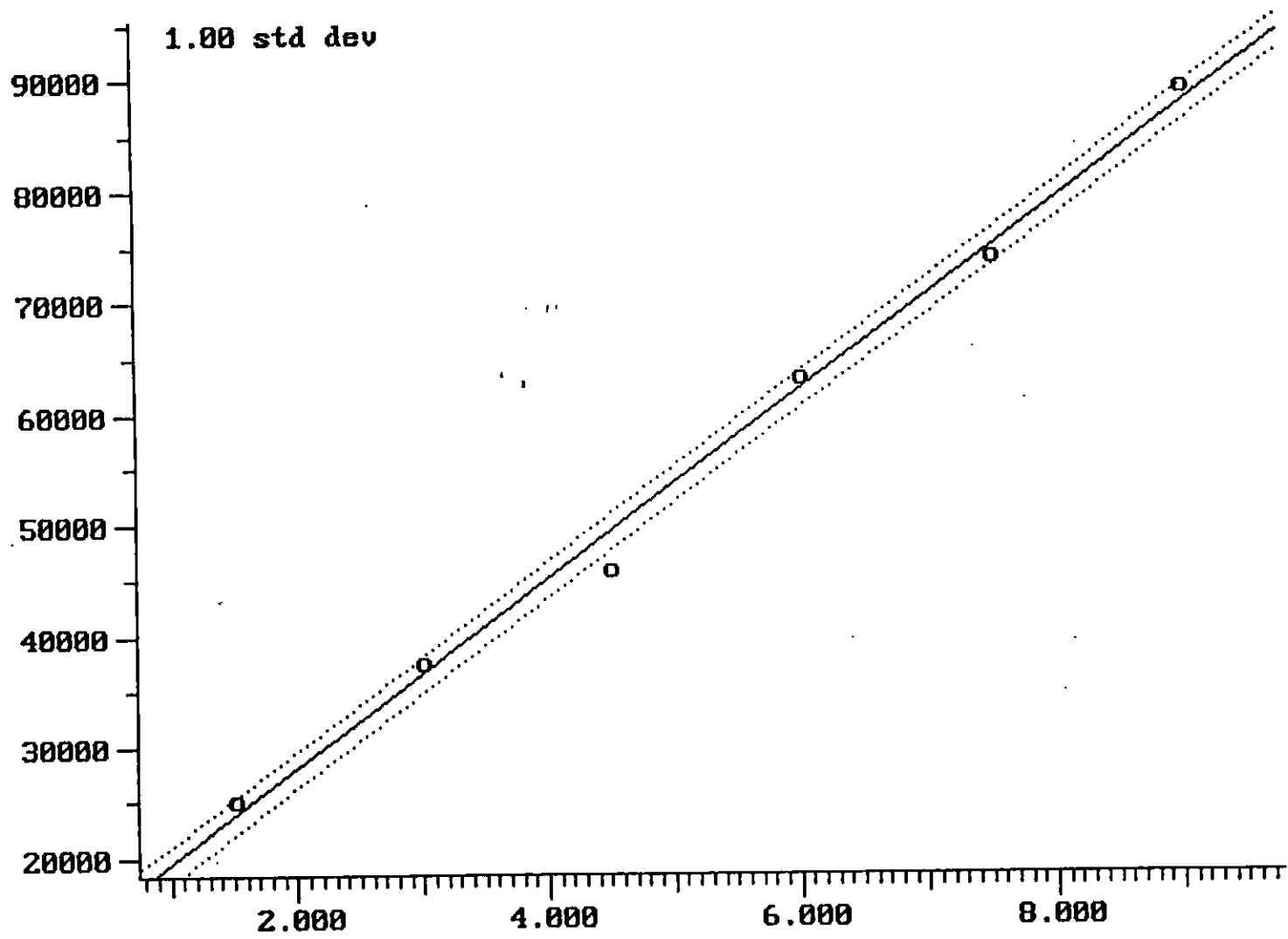
Calibration Plot (Ext Stds) Filename: T014GB
4-Ethyltoluene Compound: 37 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.998
Standard Deviation: 1.225
(Lin:Lin)



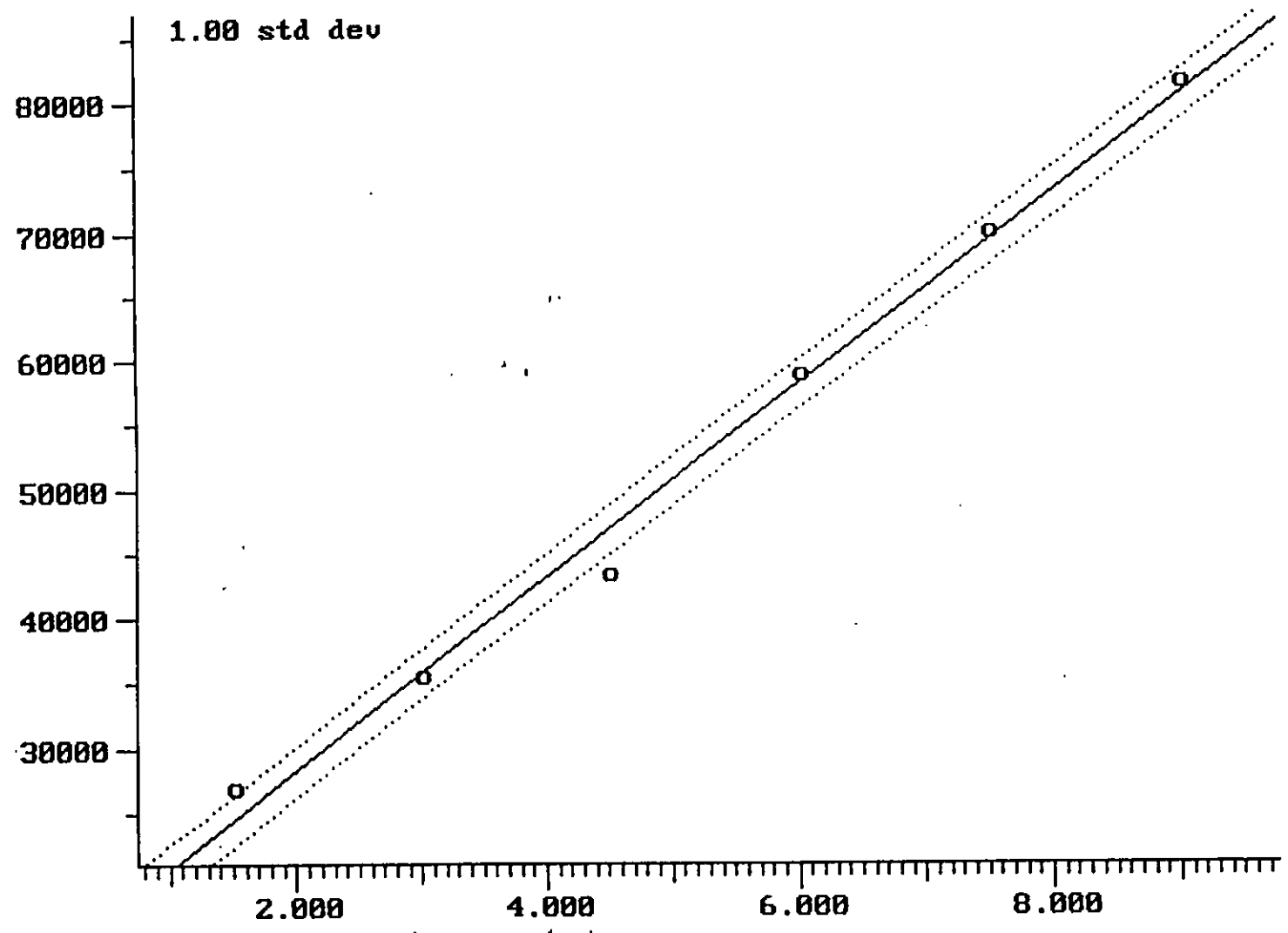
Calibration Plot (Ext Stds) Filename: T014GB
1,3,5-Trimethylbenzene Compound: 38 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

Correlation Coeff: 0.997
Standard Deviation: 1.725
(Lin/Lin)

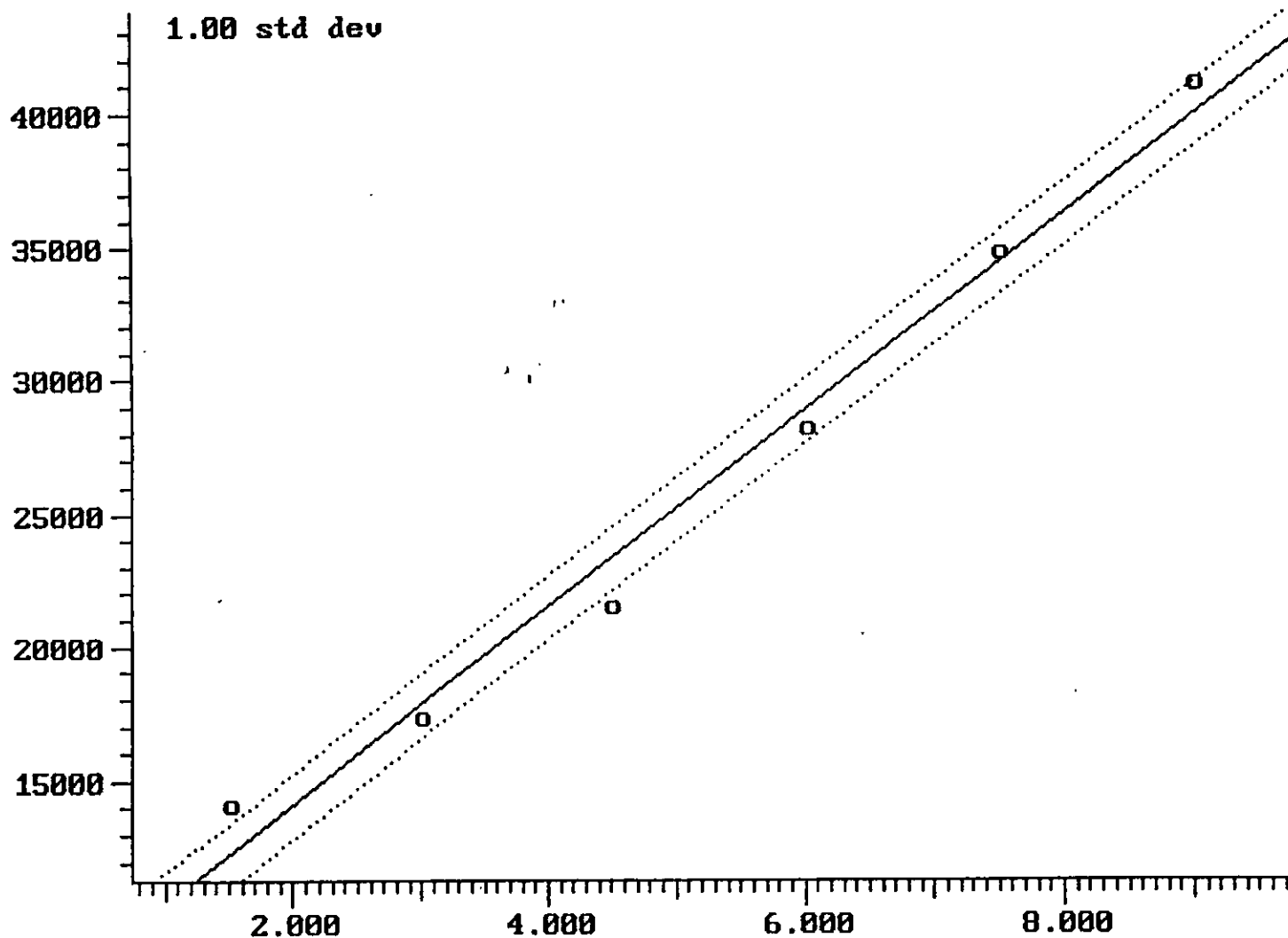


Calibration Plot (Ext Stds) Filename: T014GB
1,2,4-Trimethylbenzene Compound: 39 of 45
(Peak Area of Sample) vs (Amount of Sample Injected)

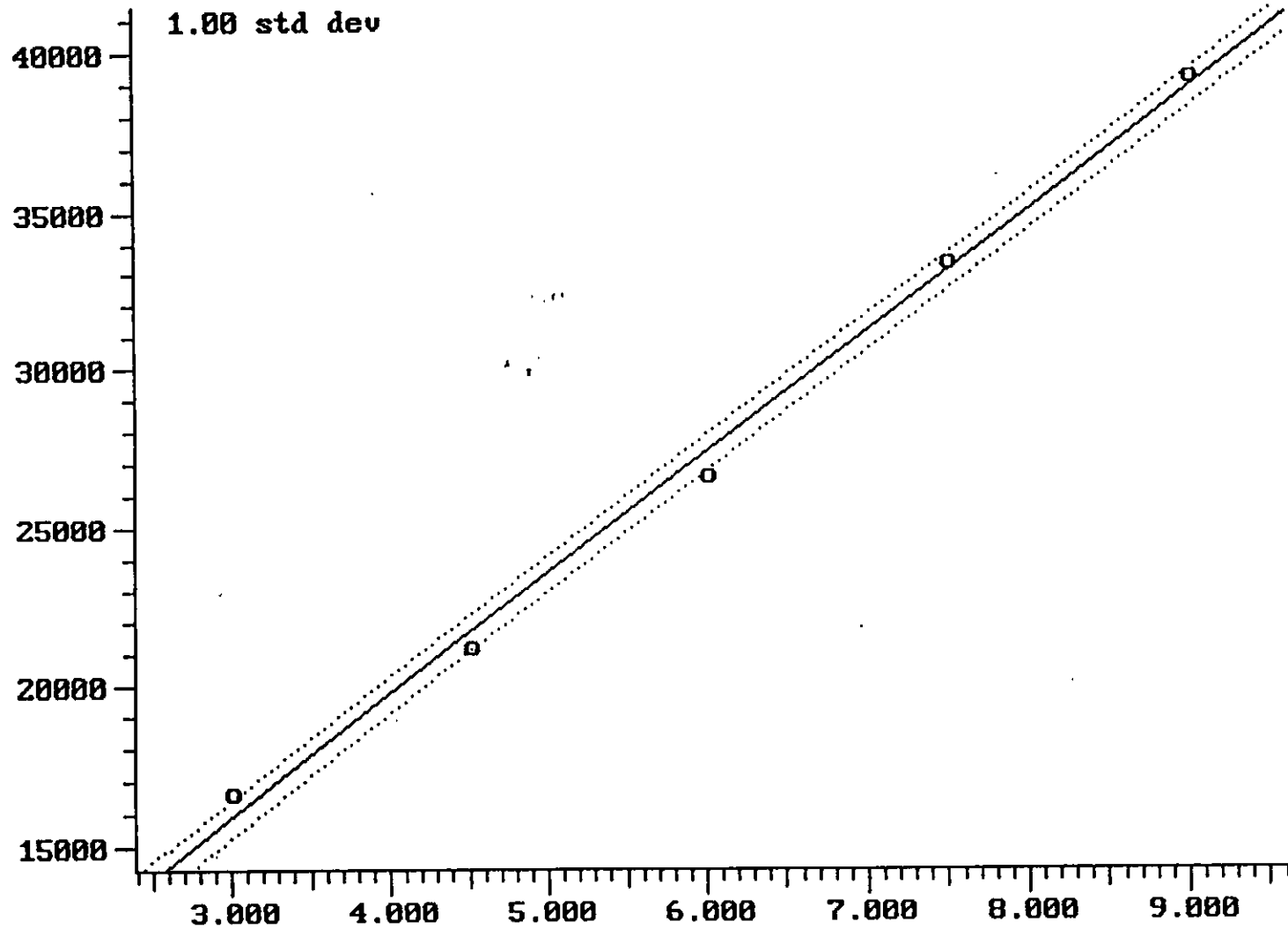
Correlation Coeff: 0.995
Standard Deviation: 1.830
(Lin/Lin)



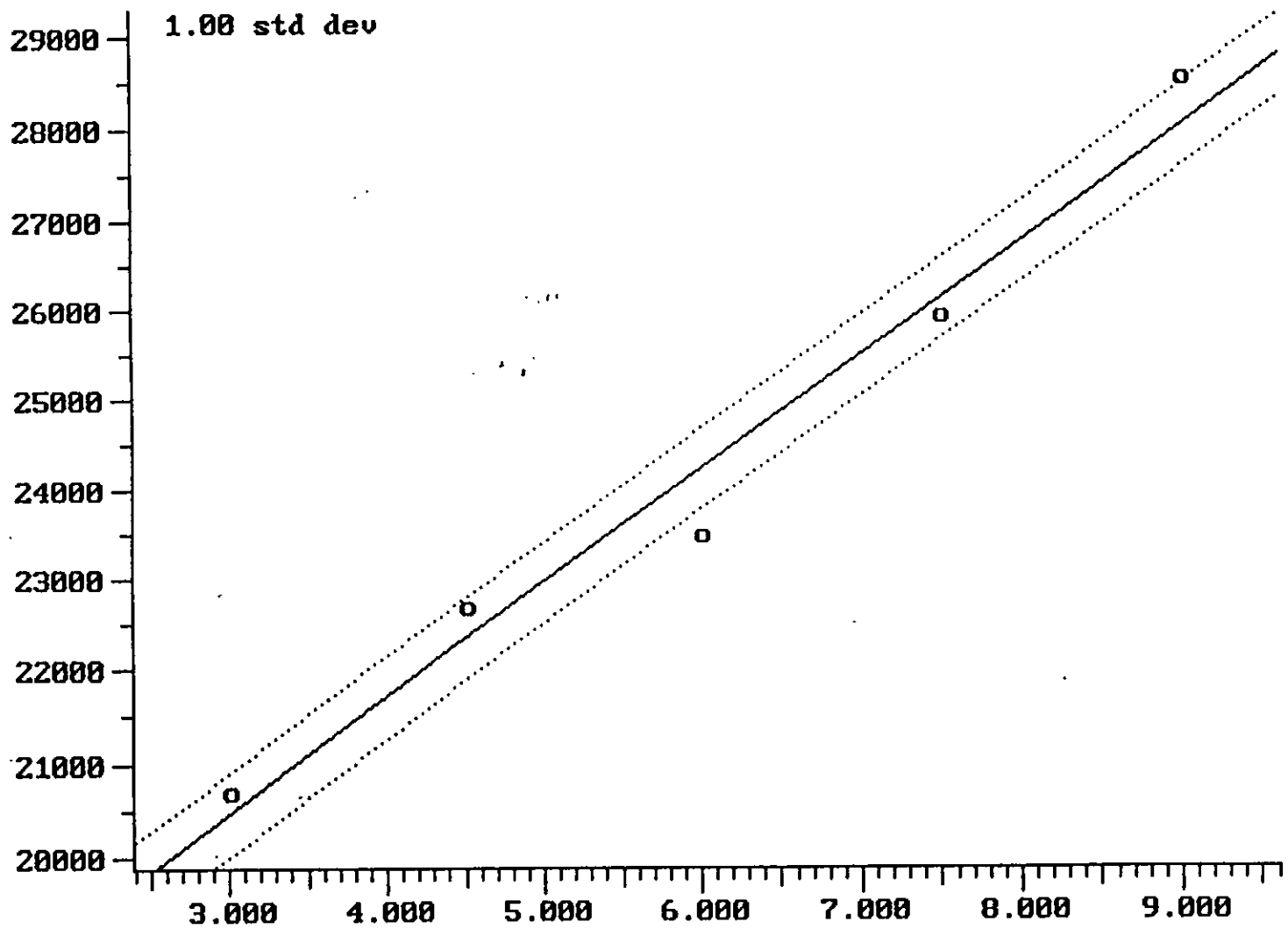
Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.992
1,3-Dichlorobenzene Compound: 40 of 45 Standard Deviation: 1.216
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin!Lin)



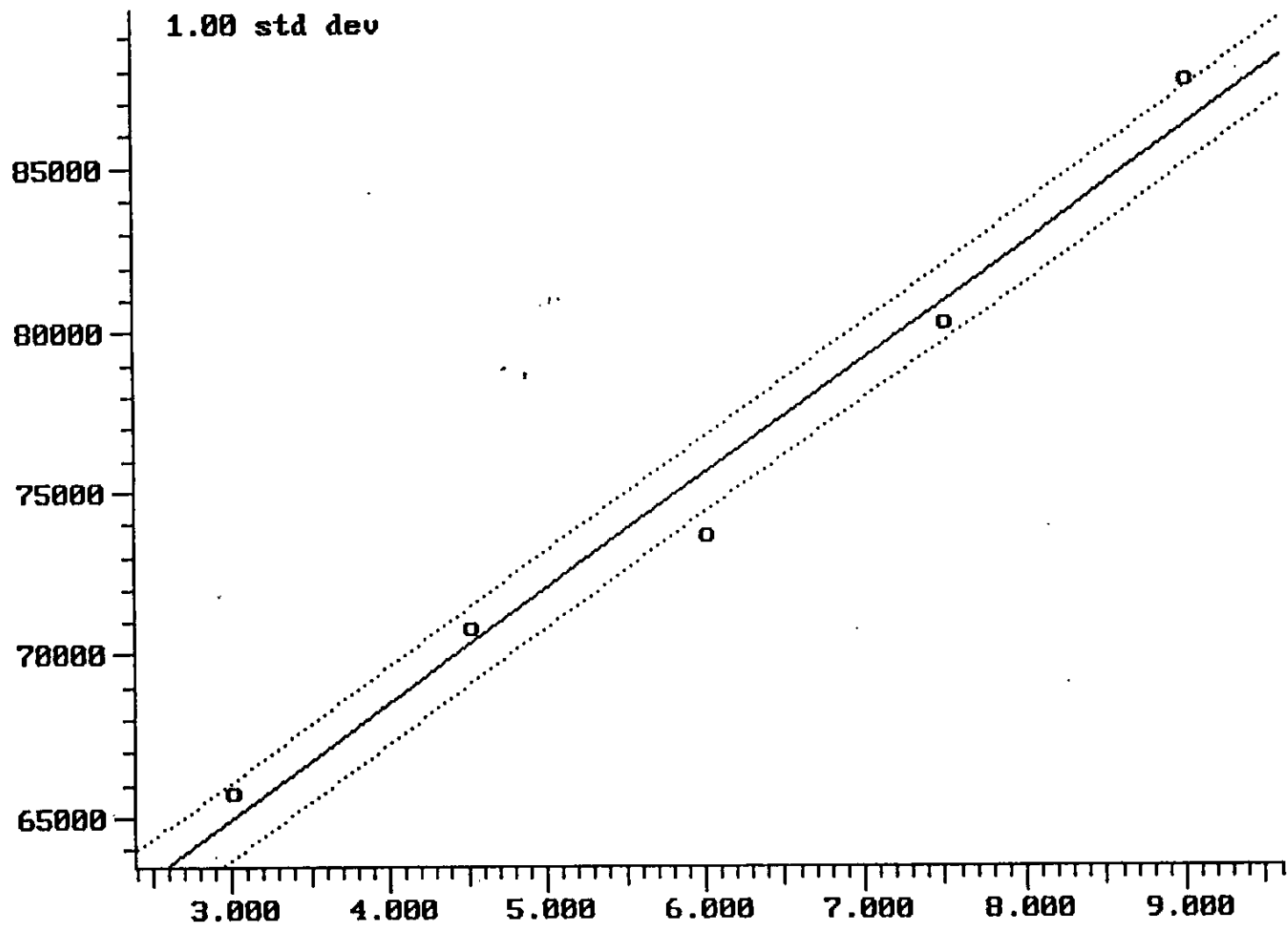
Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.998
1,2-Dichlorobenzene Compound: 42 of 45 Standard Deviation: 0.546
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)



Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.986
1,2,4-Trichlorobenzene Compound: 43 of 45 Standard Deviation: 0.458
(Peak Area of Sample) vs (Amount of Sample Injected) (LinLin)



Calibration Plot (Ext Stds) Filename: T014GB Correlation Coeff: 0.987
Naphthalene Compound: 44 of 45 Standard Deviation: 1.213
(Peak Area of Sample) vs (Amount of Sample Injected) (Lin/Lin)



SECTION 8.0
CHAIN OF CUSTODY AND FIELD SAMPLING FORMS

AirRecon®

Chain of Custody Record

5 Johnson Drive, P.O. Box 130
 Raritan, NJ 08869
 (908) 526-1000 / FAX (908) 526-7886

AirRecon Job #: 311-00142-00-000

City/State: _____

Project Manager: JSG ext. _____

P. O. Number: _____

Samples Submitted to: AirRecon Lab
 Other _____

- Container Type
- B. Gas Bag
 - D. Petri Dish
 - E. Method 25 Tank
 - F. Method 25 Trap
 - G. Glass
 - N. NIOSH Tube
 - P. Plastic Bottle
 - S. Summa Canister
 - T. VOST Tube
 - V. VOC Vial
 - X. XAD-2 Trap
 - O. Other _____

- Sample Type
- 1. Liquid
 - 2. Gas
 - 3. Filter
 - 4. Charcoal
 - 5. Resin
 - 6. Other _____

Page 1 of 1

| | DUE | E |
|------------------------------|-----|---|
| Emergency Rush ≤ 5 Bus. Days | | |
| Rush ≤ 10 Business Days | | |
| Normal ≤ 15 Business Days | | |

Requested Analysis / Method

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|---------------------------------------|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|
| X m19 list of 42 compounds (provided) | | | | | | | | | | | | | | | | | | | | | | |
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| Sample Identification | Filter # | Sample Type | Container | | Sampling Information | | Reagents and/or Preservatives | Lab ID # |
|-----------------------|----------|-------------|-----------|------|----------------------|------|-------------------------------|----------|
| | | | Type | Size | Date | Tech | | |
| Bag #1 | | 2 | B | 90L | 8/14/00 | JSG | | |
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Special Instructions: _____

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|----------------------------------|----------------------|-----------------------------|----------------------|
| Submitted By: <u>John Sumner</u> | Date: <u>8/14/00</u> | Received By: _____ | Date: _____ |
| Relinquished By: _____ | Date: _____ | Received By: _____ | Date: _____ |
| Relinquished By: _____ | Date: _____ | Received By: <u>Thom...</u> | Date: <u>8/15/00</u> |

QA/QC Report Package

Compliance NJDEP

Non-compliance Method Normal