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Washington State
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Report

Phase II Remedial Investigation Paxton Sales Corporation Yakima, Washington

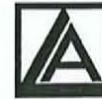
May 25, 1995

Prepared for

**Velikanje, Moore & Shore Inc. P.S.
Yakima, Washington**



Prepared by



LANDAU ASSOCIATES, INC.

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Washington State
Department of Ecology.

Mr. Mark Fickes
Velikanje, Moore & Shore, Inc. P.S.
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P. O. Box C2550
Yakima, WA 98907

**RE: STATUS REPORT
PAXTON SALES CORPORATION REMEDIAL INVESTIGATION
YAKIMA, WASHINGTON**

INTRODUCTION

This status report describes Phase II Remedial Investigation (RI) services completed at the Paxton Sales Corporation (Paxton) site between December, 1994 and April, 1995. Paxton is located at 108 West Mead, as shown on Figure 1.

The first phase of the RI is reported in our site history report (Landau Associates 1994a). Landau Associates completed Phase II field work in accordance to the work plan submitted to Velikanje, Moore & Shore, Inc. on December 27, 1994 (Landau Associates 1994b) with the following exceptions:

- One-inch brass rings were used to line the inside of the Dames & Moore soil sampler while collecting soil samples from MW-2. Analytical results suggest that the brass rings did not interfere with or elevate results of the priority pollutant metal analysis.
- A soil duplicate for all requested analyses was proposed in the work plan; however, due to a lack of soil recovery in the sampling attempts, only a duplicate for volatile organic compounds (VOC) analysis was collected from MW-2.
- The upgradient well (originally designated as MW-1) was deleted from the work plan based on low field screening results of cross- and assumed downgradient wells.
- Paxton Sales' personnel are currently measuring groundwater levels in the monitoring wells on a bi-weekly basis and Landau Associates is processing the data to determine groundwater flow and gradients.

TRIM SOL CUTTING SOLUTION

Paxton has used a cutting solution under the brand name "TrimSol" since it began operations in 1966, with the exception of a one- to two-year period in the mid-1980s when diethanolamine was used. The primary components of TrimSol are described in: 1) a Preliminary Assessment Report prepared by Ecology and Environment, Inc. (Ecology and Environment, 1988); 2) MSDS sheet and other information provided by the manufacturer, Master Chemical Corporation, and 3) laboratory analysis performed by Analytical Resources Inc., (ARI) in November, 1994. The manufacturers information and analytical results were provided to Velikanje, Moore & Shore, Inc. and the Washington State Department of Ecology (Ecology) in correspondence dated December 14, 1994.

According to the manufacturer, TrimSol is a chlorinated alkene polymer that contains no chlorinated solvents (e.g. tetrachloroethene [PCE]). The analytical work performed by ARI supports that statement, as no PCE was detected in the analyzed sample of TrimSol (at a 2 mg/L detection unit). Based on the above, it does not appear that TrimSol could be a source of PCE to either site soil or groundwater.

SOIL SAMPLING, SITE GEOLOGY AND HYDROGEOLOGY

Site soil conditions were explored by drilling three borings, during the period of January 3 to 15, 1995, using a cable-tool drill rig. All three soil borings were converted to monitoring wells upon completion of soil sampling (borings/wells identified as MW-1, MW-2 and MW-3). The borings were drilled approximately 10 ft into the saturated zone at the time of drilling. The locations of the monitoring wells are shown on Figure 2, which also shows other general site features.

Soil samples were attempted every 2.5 ft during the drilling process to identify site geology and for chemical analysis. Native soil consists of very dense gravel with varying amounts of sand and silt. Logs for the borings/wells are presented in Appendix A.

Samples were field screened for initial assessment of potential compounds of concern based on observations and photoionization detector (PID) readings. No unusual odors, discoloration or PID readings were noted during drilling/sampling. One soil sample was selected from each well for chemical analysis based on the results of field screening. A duplicate volatile organic compound (VOC) sample was collected from MW-2. Soil samples selected for chemical analysis were placed in laboratory supplied containers and labeled appropriately.

Soil samples were analyzed for total petroleum hydrocarbons [TPH] (WTPH-HCID); VOCs (EPA Method 8260, [only Method 8010 listed compounds reported]); base/neutral/acids and semi-volatile organic compounds (EPA Method 8270); cyanide (Method 335.2); and priority pollutant metals (EPA Method 6010/7000).

During drilling, groundwater levels were 20-21 ft below the surface. Water levels later stabilized at about 18-20 ft below the surface. Soil to approximately 10 ft above the static water level showed evidence of seasonal groundwater fluctuations.

GROUNDWATER SAMPLING

The monitoring wells were installed and developed according to WAC 173-160. The initial round of groundwater sampling was completed on January 23, 1995 (during lower winter-time groundwater levels). A duplicate sample was collected from MW-1. Groundwater samples were analyzed for VOCs (EPA Method 8260 [8010 list reported]); diesel-range total petroleum hydrocarbons (WTPH-D); priority pollutant metals (EPA Methods 6010/7000); base/neutral/acids and semivolatile organic compounds (EPA Method 8270); and cyanide (Method 335.2). No unusual odors or discoloration were noted during well purging and sampling.

GROUNDWATER ELEVATION DATA

Monitoring of site ground water levels occurred about weekly from January 23, 1995 through March 1, 1995, and is currently ongoing on about a bi-weekly schedule. Readings are being taken by measuring the depth to groundwater with an electrical sounding tape. The depth measurement is read from a surveyed point on the PVC well casing; groundwater elevations are calculated based on a site datum of elevation 100.00 ft. A summary of groundwater elevation data is presented in Table 1.

From the groundwater information obtained, Landau Associates has determined that net groundwater flow is to the east at an average gradient of about 0.016 ft per ft. Groundwater elevations and flow direction from the February 27, 1995 readings are shown on Figure 3. Higher, summertime groundwater levels from May 22, 1995 are shown on Figure 4. Figures 5, 6 and 7 present plots of groundwater elevation data. These data are generally consistent with regional information collected by Ecology in 1993 (Ecology 1993a), although the flow direction at Paxton shows a stronger easterly component than noted by Ecology.

ANALYTICAL PROGRAM

Analytical Resources, Inc. (ARI), based in Seattle, Washington, performed the soil and groundwater sample chemical analyses. ARI maintains a comprehensive internal quality assurance program. Landau Associates also performed a focused data validation to provide a reasonable degree of confidence in project data. The focused data validation was completed in compliance with the *National Functional Guidelines for Organic Data Review* (EPA 1991). A February 27, 1995 Landau Associates technical memorandum provides the results of the data validation, and is included as Appendix B.

SOIL QUALITY

Summaries of chemical analyses results for soil samples are presented in Tables 2 through 6; the laboratory reports are included in Appendix C. The results indicate that VOCs, petroleum hydrocarbons, semivolatile organic compounds, and cyanide were not detected in site soil. The metals results indicate that arsenic, beryllium, chromium, copper, lead, nickel, selenium, thallium, and zinc are present in site soil, but occur at typical background concentrations. The pattern of occurrence does not suggest that the dry well at Paxton is a source of metals to site soil.

GROUNDWATER QUALITY

Summaries of groundwater chemical analyses are presented in Table 7; the laboratory reports are included in Appendix C. Semivolatile organic compounds and diesel range hydrocarbons were not detected in site groundwater. VOC analyses indicate that PCE was present in all groundwater samples at the following concentrations: MW-1 = 2.3 $\mu\text{g}/\text{L}$; MW-1A, (duplicate) = 2.2 $\mu\text{g}/\text{L}$; MW-2 = 1.7 $\mu\text{g}/\text{L}$; and MW-3 = 3.1 $\mu\text{g}/\text{L}$. These values are below the MCL and Method A cleanup level of 5 $\mu\text{g}/\text{L}$ and the 4 $\mu\text{g}/\text{L}$ advisory level established by Ecology. The current pattern of occurrence does not suggest that the dry well is the source of PCE to site groundwater.

Cyanide was detected in site groundwater at MW-1; however, the data validation qualified the cyanide data because the relative percent difference between the sample and field duplicate exceeded the established control limits. The reported laboratory values for cyanide in groundwater at MW-1 are well below published MTCA Method B values (CLARC II, January '95).

The metals results indicate that arsenic, beryllium, chromium, copper, lead, mercury, nickel, selenium, thallium, and zinc are present in site groundwater. The laboratory data for arsenic has been rejected because the field duplicate results were outside of control limits. However, arsenic is commonly found in shallow groundwater throughout the Yakima area due to the past application of arsenic containing pesticides. Accordingly, the possible presence of arsenic in site groundwater is not currently viewed as cause for additional study at this time (since Paxton is not a known user [source] of arsenic).

Beryllium (as a non-carcinogen), chromium, copper, mercury, nickel, selenium and zinc were either not detected, or reported at levels below MTCA Method B values.

Lead was found in site groundwater in all 3 wells at concentrations above the MTCA Method A cleanup level. Lead in shallow groundwater is also common in the Yakima area, again primarily due to past use of pesticides. Paxton reports no past use of lead, and the pattern of occurrence does not suggest the dry well as a source. Accordingly, the presence of lead in shallow groundwater is not currently viewed as cause for additional study at this time.

Thallium was found in site groundwater at MW-1 and MW-2 at concentrations slightly above the MTCA Method B value. Levels are low and there is no known local man-caused source of thallium; therefore, the presence of thallium in shallow groundwater is not currently viewed as cause for additional study at this time.

CMX CORPORATION DATA

CMX Corporation (CMX), located at 206 West Mead, is west and generally up- or cross-gradient to Paxton. In November 1992, Ecology installed a monitoring well at CMX, collected one groundwater sample for analysis, and reported the results in February 1993 (Ecology 1993). The text of that report states "groundwater samples non-detect" (for PCE); however, Ecology's data summary table and laboratory report for that sample show a PCE concentration of 1.8 $\mu\text{g}/\text{L}$. Assuming the laboratory report is correct, it appears that low levels of PCE are present in shallow groundwater in the up-gradient direction to Paxton.

FINDINGS AND RECOMMENDATIONS

The following findings and recommendations are based on Phase II RI data, the earlier Phase I history report and other relevant documents:

- Geologic conditions at Paxton are consistent with regional findings.

- The depth to groundwater, flow direction and gradient at Paxton are generally consistent with regional finding; however, flow direction at the site appears to be in a more easterly direction than shown in Ecology, 1993. As a result, MW-2, which was located in the assumed down-gradient direction, appears to be in a cross-and downgradient direction to the Paxton drywell.
- The TrimSol cutting solution, earlier suspected of being a source of PCE, contains no PCE.
- Field observations and laboratory analytical results do not suggest that site soil conditions have been impacted by Paxton's operations.
- Groundwater quality data on samples taken during winter-time "lows" (January 1995) indicate that site groundwater contains concentrations of PCE below the 5 $\mu\text{g}/\text{L}$ MCL and MTCA Method A cleanup level and the 4 $\mu\text{g}/\text{L}$ Ecology advisory level. The site values (1.7 $\mu\text{g}/\text{L}$ to 3.1 $\mu\text{g}/\text{L}$) are in line with the 1.8 $\mu\text{g}/\text{L}$ 1992 value from the CMX site, suggesting probable background levels. The pattern of occurrence of PCE in groundwater at Paxton does not suggest that the dry well is a source.
- The occurrence of cyanide in site groundwater shows a possible link between site operations and groundwater in the monitoring well nearest the dry well (MW-1). However, the qualified value reported by the laboratory is well below the MTCA Method B cleanup level for cyanide (as a non-carcinogen).
- The occurrence of lead, arsenic and other metals in site groundwater is consistent with background levels, or that which could be attributed to past agricultural practices. The pattern of occurrence does not suggest the Paxton dry well as a source.
- Since one round of groundwater samples is typically not adequate for drawing final conclusions, Landau Associates recommends additional groundwater sampling. The next sampling should be performed after water levels rise in the spring (e.g. June '95) so that "high" and "low" groundwater level samples can be evaluated.
- Once the next groundwater sample results have been obtained, the need for future site activities, if any, can be determined.

LANDAU ASSOCIATES, INC.

By:



William D. Evans, CPG
Associate

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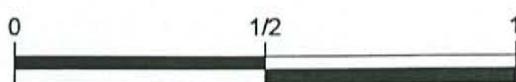
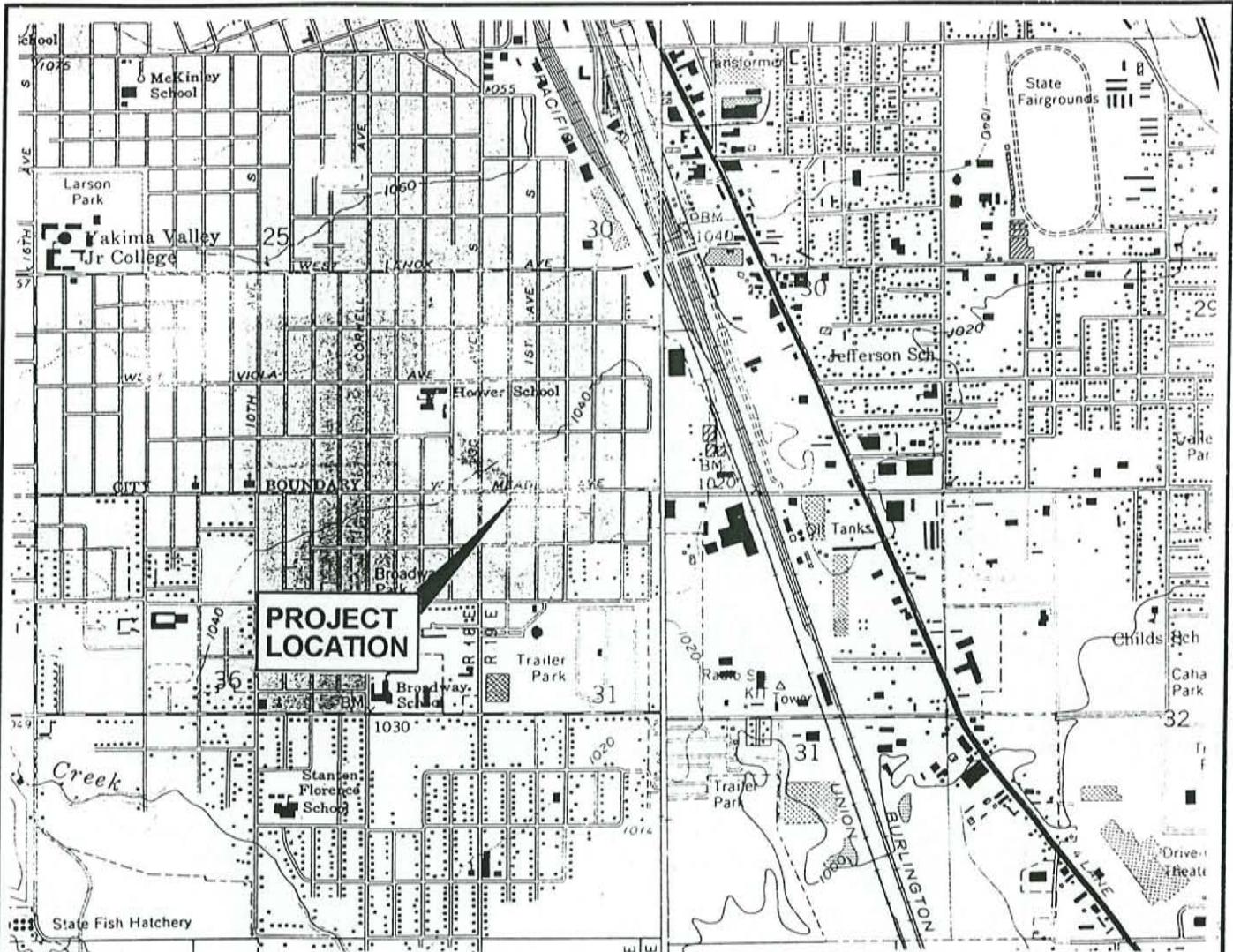
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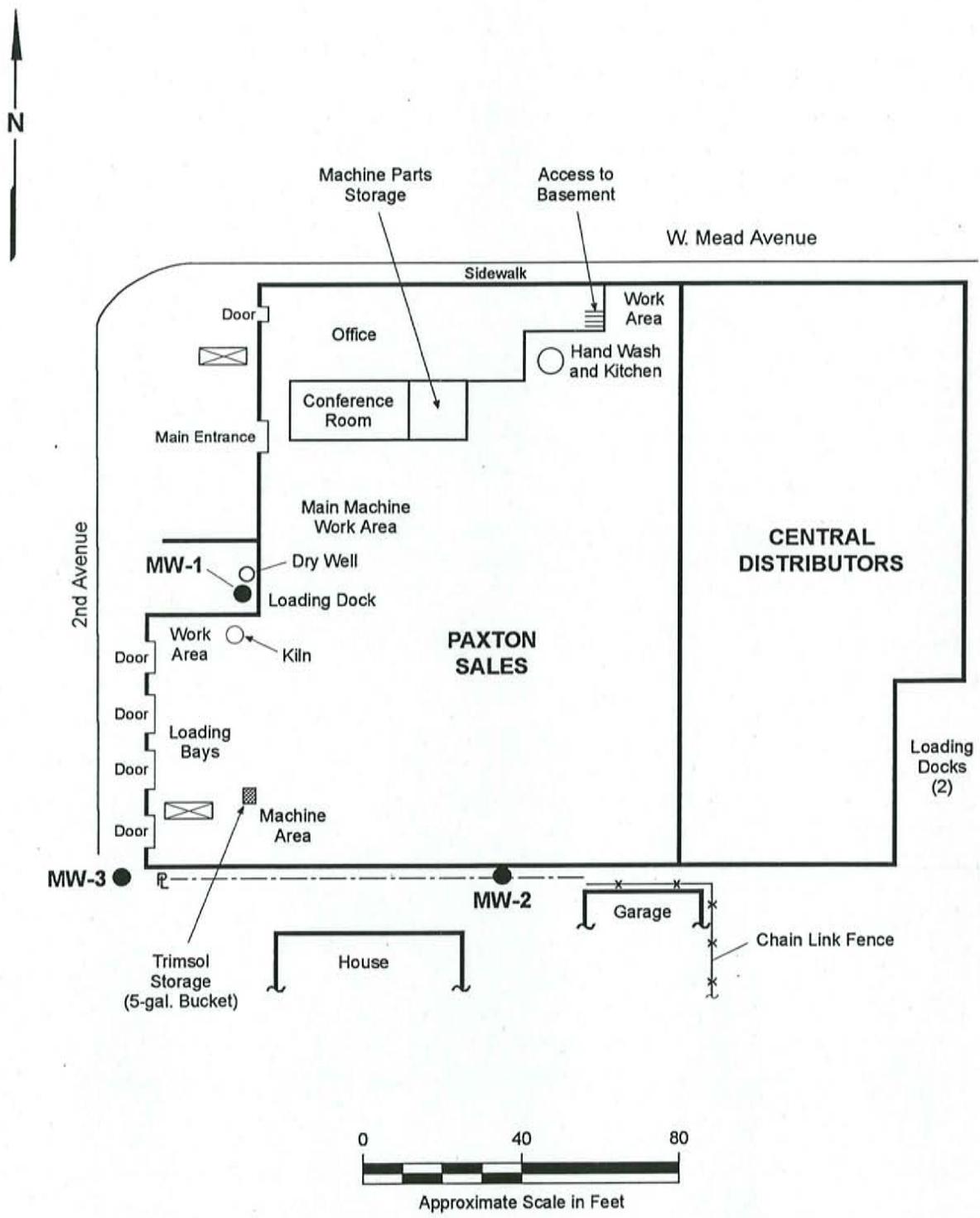
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N



**KEY**

MW-1 Monitoring Well Location and Identification



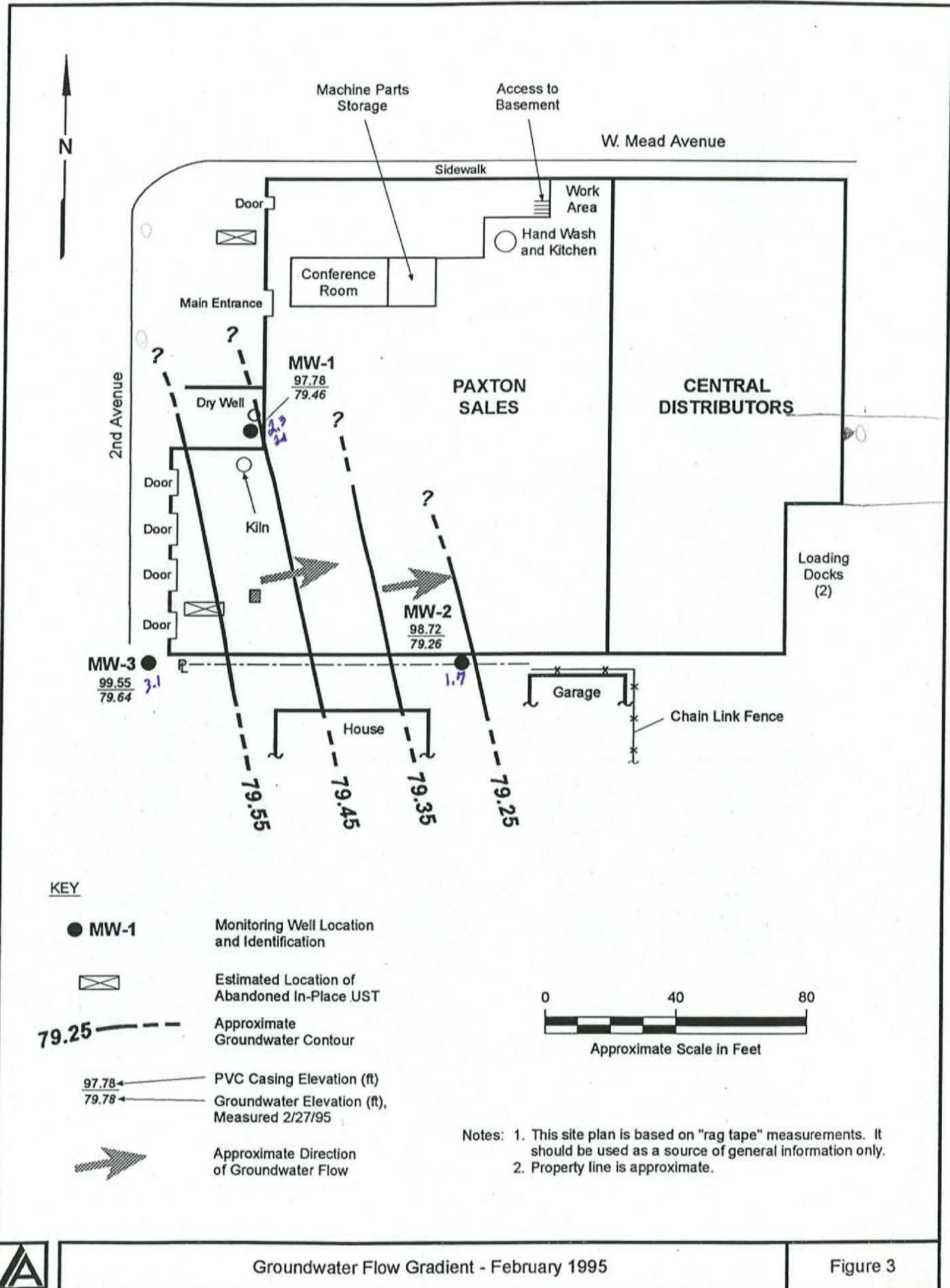
Estimated Location of Abandoned In-Place UST

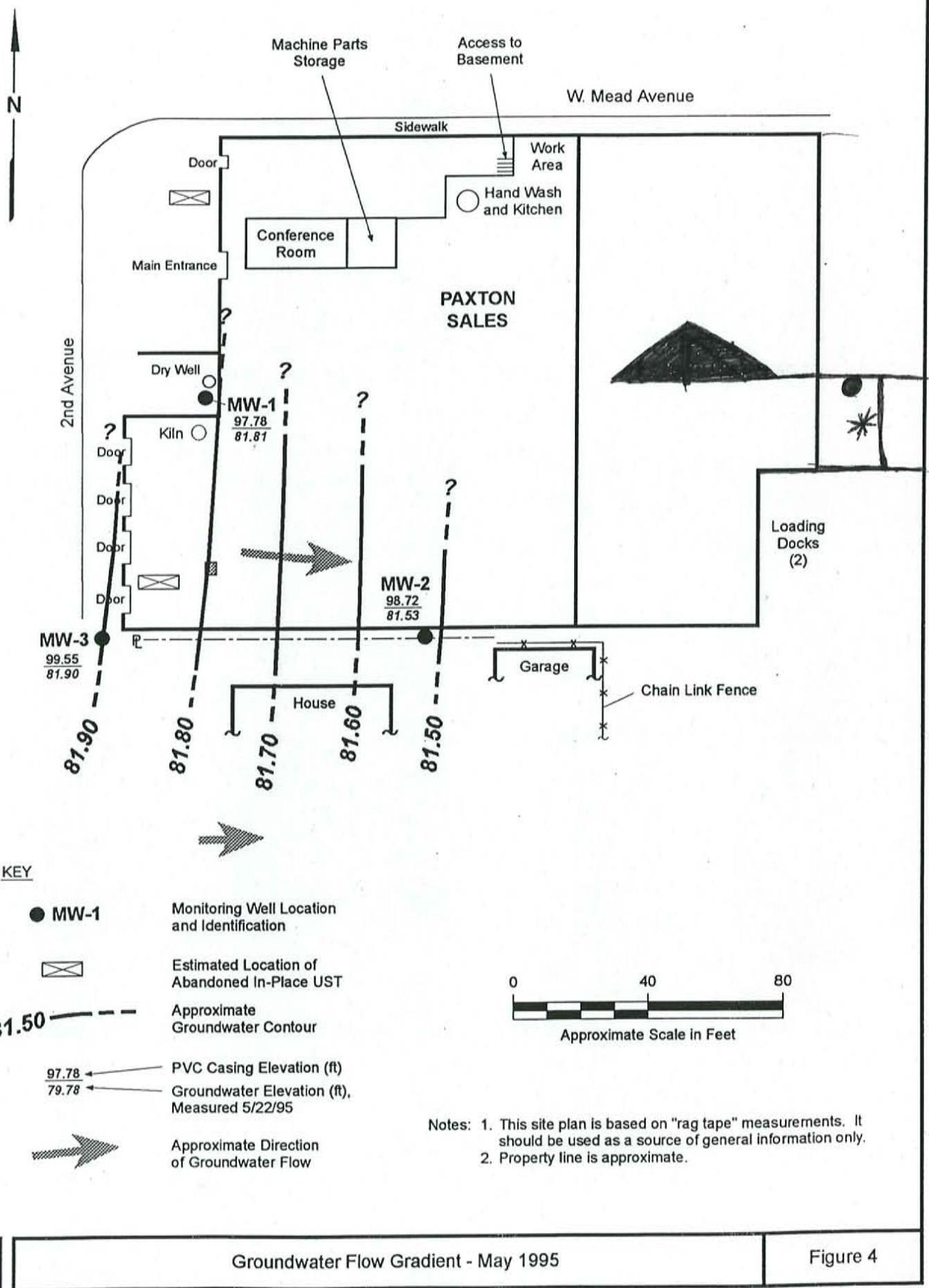
Notes: 1. This site plan is based on "rag tape" measurements. It should be used as a source of general information only.
 2. Property line is approximate.

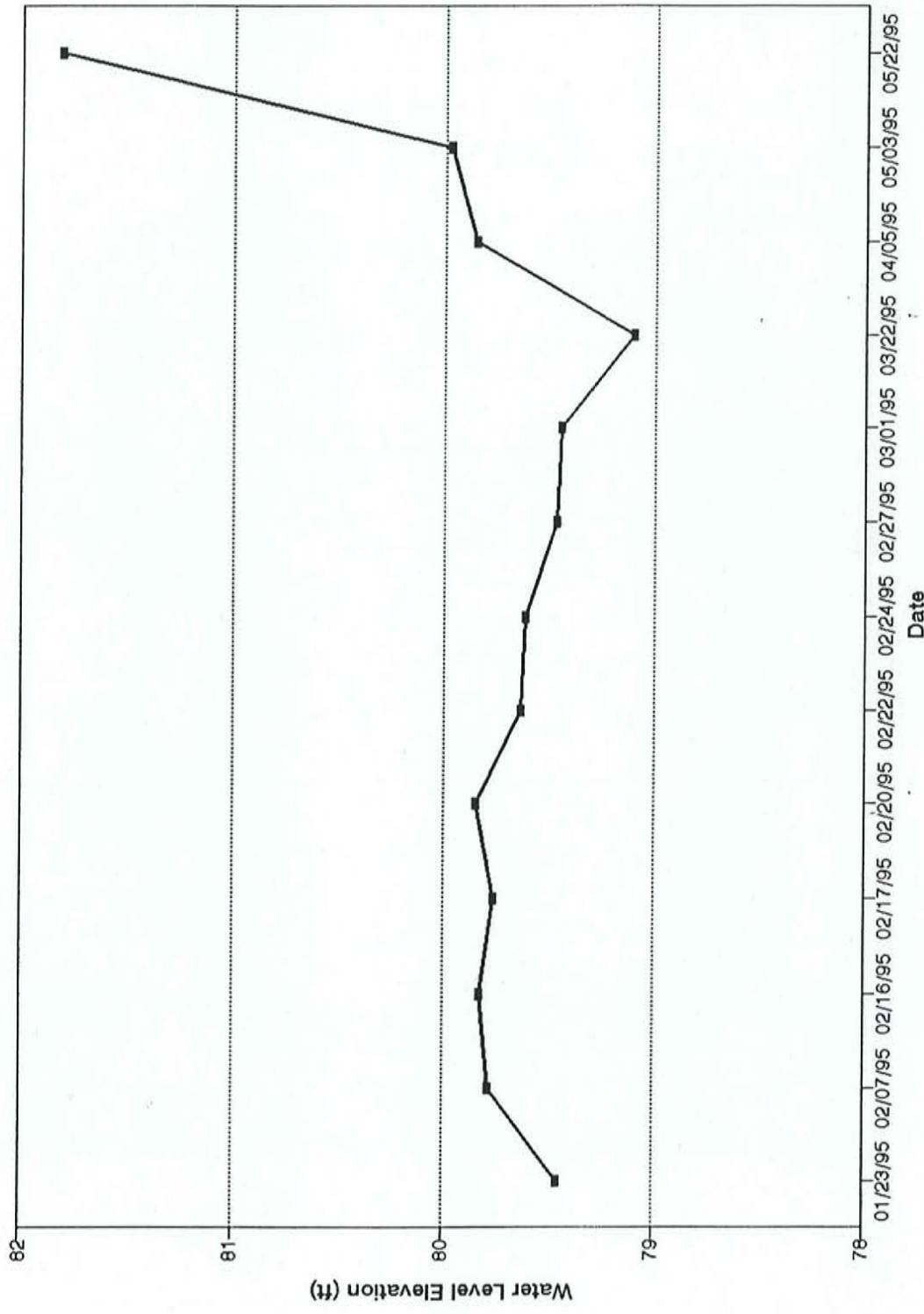


Site Plan Showing Monitoring Well Locations

Figure 2



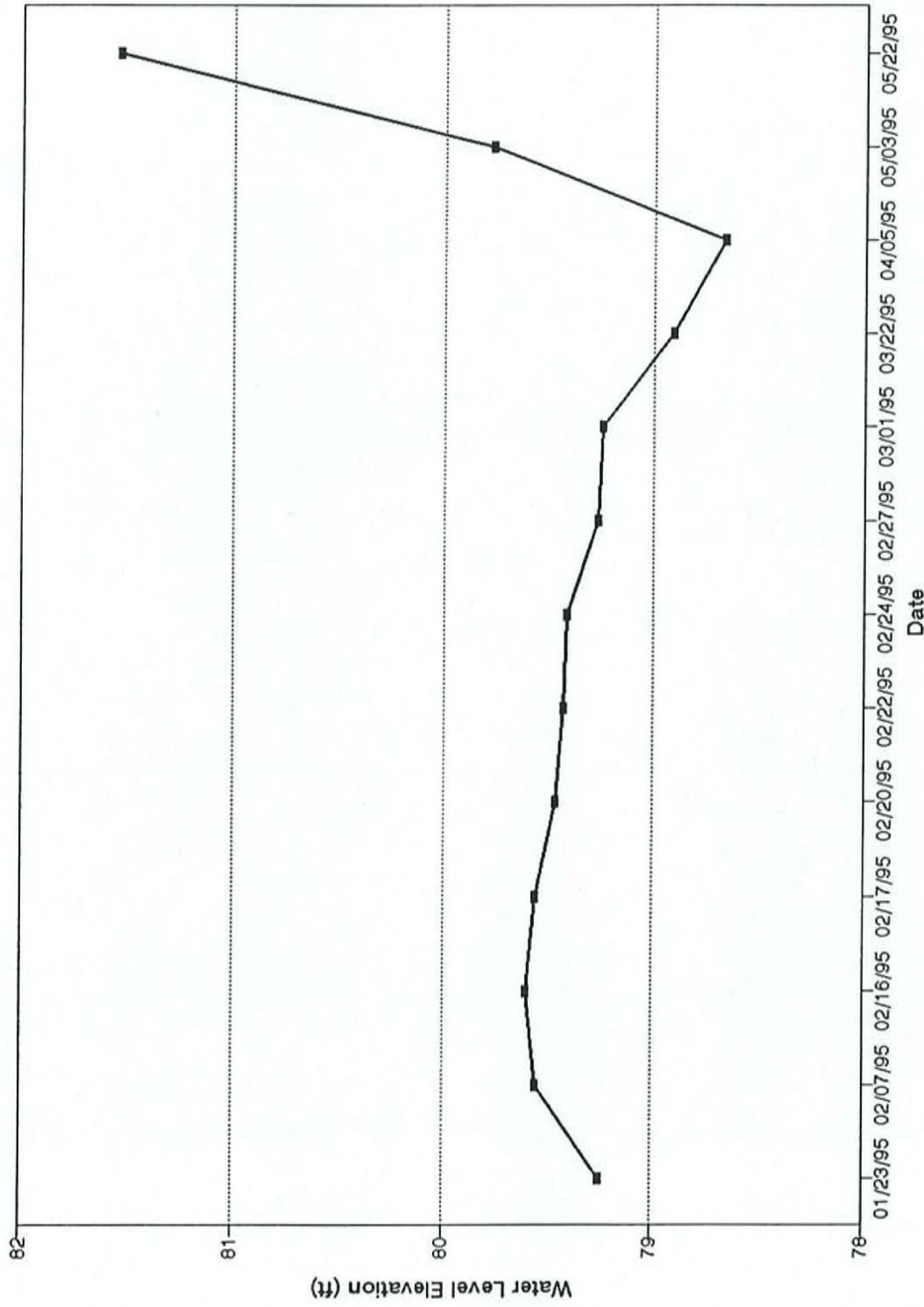




A

Summary of Groundwater Elevations
MW-1

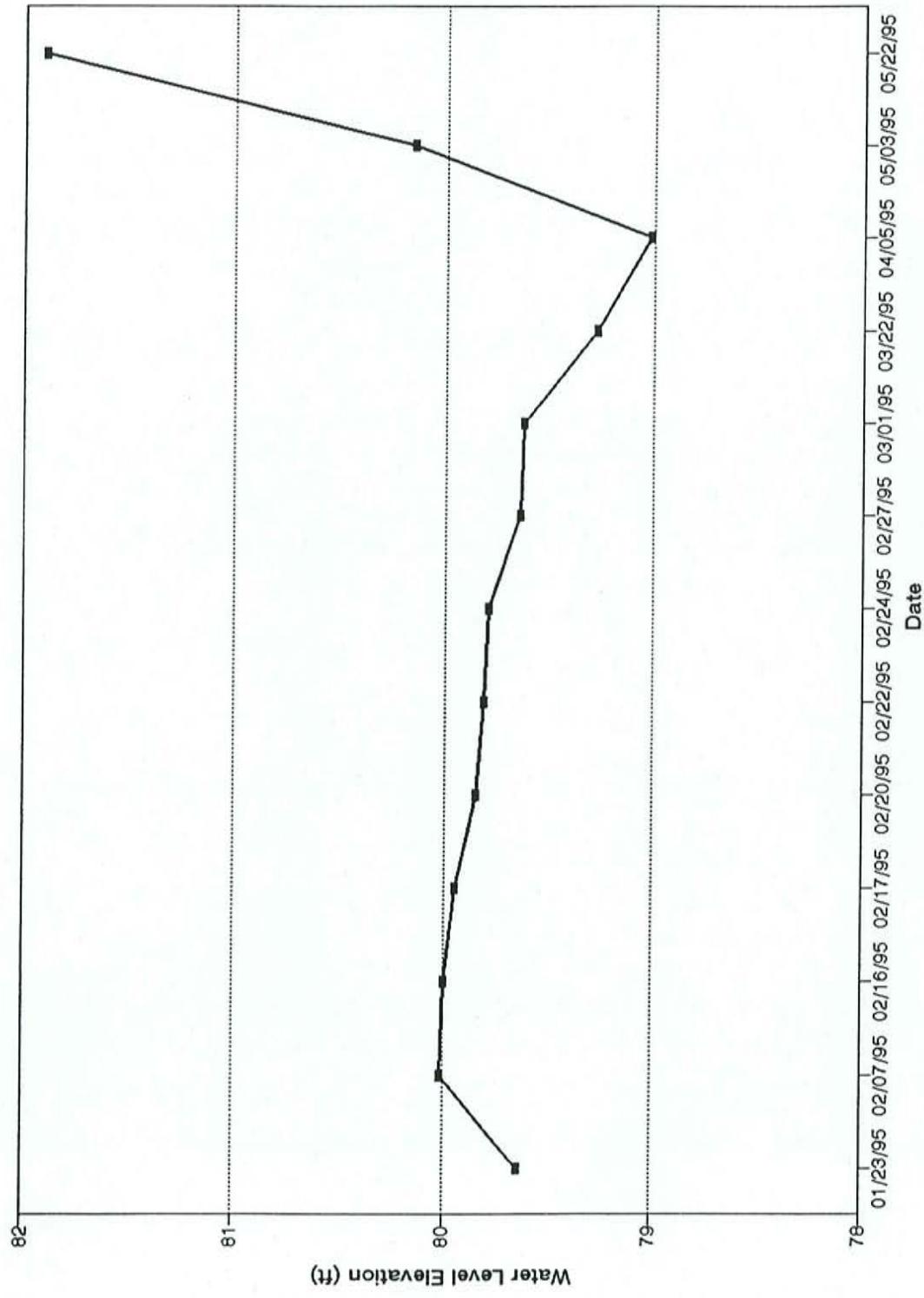
Figure 5



A

Summary of Groundwater Elevations
MW-2

Figure 6



A

Summary of Groundwater Elevations
MW-3

Figure 7

TABLE 1
SUMMARY OF GROUNDWATER ELEVATIONS
PAXTON SALES

| Well | Measurement Date | Depth to Water | Well Casing Elevation | Water Elevation |
|------|------------------|----------------|-----------------------|-----------------|
| MW-1 | 01/23/95 | 18.33 | 97.78 | 79.45 |
| | 02/07/95 | 18.00 | 97.78 | 79.78 |
| | 02/16/95 | 17.96 | 97.78 | 79.82 |
| | 02/17/95 | 18.02 | 97.78 | 79.76 |
| | 02/20/95 | 18.13 | 97.78 | 79.84 |
| | 02/22/95 | 18.15 | 97.78 | 79.63 |
| | 02/24/95 | 18.17 | 97.78 | 79.61 |
| | 02/27/95 | 18.32 | 97.78 | 79.46 |
| | 03/01/95 | 18.34 | 97.78 | 79.44 |
| | 03/22/95 | 18.68 | 97.78 | 79.10 |
| | 04/05/95 | 18.93 | 97.78 | 79.85 |
| | 05/03/95 | 17.81 | 97.78 | 79.97 |
| MW-2 | 05/22/95 | 15.97 | 97.78 | 81.81 |
| | 01/23/95 | 19.47 | 98.72 | 79.25 |
| | 02/07/95 | 19.17 | 98.72 | 79.55 |
| | 02/16/95 | 19.12 | 98.72 | 79.60 |
| | 02/17/95 | 19.16 | 98.72 | 79.56 |
| | 02/20/95 | 19.26 | 98.72 | 79.46 |
| | 02/22/95 | 19.30 | 98.72 | 79.42 |
| | 02/24/95 | 19.31 | 98.72 | 79.41 |
| | 02/27/95 | 19.46 | 98.72 | 79.26 |
| | 03/01/95 | 19.48 | 98.72 | 79.24 |
| | 03/22/95 | 19.82 | 98.72 | 78.90 |
| | 04/05/95 | 20.06 | 98.72 | 78.66 |
| MW-3 | 05/03/95 | 18.95 | 98.72 | 79.77 |
| | 05/22/95 | 17.19 | 98.72 | 81.53 |
| | 01/23/95 | 19.91 | 99.55 | 79.64 |
| | 02/07/95 | 19.54 | 99.55 | 80.01 |
| | 02/16/95 | 19.55 | 99.55 | 80.00 |
| | 02/17/95 | 19.61 | 99.55 | 79.94 |
| | 02/20/95 | 19.71 | 99.55 | 79.84 |
| | 02/22/95 | 19.74 | 99.55 | 79.81 |
| | 02/24/95 | 19.76 | 99.55 | 79.79 |
| | 02/27/95 | 19.91 | 99.55 | 79.64 |
| | 03/01/95 | 19.93 | 99.55 | 79.62 |
| | 03/22/95 | 20.28 | 99.55 | 79.27 |
| | 04/05/95 | 20.54 | 99.55 | 79.01 |
| | 05/03/95 | 19.40 | 99.55 | 80.15 |
| | 05/22/95 | 17.65 | 99.55 | 81.90 |

TABLE 2
PAXTON SALES CORPORATION
REMEDIAL INVESTIGATION
SOIL ANALYSIS DATA
VOLATILE ORGANIC COMPOUNDS (ug/kg)

| ANALYTE | MW-1-5.1 | MW-1-7.5 | MW-1-22.3 | MW-1-30.0 | MW-2-20.2 | MW-2-120.2* | MW-3-14.6 |
|---------------------------|----------|----------|-----------|-----------|-----------|-------------|-----------|
| Chromomethane | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.1 U |
| Vinyl Chloride | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.1 U |
| Chloroethane | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.1 U |
| Methylene Chloride | 2.4 U | 2.2 U | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.1 U |
| 1,1-Dichloroethene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| 1,1-Dichloroethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| trans-1,2-Dichloroethene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| Chloroform | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| 1,2-Dichloroethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| 1,1,1-Trichloroethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| Carbon Tetrachloride | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| Bromodichloromethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| 1,2-Dichloropropane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| Trichloroethene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| Dibromochloromethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| 1,1,2-Trichloroethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| trans-1,3-Dichloropropene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| 2-Chloroethylvinylether | 5.4 U | 5.4 U | 5.6 U | 5.2 U | 5.4 U | 5.5 U | 5.2 U |
| Bromoform | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| Tetrachloroethene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| 1,1,2,2-Tetrachloroethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| Chlorobenzene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.1 U | 1.1 U |
| Trichlorofluoromethane | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.2 U | 2.2 U | 2.1 U |
| 1,2-Dichlorobenzene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |
| 1,3-Dichlorobenzene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.0 U | 1.0 U |

TABLE 2
 PAXTON SALES CORPORATION
 REMEDIAL INVESTIGATION
 SOIL ANALYSIS DATA
 VOLATILE ORGANIC COMPOUNDS (ug/kg)

| ANALYTE | MW-1-5.1 | MW-1-7.5 | MW-1-22.3 | MW-1-30.0 | MW-2-20.2 | MW-2-120.2* | MW-3-14.6 |
|---------------------------|----------|----------|-----------|-----------|-----------|-------------|-----------|
| 1,4-Dichlorobenzene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.1 U | 1.0 U |
| Bromoethane | 2.2 U | 2.1 U | 2.2 U | 2.1 U | 2.2 U | 2.2 U | 2.1 U |
| Dibromomethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.1 U | 1.0 U |
| 1,1,1,2-Tetrachloroethane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.1 U | 1.0 U |
| 1,2,3-Trichloropropane | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.1 U | 1.0 U |
| Bromobenzene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.1 U | 1.0 U |
| 2-Chlorotoluene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.1 U | 1.0 U |
| 4-Chlorotoluene | 1.1 U | 1.1 U | 1.1 U | 1.0 U | 1.1 U | 1.1 U | 1.0 U |

Notes:

* = MW-2-120.2 is a duplicate of MW-2-20.2.

U = Analyte not detected at the detection limit reported.

file: paxvoc.wq1
 2-21-95
 mlm

TABLE 3
 PAXTON SALES CORPORATION
 REMEDIAL INVESTIGATION
 SOIL ANALYSIS DATA
 TPH-HCID

| ANALYTE | MW-1-5.2 | MW-1-22.4 | MW-1-25.0 | MW-1-27.5 | MW-2-20.8 | MW-3-15 |
|--------------|----------|-----------|-----------|-----------|-----------|---------|
| Gas Range | (mg/kg) | 20 U | 20 U | 20 U | 20 U | 10 U |
| Diesel Range | (mg/kg) | 25 U | 25 U | 25 U | 25 U | 10 U |
| Oil Range | (mg/kg) | 50 U | 50 U | 50 U | 50 U | 25 U |

Notes:

U = Analyte not detected at detection limit reported.

file: paxph.wq1
 2-21-95
 mlm

TABLE 4
 PAXTON SALES CORPORATION
 REMEDIAL INVESTIGATION
 SOIL ANALYSIS DATA
 SEMIVOLATILE ORGANIC COMPOUNDS (ug/kg)

| ANALYTE | MW-1-17.6 | MW-1-20.2 | MW-2-20.4 | MW-3-20 |
|-------------------------------|-----------|-----------|-----------|---------|
| Phenol | 150 U | 150 U | 150 U | 140 U |
| Bis-(2-chloroethyl) Ether | 150 U | 150 U | 150 U | 140 U |
| 2-Chlorophenol | 74 U | 76 U | 73 U | 70 U |
| 1,3-Dichlorobenzene | 74 U | 76 U | 73 U | 70 U |
| 1,4-Dichlorobenzene | 74 U | 76 U | 73 U | 70 U |
| Benzyl Alcohol | 370 U | 380 U | 370 U | 350 U |
| 1,2-Dichlorobenzene | 74 U | 76 U | 73 U | 70 U |
| 2-Methylphenol | 150 U | 150 U | 150 U | 140 U |
| 2,2'-Oxybis (1-Chloropropane) | 74 U | 76 U | 73 U | 70 U |
| 4-Methylphenol | 74 U | 76 U | 73 U | 70 U |
| N-Nitroso-Di-N-Propylamine | 150 U | 150 U | 150 U | 140 U |
| Hexachloroethane | 150 U | 150 U | 150 U | 140 U |
| Nitrobenzene | 74 U | 76 U | 73 U | 70 U |
| Isophorone | 74 U | 76 U | 73 U | 70 U |
| 2-Nitrophenol | 370 U | 380 U | 370 U | 350 U |
| 2,4-Dimethylphenol | 220 U | 230 U | 220 U | 210 U |
| Benzoic Acid | 740 U | 760 U | 730 U | 700 U |
| bis (2-chloroethoxy) Methane | 74 U | 76 U | 73 U | 70 U |
| 2,4-Dichlorophenol | 220 U | 230 U | 220 U | 210 U |
| 1,2,4-Trichlorobenzene | 74 U | 76 U | 73 U | 70 U |
| Naphthalene | 74 U | 76 U | 73 U | 70 U |
| 4-Chloroaniline | 220 U | 230 U | 220 U | 210 U |
| Hexachlorobutadiene | 150 U | 150 U | 150 U | 140 U |
| 4-Chloro-3-methylphenol | 150 U | 150 U | 150 U | 140 U |
| 2-Methylnaphthalene | 74 U | 76 U | 73 U | 70 U |
| Hexachlorocyclopentadiene | 370 U | 380 U | 370 U | 350 U |
| 2,4,6-Trichlorophenol | 370 U | 380 U | 370 U | 350 U |
| 2,4,5-Trichlorophenol | 370 U | 380 U | 370 U | 350 U |
| 2-Chloronaphthalene | 74 U | 76 U | 73 U | 70 U |
| 2-Nitroaniline | 370 U | 380 U | 370 U | 350 U |
| Dimethylphthalate | 74 U | 76 U | 73 U | 70 U |
| Acenaphthylene | 74 U | 76 U | 73 U | 70 U |
| 3-Nitroaniline | 440 U | 460 U | 440 U | 420 U |
| Acenaphthene | 74 U | 76 U | 73 U | 70 U |
| 2,4-Dinitrophenol | 740 U | 760 U | 730 U | 700 U |
| 4-Nitrophenol | 370 U | 380 U | 370 U | 350 U |
| Debenzofuran | 74 U | 76 U | 73 U | 70 U |
| 2,6-Dinitrotoluene | 370 U | 380 U | 370 U | 350 U |
| 2,4-Dinitrotoluene | 370 U | 380 U | 370 U | 350 U |
| Diethylphthalate | 74 U | 76 U | 73 U | 70 U |

TABLE 4
 PAXTON SALES CORPORATION
 REMEDIAL INVESTIGATION
 SOIL ANALYSIS DATA
 SEMIVOLATILE ORGANIC COMPOUNDS (ug/kg)

| ANALYTE | MW-1-17.6 | MW-1-20.2 | MW-2-20.4 | MW-3-20 |
|------------------------------|-----------|-----------|-----------|---------|
| 4-Chlorophenyl-phenylether | 74 U | 76 U | 73 U | 70 U |
| Fluorene | 74 U | 76 U | 73 U | 70 U |
| 4-Nitroaniline | 370 U | 380 U | 370 U | 350 U |
| 4,6-Dinitro-2-Methylphenol | 740 U | 760 U | 730 U | 700 U |
| N-Nitrosodiphenylamine | 74 U | 76 U | 73 U | 70 U |
| 4-Bromophenyl-phenylether | 74 U | 76 U | 73 U | 70 U |
| Hexachlorobenzene | 74 U | 76 U | 73 U | 70 U |
| Pentachlorophenol | 370 U | 380 U | 370 U | 350 U |
| Phenanthrene | 74 U | 76 U | 73 U | 70 U |
| Carbazole | 74 U | 76 U | 73 U | 70 U |
| Anthracene | 74 U | 76 U | 73 U | 70 U |
| Di-n-Butylphthalate | 74 U | 76 U | 73 U | 70 U |
| Fluoranthene | 74 U | 76 U | 73 U | 70 U |
| Pyrene | 74 U | 76 U | 73 U | 70 U |
| Butylbenzylphthalate | 74 U | 76 U | 73 U | 70 U |
| 3,3'-Dichlorobenzidine | 370 U | 380 U | 370 U | 350 U |
| Benzo (a) anthracene | 74 U | 76 U | 73 U | 70 U |
| bis (2-Ethylhexyl) phthalate | 74 U | 76 U | 73 U | 70 U |
| Chrysene | 74 U | 76 U | 73 U | 70 U |
| Di-n-Octyl phthalate | 74 U | 76 U | 73 U | 70 U |
| Benzo (b) fluoranthene | 74 U | 76 U | 73 U | 70 U |
| Benzo (k) fluoranthene | 74 U | 76 U | 73 U | 70 U |
| Benzo (a) pyrene | 74 U | 76 U | 73 U | 70 U |
| Indeno (1,2,3-cd) pyrene | 74 U | 76 U | 73 U | 70 U |
| Dibenz (a,h) anthracene | 74 U | 76 U | 73 U | 70 U |
| Benzo (g,h,i) perylene | 74 U | 76 U | 73 U | 70 U |

Notes:

U = Analyte not detected at detection limit reported.

file: paxsvoc.wq1

2-21-95

mlm

TABLE 5
PAXTON SALES CORPORATION
REMEDIAL INVESTIGATION
SOIL ANALYSIS DATA
CYANIDE, TOTAL SOLIDS

| ANALYTE | | MW-1-10.0 | MW-1-30.1 | MW-2-20.5 | MW-3-19.8 |
|--------------|---------|-----------|-----------|-----------|-----------|
| Cyanide | (mg/kg) | 0.18 U | 0.20 U | 0.11 U | 0.20 U |
| Total Solids | (%) | 92.3 | 91.3 | 88.9 | 92.9 |

Notes:

U = Analyte not detected at detection limit reported.

file: paxcyan.wq1
2-21-95
mlm

TABLE 6
PAXTON SALES CORPORATION
REMEDIAL INVESTIGATION
SOIL ANALYSIS DATA
METALS (mg/kg)

| ANALYTE | MW-1-5.3 | MW-1-15.2 | MW-2-20.3 | MW-3-14.8 |
|--------------------------------|----------|-----------|-----------|-----------|
| Antimony | 0.1 U | 0.1 U | 0.1 U | 0.1 U |
| Arsenic - ^{1,4 mg/kg} | 3.4 | 2.7 | 3.1 | 3.5 |
| Beryllium | 0.3 | 0.3 | 0.3 | 0.2 |
| Cadmium | 0.2 U | 0.2 U | 0.2 U | 0.2 |
| Chromium - ^{II or VI} | 15.5 | 16.9 | 24.6 | 14.7 |
| Copper | 17.2 | 22.7 | 21.9 | 27.8 |
| Lead | 3.7 | 5.0 | 3.9 | 3.6 |
| Mercury | 0.05 U | 0.05 U | 0.04 U | 0.04 U |
| Nickel | 11 | 14 | 16 | 11 |
| Selenium | 0.1 U | 0.1 U | 0.1 | 0.2 |
| Silver | 0.3 U | 0.3 U | 0.3 U | 0.3 U |
| Thallium | 0.2 | 0.2 | 0.5 U | 0.2 |
| Zinc | 47.9 | 49.4 | 53.2 | 49.1 |

Notes:

U = Analyte not detected at detection limit reported.

file: paxmetl.wq1
2-21-95
mlm

TABLE 7
 PAXTON SALES CORPORATION
 REMEDIAL INVESTIGATION
 GROUNDWATER ANALYSIS DATA

| ANALYTE | MW-1-123 | MW-1A-123* | MW-2-123 | MW-3-123 |
|--|----------|------------|----------|----------|
| <u>Volatile Organic Compounds (ug/L)</u> | | | | |
| Chloromethane | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| Vinyl Chloride | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| Chloroethane | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| Methylene Chloride | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| 1,1-Dichloroethene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,1-Dichloroethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| trans-1,2-Dichloroethene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Chloroform | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,2-Dichloroethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,1,1-Trichloroethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Carbon Tetrachloride | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Bromodichloromethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,2-Dichloropropane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Trichloroethene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Dibromochloromethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,1,2-Trichloroethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| trans-1,3-Dichloropropene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 2-Chloroethylvinylether | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| Bromoform | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Tetrachloroethene | 2.3 | 2.2 | 1.7 | 3.1 |
| 1,1,2,2-Tetrachloroethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Chlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Trichlorofluoromethane | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| 1,2-Dichlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,3-Dichlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,4-Dichlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Bromoethane | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| Dibromomethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,1,1,2-Tetrachloroethane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,2,3-Trichloropropane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Bromobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 2-Chlorotoluene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 4-Chlorotoluene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |

TABLE 7
PAXTON SALES CORPORATION
REMEDIAL INVESTIGATION
GROUNDWATER ANALYSIS DATA

| ANALYTE | MW-1-123 | MW-1A-123* | MW-2-123 | MW-3-123 |
|--|----------|------------|----------|----------|
| <u>Semivolatile Organic Compounds (ug/L)</u> | | | | |
| Phenol | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| Bis-(2-chloroethyl) Ether | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| 2-Chlorophenol | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,3-Dichlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 1,4-Dichlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Benzyl Alcohol | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| 1,2-Dichlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 2-Methylphenol | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| 2,2'-Oxybis (1-Chloropropane) | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 4-Methylphenol | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| N-Nitroso-Di-N-Propylamine | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| Hexachloroethane | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| Nitrobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Isophorone | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 2-Nitrophenol | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| 2,4-Dimethylphenol | 3.0 U | 3.0 U | 3.0 U | 3.0 U |
| Benzoic Acid | 10 U | 10 U | 10 U | 10 U |
| bis (2-chloroethoxy) Methane | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 2,4-Dichlorophenol | 3.0 U | 3.0 U | 3.0 U | 3.0 U |
| 1,2,4-Trichlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Naphthalene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 4-Chloroaniline | 3.0 U | 3.0 U | 3.0 U | 3.0 U |
| Hexachlorobutadiene | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| 4-Chloro-3-methylphenol | 2.0 U | 2.0 U | 2.0 U | 2.0 U |
| 2-Methylnaphthalene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Hexachlorocyclopentadiene | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| 2,4,6-Trichlorophenol | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| 2,4,5-Trichlorophenol | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| 2-Chloronaphthalene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 2-Nitroaniline | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| Dimethylphthalate | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Acenaphthylene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 3-Nitroaniline | 6.0 U | 6.0 U | 6.0 U | 6.0 U |
| Acenaphthene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 2,4-Dinitrophenol | 10 U | 10 U | 10 U | 10 U |
| 4-Nitrophenol | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| Dibenzofuran | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 2,6-Dinitrotoluene | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| 2,4-Dinitrotoluene | 5.0 U | 5.0 U | 5.0 U | 5.0 U |

TABLE 7
 PAXTON SALES CORPORATION
 REMEDIAL INVESTIGATION
 GROUNDWATER ANALYSIS DATA

| ANALYTE | MW-1-123 | MW-1A-123* | MW-2-123 | MW-3-123 |
|--|----------|------------|----------|----------|
| Semivolatile Organic Compounds (ug/L) (continued) | | | | |
| Diethylphthalate | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 4-Chlorophenyl-phenylether | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Fluorene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 4-Nitroaniline | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| 4,6-Dinitro-2-Methylphenol | 10 U | 10 U | 10 U | 10 U |
| N-Nitrosodiphenylamine | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 4-Bromophenyl-phenylether | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Hexachlorobenzene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Pentachlorophenol | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| Phenanthrene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Carbazole | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Anthracene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Di-n-Butylphthalate | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Fluoranthene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Pyrene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Butylbenzylphthalate | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| 3,3'-Dichlorobenzidine | 5.0 U | 5.0 U | 5.0 U | 5.0 U |
| Benzo (a) anthracene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| bis (2-Ethylhexyl) phthalate | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Chrysene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Di-n-Octyl phthalate | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Benzo (b) fluoranthene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Benzo (k) fluoranthene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Benzo (a) pyrene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Indeno (1,2,3-cd) pyrene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Dibenz (a,h) anthracene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |
| Benzo (g,h,i) perylene | 1.0 U | 1.0 U | 1.0 U | 1.0 U |

TABLE 7
PAXTON SALES CORPORATION
REMEDIAL INVESTIGATION
GROUNDWATER ANALYSIS DATA

| ANALYTE | MW-1-123 | MW-1A-123* | MW-2-123 | MW-3-123 |
|----------------------|----------|------------|----------|----------|
| METALS (mg/L) | | | | |
| Antimony | 0.001 UJ | 0.001 UJ | 0.001 UJ | 0.001 UJ |
| Arsenic | --- R | --- R | --- R | --- R |
| Beryllium | 0.003 | 0.002 | 0.003 | 0.001 |
| Cadmium | 0.002 U | 0.002 U | 0.002 U | 0.002 U |
| Chromium | 0.106 J | 0.085 J | 0.127 J | 0.042 J |
| Copper | 0.210 | 0.185 | 0.289 | 0.074 |
| Lead | 0.038 | 0.034 | 0.046 | 0.015 |
| Mercury | 0.0005 | 0.0005 | 0.0005 | 0.001 U |
| Nickel | 0.09 | 0.08 | 0.10 | 0.04 |
| Selenium | 0.005 U | 0.005 U | 0.007 | 0.007 |
| Silver | 0.003 U | 0.003 U | 0.003 U | 0.003 U |
| Thallium | 0.002 | 0.002 | 0.002 | 0.001 U |
| Zinc | 0.343 | 0.289 | 0.403 | 0.122 |
| WTPH-D (mg/L) | 0.25 U | 0.25 U | 0.25 U | 0.25 U |
| CYANIDE (mg/L) | 0.020 J | 0.031 J | 0.005 UJ | 0.005 UJ |

Notes:

* = MW-1A-123 is a duplicate of MW-1-123.

J = Estimated concentration.

U = Undetected at the reported detection limit.

UJ = Undetected at the estimated detection limit listed.

R = Data rejected by quality assurance review.

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2-21-95

mlm

Appendix A

Soil Boring and Well Logs

Soil Classification System

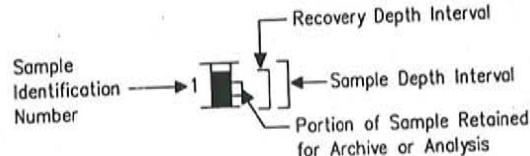
| MAJOR DIVISIONS | | GRAPHIC SYMBOL | USCS LETTER SYMBOL ⁽¹⁾ | TYPICAL DESCRIPTIONS ⁽²⁾⁽³⁾ |
|--|--|---|-----------------------------------|--|
| COARSE-GRAINED SOIL <small>(More than 50% of material is larger than No.200 sieve size)</small> | GRAVEL AND GRAVELLY SOIL <small>(More than 50% of coarse fraction retained on No.4 sieve)</small> | CLEAN GRAVEL <small>(Little or no fines)</small> | GW | Well-graded gravel; gravel/sand mixture(s); little or no fines |
| | | GRAVEL WITH FINES <small>(Appreciable amount of fines)</small> | GP | Poorly graded gravel; gravel/sand mixture(s); little or no fines |
| | SAND AND SANDY SOIL <small>(More than 50% of coarse fraction passed through No.4 sieve)</small> | CLEAN SAND <small>(Little or no fines)</small> | GM | Silty gravel; gravel/sand/silt mixture(s) |
| | | SAND WITH FINES <small>(Appreciable amount of fines)</small> | GC | Clayey gravel; gravel/sand/clay mixture(s) |
| | SILT AND CLAY <small>(Liquid Limit less than 50)</small> | CLEAN SILT | SW | Well-graded sand; gravelly sand; little or no fines |
| | | SILT WITH FINES | SP | Poorly graded sand; gravelly sand; little or no fines |
| | | SILT CLAY | SM | Silty sand; sand/silt mixture(s) |
| | HIGHLIGHTED SILT AND CLAY <small>(Liquid Limit greater than 50)</small> | CLAY | SC | Clayey sand; sand/clay mixture(s) |
| | | CLAY | ML | Inorganic silt and very fine sand; rock flour; silty or clayey fine sand or clayey silt with slight plasticity |
| FINE-GRAINED SOIL <small>(More than 50% of material is smaller than No.200 sieve size)</small> | SILT AND CLAY <small>(Liquid Limit greater than 50)</small> | CLAY | CL | Inorganic clay of low to medium plasticity; gravelly clay; sandy clay; silty clay; lean clay |
| | | ORGANIC SILT | OL | Organic silt; organic, silty clay of low plasticity |
| | | ORGANIC CLAY | MH | Inorganic silt; micaceous or diatomaceous fine sand or silty soil |
| | HIGHLY ORGANIC SOIL | ORGANIC CLAY | CH | Inorganic clay of high plasticity; fat clay |
| | | PEAT | OH | Organic clay of medium to high plasticity; organic silt |
| | OTHER | PEAT | PT | Peat; humus; swamp soil with high organic content |
| | | ASPHALT OR CONCRETE | AC | Pavement; Asphalt or Concrete |

Notes:

1. USCS letter symbols correspond to the symbols used by the Unified Soil Classification System and ASTM Classification methods. Dual letter symbols (e.g., SM-SP) for a sand or gravel indicate a soil with an estimated 5-15% fines. Multiple letter symbols (e.g., ML/CL) indicate borderline or multiple soil classifications.
2. Soil classifications are based on the general approach presented in the *Standard Practice for Description and Identification of Soils (Visual-Manual Procedure)*, as outlined in ASTM D2488. Where laboratory index testing has been conducted, soil classifications are based on the *Standard Test Method for Classification of Soils for Engineering Purposes*, as outlined in ASTM D2487.
3. Soil description terminology is based on visual estimates (in the absence of laboratory test data) of the percentages of each soil type and is defined as follows: Primary Constituent: >50% - "GRAVEL," "SAND," "SILT," "CLAY," etc.
 Secondary Constituents: >30% and ≤50% - "very gravelly," "very sandy," "very silty," etc.
 >15% and ≤30% - "gravelly," "sandy," "silty," etc.
 Additional Constituents: >5% and ≤15% - "with gravel," "with sand," "with silt," etc.
 ≤5% - "trace gravel," "trace sand," "trace silt," etc., or not noted.

Key

SAMPLE NUMBER & INTERVAL



TEST DATA

| Code | Description |
|------|-------------|
|------|-------------|

SAMPLER TYPE

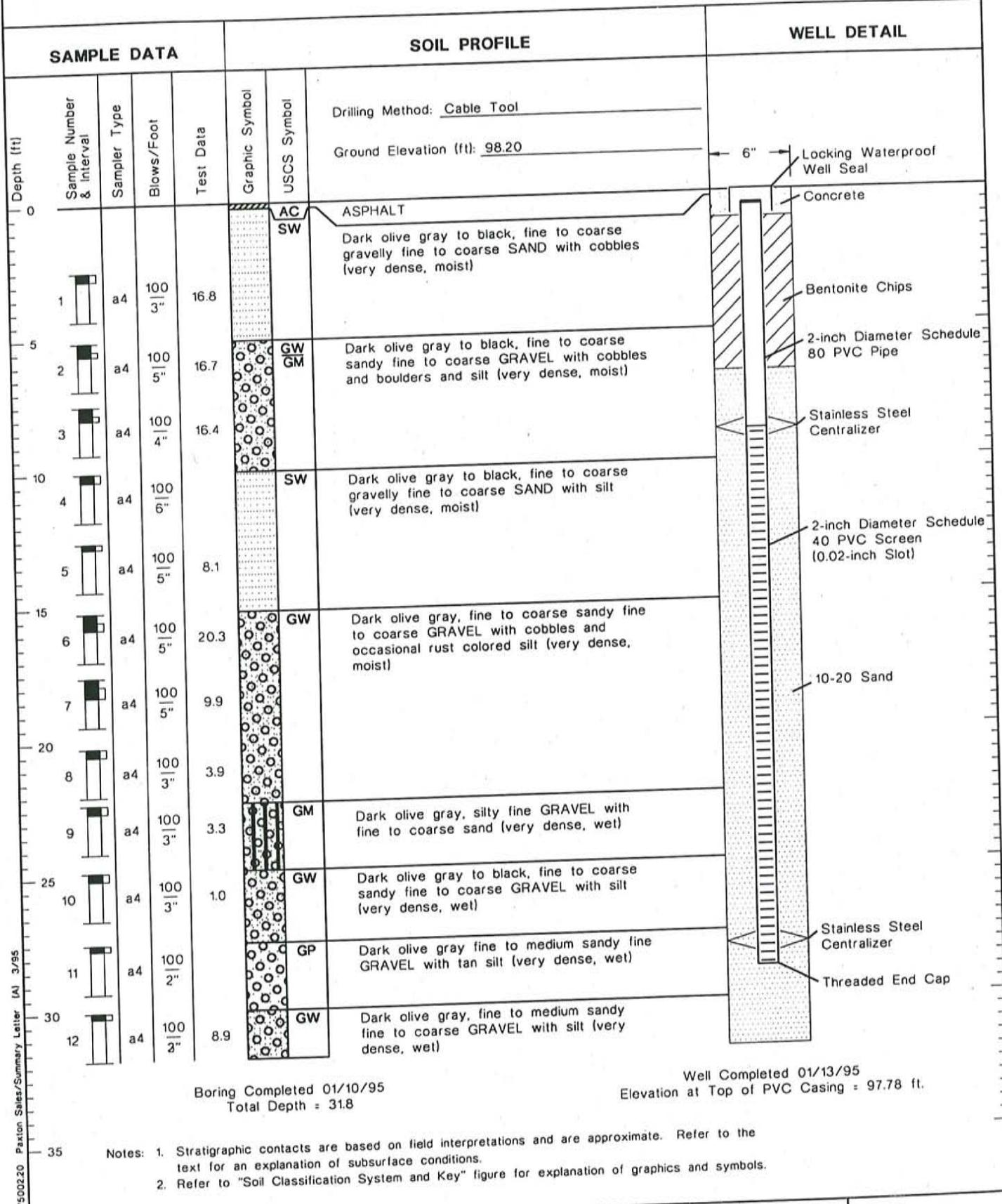
| Code | Description |
|-------|--|
| a | 3.25-inch O.D., 2.42-inch I.D. Split Spoon Sampler |
| b | 2.00-inch O.D., 1.50-inch I.D. Split Spoon Sampler |
| c | Shelby Tube |
| d | Grab Sample |
| e | 3.00-inch I.D. Core Barrel Sampler |
| 1 | 300-lb Hammer, 30-inch Drop |
| 2 | 140-lb Hammer, 30-inch Drop |
| 3 | Pushed |
| 4 | 350-lb. Hammer, 30-inch Drop |
| OTHER | |

ATD

Approximate Water Elevation At Time of Drilling (ATD)
or On Date Noted



MW-1



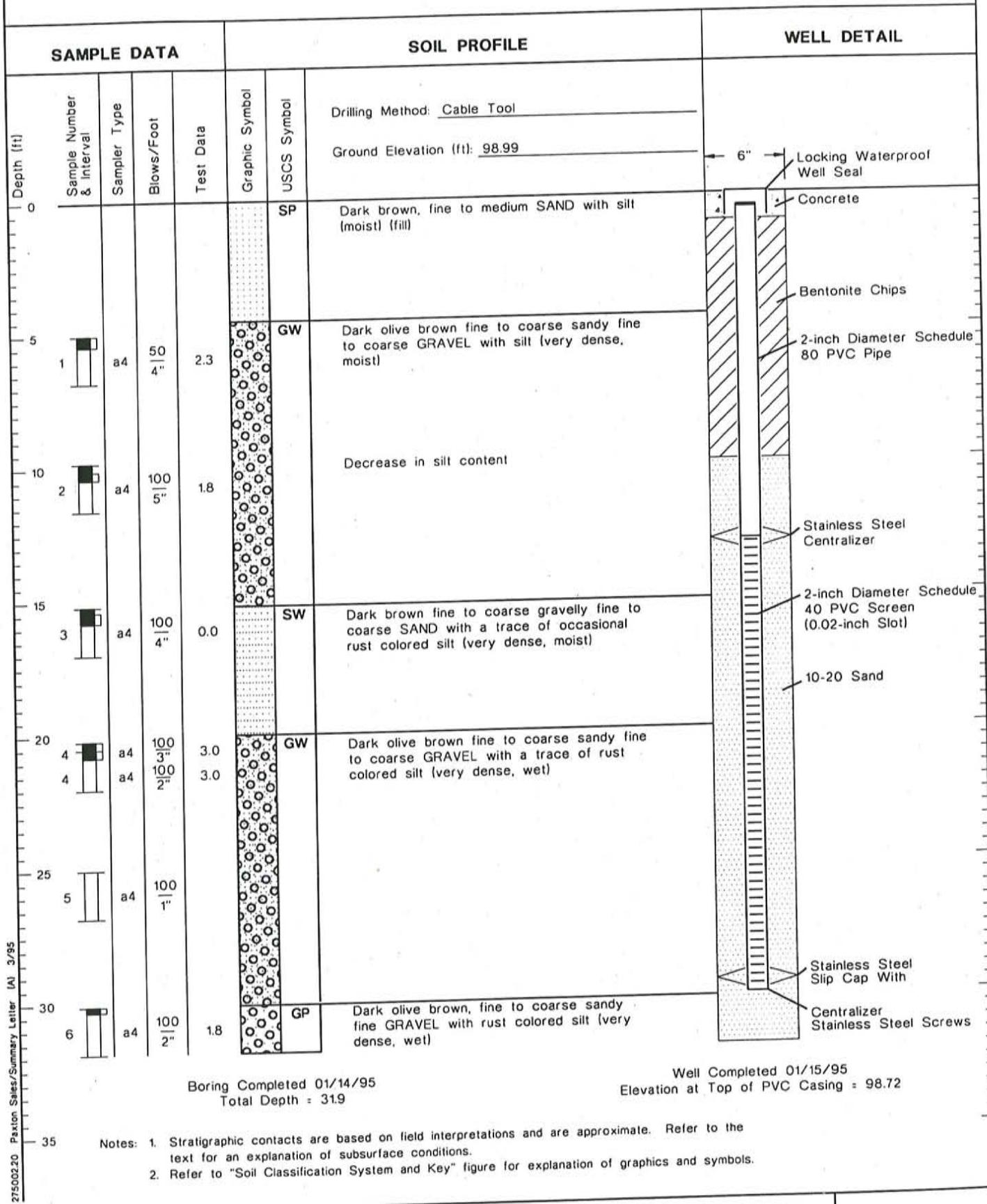
27500220



Boring and Monitoring Well MW-1

Figure A-2

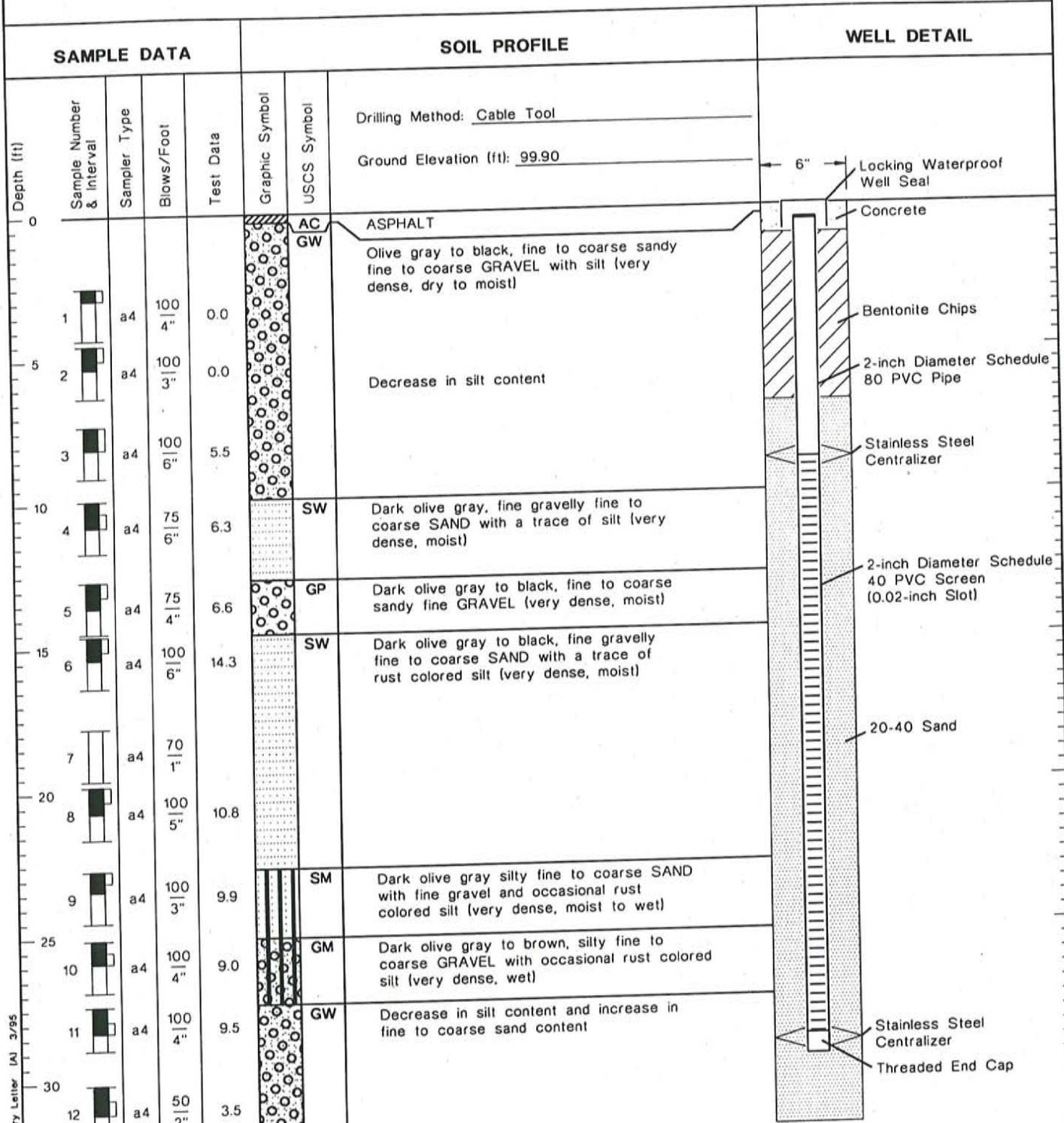
MW-2



Boring and Monitoring Well MW-2

Figure A-3

MW-3



Boring Completed 01/09/95
Total Depth = 31.9

Well Completed 01/09/95
Elevation at Top of PVC Casing = 99.55

- Notes: 1. Stratigraphic contacts are based on field interpretations and are approximate. Refer to the text for an explanation of subsurface conditions.
2. Refer to "Soil Classification System and Key" figure for explanation of graphics and symbols.



Boring and Monitoring Well MW-3

Figure A-4

Appendix B

Data Validation Technical Memorandum



LANDAU
ASSOCIATES,
INC.

Environmental and Geotechnical Services

TECHNICAL MEMORANDUM

TO: Bill Evans, Project Manager, Landau Associates, Inc.

FROM: Mandie MacDonald, Landau Associates, Inc. *mwm*

DATE: March 1, 1995

RE: **PAXTON SALES CORPORATION
REMEDIAL INVESTIGATION
LABORATORY DATA VALIDATION**

This memorandum provides the results of a focused data validation associated with 25 soil samples, 4 water samples, and 1 trip blank collected during the remedial investigation at the Paxton Sales Corporation site. Samples were analyzed by Analytical Resources Inc. (ARI), Seattle, Washington. This data quality validation covers ARI data packages J442, J468, J477, and J547. Samples were analyzed for one or more of the following: volatile organic compounds (EPA Method 8260), semivolatile organic compounds (EPA Method 8270), metals (WPA Method 6010/7000 series), total petroleum hydrocarbons (TPH) (Washington Method WTPH-HCID), diesel range petroleum hydrocarbons (Washington Method WTPH-d), and cyanide (EPA method 335.2). Sample data were evaluated in accordance with portions of the *National Functional Guidelines for Organic Data Review* (EPA 1991). The following parameters were evaluated:

- Chain-of-custody records
- Holding time
- Blank results (laboratory method and field trip)
- Surrogate recoveries (organic analyses only)
- Matrix spikes and matrix spike duplicates
- Laboratory control and standard reference samples
- Laboratory duplicates
- Field duplicates
- Completeness and overall data quality.

CHAIN-OF-CUSTODY RECORDS

Signed chain-of-custody records were attached to each of the data packages. Two of the seven bottles containing one of the water samples were broken during transit, but the lab was able to complete all requested analyses using the remaining bottles for that sample. All other samples were received by the laboratory in good condition.

VOLATILE ORGANIC COMPOUND ANALYSES

HOLDING TIMES

Acceptable without further qualification. The specified holding time between sample collection and analysis was met for all samples. No data qualifiers are needed.

BLANK RESULTS

Laboratory Method Blanks

Acceptable with the exceptions described below. At least one method blank was analyzed with each batch of samples. Methylene chloride was reported in the method blank associated with data package J442, but not in any of the associated samples. Several compounds were reported in the first method blank associated with data package J468. A second method blank was run and all volatile organic compounds were reported as undetected; however, the laboratory elected to qualify its sample results based on the first method blank. Reviewing sample results based on the first method blank resulted in low levels of methylene chloride (reported in two samples) being qualified as undetected. No other data qualifiers are needed. Data qualifiers are summarized in Table 1.

Field Trip Blanks

Acceptable without further qualification. One trip blank was submitted for analysis with the water samples (ARI J547). All volatile organic compounds were reported as undetected in the trip blank. No data qualifiers are needed.

SURROGATE RECOVERIES

Acceptable without further qualification. Surrogate recoveries were reported within control limits for all surrogate spiking compounds in all samples. No data qualifiers are needed.

MATRIX SPIKE AND MATRIX SPIKE DUPLICATE RESULTS

Matrix spike and matrix spike duplicate (MS/MSD) analyses were not performed; however, the laboratory did perform blank spike analyses with each sample set. All spike recoveries were reported within MS control limits. No data qualifiers are needed.

LABORATORY CONTROL/STANDARD REFERENCE SAMPLES

No laboratory control or standard reference samples were analyzed for volatile organic compounds; however, the blank spike samples described above can be considered laboratory control samples for calibration verification purposes. No data qualifiers are needed.

LABORATORY DUPLICATES

Not performed or required for volatile organic compound analysis.

FIELD DUPLICATES

Two blind field duplicate pairs were submitted for volatile organic compound analysis, one with the water samples and one with the soil samples associated with data package J477. All volatile organic compounds were reported as undetected in both samples of the soil duplicate pair. All volatile organic compounds except tetrachloroethene were reported as undetected in the water duplicate pair. The relative percent difference (RPD) between the reported tetrachloroethene results in this pair was calculated to be within acceptable control limits. No data qualifiers are needed. Field duplicate results are summarized in Table 2.

SEMOVOLATILE ORGANIC COMPOUND ANALYSES

HOLDING TIMES

Acceptable without further qualification. The specified holding times between sample collection, extraction, and analysis were met for all samples. No data qualifiers are needed.

BLANK RESULTS

Laboratory Method Blanks

Acceptable without further qualification. At least one method blank was analyzed with each batch of samples. All semivolatile organic compounds were reported as undetected in the method blanks. No data qualifiers are needed.

Field Trip Blanks

Not required for semivolatile organic analysis.

SURROGATE RECOVERIES

Acceptable without further qualification. Surrogate recoveries were reported within control limits for all surrogate spiking compounds in all samples. No data qualifiers are needed.

MATRIX SPIKE AND MATRIX SPIKE DUPLICATE RESULTS

Acceptable with the following discussion. Matrix spike analyses were performed with data packages 468 and J477. No matrix spike duplicate analyses were performed. Recoveries of all spiking compounds except pentachlorophenol were reported within acceptable control limits. Pentachlorophenol recovery was reported below the control limit in both MS samples, indicating that sample results may be biased low; however data are not qualified based solely on matrix spike results. No data qualifiers are needed.

A blank spike sample was analyzed with data package J442. Recoveries of all spiking compounds were reported within MS/MSD control limits. No data qualifiers are needed.

LABORATORY CONTROL/STANDARD REFERENCE SAMPLES

A laboratory control sample was analyzed with data package J547. All analytes were reported within MS/MSD control limits. No data qualifiers are needed.

LABORATORY DUPLICATES

Not required for semivolatile organic compound analysis.

FIELD DUPLICATE RESULTS

A blind field duplicate was submitted for semivolatile organic compound analysis with the water samples (ARI J547). All semivolatile organic compounds were reported as undetected in both samples of the duplicate pair. No data qualifiers are needed.

METALS ANALYSES

HOLDING TIMES

Acceptable without further qualification. The specified holding time between sample collection and analysis was met for all samples. No data qualifiers are needed.

BLANK RESULTS

Laboratory Method Blanks

Acceptable with the following exceptions. At least one method blank was analyzed with each sample batch. Zinc was reported at low levels in the each of the method blanks associated with the soil samples and lead was reported in the method blank associated with the water samples. Reported detections less than five times the blank concentration were qualified as undetected (U). No sample results were affected. No further data qualifiers are needed.

Field Trip Blanks

Not required for metals analysis.

SURROGATE RECOVERIES

Not applicable to metals analysis.

MATRIX SPIKE AND MATRIX SPIKE DUPLICATE RESULTS

Acceptable with the following exceptions. Blank spike samples were analyzed with each of the soil sample batches. All spiking analytes were reported within control limits for MS analysis. An MS sample was analyzed with the water samples. The reported recovery for antimony was well below the control limit. All results for antimony were qualified as estimated (J for detects, UJ for nondetects). Data qualifiers are summarized in Table 1. No additional data qualifiers are needed.

LABORATORY DUPLICATES

One laboratory duplicate was analyzed with the water samples. Relative percent differences (RPD) between the sample and its duplicate were reported within acceptable control limits for all analytes. No data qualifiers are needed.

FIELD DUPLICATE RESULTS

One field duplicate pair was analyzed for metals with the water samples. Relative percent differences (RPD) between the duplicate samples were calculated to be within acceptable control limits for all analytes except chromium and arsenic. The RPD for chromium was slightly below the control limit; all sample results for chromium were qualified as estimated (J for detects, UJ for nondetects). The RPD for arsenic exceeded the control limit by a large enough margin that all arsenic results for the water samples were rejected (R). Data qualifiers are summarized in Table 1.

TPH ANALYSES (WTPH-HCID AND WTPH-d)

HOLDING TIMES

Acceptable without further qualification. The specified holding time between sample collection and analysis was met for all samples. No data qualifiers are needed.

BLANK RESULTS

Laboratory Method Blanks

Acceptable without further qualification. At least one laboratory method blank was analyzed with each batch of samples. All analytes were reported as undetected in the method blanks. No data qualifiers are needed.

Field Trip Blanks

Not required for TPH analysis.

SURROGATE RECOVERIES

Acceptable without further qualification. Surrogate recoveries were reported within control limits for all surrogate spiking compounds in all samples. No data qualifiers are needed.

MATRIX SPIKE/MATRIX SPIKE DUPLICATE RESULTS

Matrix spike samples were not performed with the TPH analyses; however, blank spike samples were analyzed with data packages J442, J468, and J547. Recoveries for all spiking compounds were reported within MS/MSD control limits for all samples. No data qualifiers are needed.

LABORATORY CONTROL/STANDARD REFERENCE SAMPLES

A laboratory control sample was analyzed with data package J477. Recoveries of all analytes were reported within the control limits for matrix spike samples. No data qualifiers are needed.

LABORATORY DUPLICATES

No laboratory duplicates were performed for TPH analysis.

FIELD DUPLICATE RESULTS

One blind field duplicate was analyzed with the water samples. Diesel range petroleum hydrocarbons were reported as undetected in both samples of the duplicate pair. No data qualifiers are needed.

CYANIDE ANALYSES

HOLDING TIMES

Acceptable without further qualification. The specified holding time between sample collection and analysis was met for all samples. No data qualifiers are needed.

BLANK RESULTS

Laboratory Method Blanks

Acceptable without further qualification. At least one method blank was analyzed with each batch of samples. Cyanide was reported as undetected in all of the method blanks. No data qualifiers are needed.

Field Trip Blanks

Not required for inorganic analyses.

SURROGATE RECOVERIES

Not applicable to inorganic analyses.

MATRIX SPIKE/MATRIX SPIKE DUPLICATE RESULTS

Matrix spike samples were analyzed with sample batches J442, J477, and J547. Recoveries of the spiked analytes were reported within acceptable control limits for all samples. No matrix spike duplicate samples were analyzed.

LABORATORY CONTROL/STANDARD REFERENCE SAMPLES

Standard reference material samples were analyzed with each batch of field samples. Recoveries of the standard reference concentrations were reported within the control limits for matrix spike samples. No data qualifiers are needed.

LABORATORY DUPLICATES

Laboratory duplicate samples were analyzed with sample batches J442, J477, and J547. Relative percent differences between the two samples in each duplicate pair were reported within acceptable control limits. No data qualifiers are needed.

FIELD DUPLICATE RESULTS

One blind field duplicate was submitted for cyanide analysis with the water samples. The calculated relative percent difference between the two samples of the duplicate pair exceeded the established control limit. Because of this variability between duplicate samples, all cyanide results for the water samples were qualified as estimated (J). Data qualifiers are summarized in Table 1.

OVERALL DATA QUALITY AND COMPLETENESS

All data except cyanide in the water samples are considered usable, with qualifiers were needed as described above and summarized in Table 1. The four cyanide analyses in the water samples were rejected due to variability between duplicate samples. Data completeness is therefore calculated to be 99.6%.

REFERENCES

EPA. 1991. *National Functional Guidelines for Organic Data Review*. U.S. Environmental Protection Agency.

TABLE 1
PAXTON SALES CORPORATION
REMEDIAL INVESTIGATION LABORATORY DATA VALIDATION
SUMMARY OF DATA QUALIFIERS

| Data Package | Samples Affected | Analyte | Qualifier Assigned | Reason |
|--------------|---|--------------------|---|--|
| J468 | MW-1-5.1 MW-1-7.5 | Methylene Chloride | Concentrations less than 10 times method blank concentration assigned U (Undetected) | Method blank contamination. |
| J547 | MW-1-123 MW-1A-123 MW-2-123 MW-3-123 | Cyanide | All results qualified as estimated concentration (J) | Field duplicate results outside control limits. |
| J547 | MW-1-123 MW-1A-123 MW-2-123 MW-3-123 | Antimony | All results qualified as estimated concentration (J) | Low spike recovery. |
| J547 | MW-1-123 MW-1A-123 MW-2-123 MW-3-123 | Chromium | All results qualified as estimated concentration (J) | Field duplicate results outside control limits. |
| J547 | MW-1-123 MW-1A-123 MW-2-123 MW-3-123 | Arsenic | All results rejected (R) | Field duplicate results outside control limits. |

TABLE 2
PAXTON SALES CORPORATION
REMEDIAl INVESTIGATION LABORATORY DATA VALIDATION
SUMMARY OF FIELD DUPLICATE RESULTS

| Volatile Organic Compounds | MW-2-20.2 | MW-2-120.2 | RPD |
|--------------------------------|------------|------------|-------|
| All Analytes | Undetected | Undetected | N/A |
| Volatile Organic Compounds | MW-1-123 | MW-1A-123 | RPD |
| Tetrachloroethene | 2.3 ug/L | 2.2 ug/L | 4.4% |
| All Other Analytes | Undetected | Undetected | N/A |
| Semivolatile Organic Compounds | MW-1-123 | MW-1A-123 | RPD |
| All Analytes | Undetected | Undetected | N/A |
| Metals | MW-1-123 | MW-1A-123 | RPD |
| Arsenic | .024 | .007 | 110%* |
| Beryllium | .003 | .002 | A/L |
| Chromium | .106 | .085 | 21.8% |
| Copper | .210 | .185 | 12.7% |
| Lead | .038 | .034 | 11.1% |
| Mercury | .0005 | .0005 | 0 |
| Nickel | .009 | .08 | 11.8% |
| Thallium | .002 | .002 | 0 |
| Zinc | .343 | .289 | 17.1% |
| All other metals | Undetected | Undetected | N/A |

TABLE 2
 PAXTON SALES CORPORATION
 REMEDIAL INVESTIGATION LABORATORY DATA VALIDATION
 SUMMARY OF FIELD DUPLICATE RESULTS

| WTPH-d | MW-1-123 | MW-1A-123 | RPD |
|---------------------|------------------------|-------------------------|------|
| Diesel Hydrocarbons | Undetected | Undetected | N/A |
| Cyanide | MW-1-123 0.020 mg/L | MW-1A-123 0.031 mg/L | 43%* |

Notes:

RPD = Relative percent difference.
 N/A = Not applicable, RPD cannot be calculated for undetected analytes.
 A/L = Alternate limit applies for concentrations less than 5 times the detection limit.
 * = Outside control limits.

Appendix C

Laboratory Reports



Analytical Resources, Incorporated
Analytical Chemists and Consultants

cc to MIM

January 30, 1995

Mr. Bill Evans
Landau Associates, Inc.
3600 Port of Tacoma Road
Suite 501
Tacoma, WA 98424

RECEIVED

JAN 3 1 1995

LANDAU ASSOCIATES, INC.
TACOMA

RE: Project No. 275002.23 Paxton Sales / ARI Job No. J442

Dear Mr. Evans:

Please find enclosed original results and Chain-of-Custody Record (COC NO. 2511) for the above-referenced project. Analytical Resources, Incorporated (ARI), accepted the following five soil samples on January 11, 1995:

MW-3-15

MW-3-14.8

MW-3-19.8

MW-3-20

MW-3-14.6

ARI received the samples intact with no discrepancies on the chain-of-custody documentation. The laboratory analyzed the samples for volatile organic compounds by EPA method 8260, semivolatile organics by EPA method 8270, total petroleum hydrocarbons by WDOE method WTPH-HCID, metals following the EPA method 6010/7000 series, and cyanide by EPA method 335.2.

Sample analysis was routine and no analytical complications were noted with this sample delivery group.

As always, a copy of this report and all associated raw data will remain on file with ARI. If you have any questions or require additional information, please feel free to contact me at your convenience. If I am unavailable, you can leave a message on my voice mail and I will return your call as soon as possible.

Sincerely,

ANALYTICAL RESOURCES, INC.

Bryan D. Anderson

Bryan D. Anderson
Project Manager
(206)340-2866, ext. 116

enclosures
cc: file J442

BDA/dn



ANALYTICAL
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INCORPORATED

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

QC Report No: J442-Landau Associates, Incorporated (206) 621-7523 (FAX)

TOTAL PETROLEUM HYDROCARBONS
WA HCID Method by GC/FID

Matrix: Soil

Project: 275002.23

Paxton Sales

Data Release Authorized:

Date Received: 01/11/95

Reported: 01/20/95

Calvin M. Newman

| Lab ID | Client Sample ID | Date Analyzed | Dilution Factor | Gas Range | Diesel Range | Oil Range | Surrogate Recovery |
|--------|------------------|---------------|-----------------|-----------|--------------|-----------|--------------------|
| 95-291 | Method Blank | 01/16/95 | 1:1 | 10 U | 10 U | 25 U | 92.5% |
| J442A | MW-3-15 | 01/17/95 | 1:1 | 10 U | 10 U | 25 U | 106% |

Values reported in ppm (mg/L).

Surrogate is Methyl Arachidate.

Gas value based on total peaks in the range from Toluene to C12.

Diesel value based on the total peaks in the range from C12 to C24.

Oil value based on the total peaks in the range from C24 to C32.

Data Qualifiers

- U Compound not detected at the given detection limit.
- X Value detected above linear range of instrument. Dilution required.
- J Indicates an estimated value below the calculated detection limit.
- S No value reported due to saturation of the detector. Dilution required.
- D Indicates the surrogate was not detected because of dilution of the extract.
- C Indicates a probable value which cannot be confirmed due to matrix interference.
- NR Indicates no recovery due to matrix interference and/or dilution.



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(206) 621-7523 (FAX)

TOTAL PETROLEUM HYDROCARBONS
WA HCID Method by GC/FID

Lab Sample ID: J442SB
LIMS ID: 95-291
Matrix: Soil

QC Report No: J442-Landau Associates, Incorporated
Project: 275002.23
Paxton Sales

Data Release Authorized:
Reported: 01/20/95

Date extracted: 01/16/95

LABORATORY CONTROL SAMPLE RECOVERY REPORT

| CONSTITUENT | SPIKE VALUE | SPIKE ADDED | % RECOVERY |
|---------------------------|----------------|----------------|---------------|
| LABORATORY CONTROL SAMPLE | | | |
| Diesel Range | 539 | 500 | 108% |

HCID Surrogate Recovery

| | | |
|-------------|------------------|------|
| Spike Blank | Methylarachidate | 104% |
|-------------|------------------|------|

Values reported in parts per million (mg/kg)

HCID SPIKE CONTROL LIMITS

| | |
|------------------|---------|
| Percent Recovery | 50-150% |
| Duplicate RPD | <50% |

Advisory QA Limits



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ORGANIC COMPOUND DATA REPORTING QUALIFIERS

- U** Indicates the compound was analyzed for, but not detected at the given detection limit.
- J** Indicates an estimated value when the result is less than the calculated detection limit.
- D** Indicates the surrogate/spike(s) was not detected, due to dilution of extract.
- NR** Indicates the surrogate recovery cannot be reported due to matrix interference.
- E** Indicates a value above the linear range of the detector.
Sample dilution required.
- S** Indicates no value reported due to saturation of the detector.
Dilution required.
- Y** Indicates a raised detection limit due to matrix interferences.
- NA** Indicates compound was not analyzed.
- M** Indicates an estimated value of analyte found and confirmed by analyst but with low spectral match.
- B** Indicates compound was found in the associated method blank.

ORGANICS ANALYSIS DATA SHEET
Volatile by Purge & Trap GC/MS

Page 1 of 1



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Sample No: MW-3-14.6

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J442E

LIMS ID: 95-295

Matrix: Soil

Data Release Authorized: ✓

Reported: 01/17/95

QC Report No: J442-Landau Associates, Incorporated

Project: 275002.23

Paxton Sales

Date Sampled: 01/06/95

Date Received: 01/11/95

Instrument: FINN1

Date Analyzed: 01/12/95

Sample Amount: 4.83 g dry Wt

Percent Moisture: 9.1%

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.1 U |
| 75-01-4 | Vinyl Chloride | 2.1 U |
| 75-00-3 | Chloroethane | 2.1 U |
| 75-09-2 | Methylene Chloride | 2.1 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.2 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.0 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.1 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-96-4 | Bromoethane | 2.1 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 106% |
| d8-Toluene | 97.2% |
| Bromofluorobenzene | 100% |
| d4-1,2-Dichlorobenzene | 99.5% |

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



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Sample No: Method Blank

Analytical
Chemists &
Consultants

Lab Sample ID: 011295MB
LIMS ID: 95-295
Matrix: Soil
Data Release Authorized: *MTH*
Reported: 01/17/95

QC Report No: J442-Landau Associates, Incorporated
Project: 275002.23
Paxton Sales
Date Sampled: NA
Date Received: NA

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Instrument: FINN1
Date Analyzed: 01/12/95
Sample Amount: 5.00 g dry Wt Equiv
Percent Moisture: NA

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 6.9 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.0 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 2.0 U |
| 75-69-4 | Trichlorofluoromethane | 1.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-83-9 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 93.0% |
| d8-Toluene | 97.0% |
| Bromofluorobenzene | 99.1% |
| d4-1,2-Dichlorobenzene | 96.2% |

ORGANICS ANALYSIS DATA SHEET
Volatile s by GC/MS
Page 1 of 1



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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J442SB
LIMS ID: 95-295
Matrix: Soil
Data Release Authorized: ✓
Reported: 01/18/95
Date Analyzed: 01/12/95

QC Report No: J442-Landau Associates, Incorporated
Project: 275002.23
Paxton Sales
Date Received: 01/11/95

| LABORATORY CONTROL SAMPLE CONSTITUENT | SPIKE VALUE | SPIKE AMT | % RECOVERY |
|--|----------------|--------------|---------------|
| Chloromethane | 68.4 | 50.0 | 137% |
| Vinyl Chloride | 67.0 | 50.0 | 134% |
| Chloroethane | 50.2 | 50.0 | 100% |
| Methylene Chloride | 45.6 | 50.0 | 91.2% |
| 1,1-Dichloroethene | 54.5 | 50.0 | 109% |
| 1,1-Dichloroethane | 47.4 | 50.0 | 94.8% |
| trans-1,2-Dichloroethene | 50.3 | 50.0 | 101% |
| Chloroform | 44.8 | 50.0 | 89.6% |
| 1,2-Dichloroethane | 47.3 | 50.0 | 94.6% |
| 1,1,1-Trichloroethane | 44.8 | 50.0 | 89.6% |
| Carbon Tetrachloride | 48.2 | 50.0 | 96.4% |
| Bromodichloromethane | 47.3 | 50.0 | 94.6% |
| 1,2-Dichloropropane | 50.1 | 50.0 | 100% |
| Trichloroethene | 47.2 | 50.0 | 94.4% |
| Dibromochloromethane | 46.0 | 50.0 | 92.0% |
| 1,1,2-Trichloroethane | 46.1 | 50.0 | 92.2% |
| trans-1,3-Dichloropropene | 45.6 | 50.0 | 91.2% |
| 2-Chloroethylvinylether | 49.8 | 50.0 | 99.6% |
| Bromoform | 46.7 | 50.0 | 93.4% |
| Tetrachloroethene | 46.1 | 50.0 | 92.2% |
| 1,1,2,2-Tetrachloroethane | 47.1 | 50.0 | 94.2% |
| Chlorobenzene | 45.6 | 50.0 | 91.2% |
| Trichlorofluoromethane | 44.3 | 50.0 | 88.6% |
| 1,2-Dichlorobenzene | 45.6 | 50.0 | 91.2% |
| 1,3-Dichlorobenzene | 46.2 | 50.0 | 92.4% |
| 1,4-Dichlorobenzene | 46.4 | 50.0 | 92.8% |
| Bromoethane | 52.2 | 50.0 | 104% |
| Dibromomethane | 44.6 | 50.0 | 89.2% |
| 1,1,1,2-Tetrachloroethane | 45.8 | 50.0 | 91.6% |
| 1,2,3-Trichloropropane | 47.2 | 50.0 | 94.4% |
| Bromobenzene | 47.3 | 50.0 | 94.6% |
| 2-Chlorotoluene | 48.9 | 50.0 | 97.8% |
| 4-Chlorotoluene | 45.1 | 50.0 | 90.2% |

| <u>Spike Blank Surrogate Recovery</u> | |
|---------------------------------------|-------|
| d4-1,2-Dichloroethane | 94.2% |
| d8-Toluene | 98.0% |
| Bromofluorobenzene | 98.4% |
| d4-1,2-Dichlorobenzene | 100.% |

Reported in ug/kg-dry-Wt



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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

ORGANIC COMPOUND DATA REPORTING QUALIFIERS

- U** Indicates the compound was analyzed for, but not detected at the given detection limit.
- J** Indicates an estimated value when the result is less than the calculated detection limit.
- D** Indicates the surrogate/spike(s) was not detected, due to dilution of extract.
- NR** Indicates the surrogate recovery cannot be reported due to matrix interference.
- E** Indicates a value above the linear range of the detector.
Sample dilution required.
- S** Indicates no value reported due to saturation of the detector.
Dilution required.
- Y** Indicates a raised detection limit due to matrix interferences.
- NA** Indicates compound was not analyzed.
- M** Indicates an estimated value of analyte found and confirmed by analyst but with low spectral match.
- B** Indicates compound was found in the associated method blank.



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 1 of 2

Lab Sample ID: J442MB

LIMS ID: 95-294

Matrix: Soil

Data Release Authorized: *MH*

Reported: 01/17/95

Sample No: Method Blank

Analytical
Chemists &
Consultants

QC Report No: J442-Landau Associates, Incorporated

Project: 275002.23

333 Ninth Ave. North
Seattle, WA 98109-5187

Paxton Sales

(206) 621-6490

Date Sampled: NA

(206) 621-7523 (FAX)

Date Received: NA

Date extracted: 01/13/95

Sample Amount: 30.0 g-dry-wt Equiv

Date analyzed: 01/16/95

Final Extract Volume: 2.0 mL

Instrument: FINN8

Dilution Factor: 1:1

GPC Cleanup: NO

Percent Moisture: NA

pH: NA

| CAS Number | Analyte | ug/kg |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 130 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 130 U |
| 95-57-8 | 2-Chlorophenol | 67 U |
| 541-73-1 | 1,3-Dichlorobenzene | 67 U |
| 106-46-7 | 1,4-Dichlorobenzene | 67 U |
| 100-51-6 | Benzyl Alcohol | 330 U |
| 95-50-1 | 1,2-Dichlorobenzene | 67 U |
| 95-48-7 | 2-Methylphenol | 130 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 67 U |
| 106-44-5 | 4-Methylphenol | 67 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 130 U |
| 67-72-1 | Hexachloroethane | 130 U |
| 98-95-3 | Nitrobenzene | 67 U |
| 78-59-1 | Isophorone | 67 U |
| 88-75-5 | 2-Nitrophenol | 330 U |
| 105-67-9 | 2,4-Dimethylphenol | 200 U |
| 65-85-0 | Benzoic Acid | 670 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 67 U |
| 120-83-2 | 2,4-Dichlorophenol | 200 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 67 U |
| 91-20-3 | Naphthalene | 67 U |
| 106-47-8 | 4-Chloroaniline | 200 U |
| 87-68-3 | Hexachlorobutadiene | 130 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 130 U |
| 91-57-6 | 2-Methylnaphthalene | 67 U |
| 77-47-4 | Hexachlorocyclopentadiene | 330 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 330 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 330 U |
| 91-58-7 | 2-Chloronaphthalene | 67 U |
| 88-74-4 | 2-Nitroaniline | 330 U |
| 131-11-3 | Dimethylphthalate | 67 U |
| 208-96-8 | Acenaphthylene | 67 U |
| 99-09-2 | 3-Nitroaniline | 400 U |
| 83-32-9 | Acenaphthene | 67 U |
| 51-28-5 | 2,4-Dinitrophenol | 670 U |
| 100-02-7 | 4-Nitrophenol | 330 U |
| 132-64-9 | Dibenzofuran | 67 U |
| 606-20-2 | 2,6-Dinitrotoluene | 330 U |



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2

Lab Sample ID: J442MB

LIMS ID: 95-294

Matrix: Soil

Data Release Authorized: *MWT*

Reported: 01/17/95

Sample No: Method Blank

Analytical
Chemists &
Consultants

QC Report No: J442-Landau Associates, Incorporated

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Project: 275002.23

Paxton Sales

Date Sampled: NA

(206) 621-6490

Date Received: NA

(206) 621-7523 (FAX)

Date extracted: 01/13/95

Sample Amount: 30.0 g-dry-wt Equiv

Date analyzed: 01/16/95

Final Extract Volume: 2.0 mL

Instrument: FINN8

Dilution Factor: 1:1

GPC Cleanup: NO

Percent Moisture: NA

pH: NA

| CAS Number | Analyte | ug/kg |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 330 U |
| 84-66-2 | Diethylphthalate | 67 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 67 U |
| 86-73-7 | Fluorene | 67 U |
| 100-01-6 | 4-Nitroaniline | 330 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 670 U |
| 86-30-6 | N-Nitrosodiphenylamine | 67 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 67 U |
| 118-74-1 | Hexachlorobenzene | 67 U |
| 87-86-5 | Pentachlorophenol | 330 U |
| 85-01-8 | Phenanthrene | 67 U |
| 86-74-8 | Carbazole | 67 U |
| 120-12-7 | Anthracene | 67 U |
| 84-74-2 | Di-n-Butylphthalate | 67 U |
| 206-44-0 | Fluoranthene | 67 U |
| 129-00-0 | Pyrene | 67 U |
| 85-68-7 | Butylbenzylphthalate | 67 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 330 U |
| 56-55-3 | Benzo(a)anthracene | 67 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 67 U |
| 218-01-9 | Chrysene | 67 U |
| 117-84-0 | Di-n-Octyl phthalate | 67 U |
| 205-99-2 | Benzo(b)fluoranthene | 67 U |
| 207-08-9 | Benzo(k)fluoranthene | 67 U |
| 50-32-8 | Benzo(a)pyrene | 67 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 67 U |
| 53-70-3 | Dibenz(a,h)anthracene | 67 U |
| 191-24-2 | Benzo(g,h,i)perylene | 67 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 63.6% | d5-Phenol | 57.0% |
| 2-Fluorobiphenyl | 74.8% | 2-Fluorophenol | 50.5% |
| d14-p-Terphenyl | 86.3% | 2,4,6-Tribromophenol | 49.8% |
| d4-1,2-Dichlorobenzene | 68.2% | d4-2-Chlorophenol | 60.9% |



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 1 of 2

Lab Sample ID: J442D

LIMS ID: 95-294

Matrix: Soil

Data Release Authorized: MTT

Reported: 01/17/95

Sample No: MW-3-20

Analytical
Chemists &
Consultants

QC Report No: J442-Landau Associates, Incorporated

Project: 275002.23

333 Ninth Ave. North
Seattle, WA 98109-5187

Paxton Sales

(206) 621-6490

Date Sampled: 01/06/95

(206) 621-7523 (FAX)

Date Received: 01/11/95

Date extracted: 01/13/95

Date analyzed: 01/16/95

Instrument: FINN8

GPC Cleanup: NO

Sample Amount: 28.6 g-dry-wt

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: 5.7%

pH: 6.7

| CAS Number | Analyte | ug/kg |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 140 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 140 U |
| 95-57-8 | 2-Chlorophenol | 70 U |
| 541-73-1 | 1,3-Dichlorobenzene | 70 U |
| 106-46-7 | 1,4-Dichlorobenzene | 70 U |
| 100-51-6 | Benzyl Alcohol | 350 U |
| 95-50-1 | 1,2-Dichlorobenzene | 70 U |
| 95-48-7 | 2-Methylphenol | 140 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 70 U |
| 106-44-5 | 4-Methylphenol | 70 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 140 U |
| 67-72-1 | Hexachloroethane | 140 U |
| 98-95-3 | Nitrobenzene | 70 U |
| 78-59-1 | Isophorone | 70 U |
| 88-75-5 | 2-Nitrophenol | 350 U |
| 105-67-9 | 2,4-Dimethylphenol | 210 U |
| 65-85-0 | Benzoic Acid | 700 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 70 U |
| 120-83-2 | 2,4-Dichlorophenol | 210 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 70 U |
| 91-20-3 | Naphthalene | 70 U |
| 106-47-8 | 4-Chloroaniline | 210 U |
| 87-68-3 | Hexachlorobutadiene | 140 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 140 U |
| 91-57-6 | 2-Methylnaphthalene | 70 U |
| 77-47-4 | Hexachlorocyclopentadiene | 350 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 350 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 350 U |
| 91-58-7 | 2-Chloronaphthalene | 70 U |
| 88-74-4 | 2-Nitroaniline | 350 U |
| 131-11-3 | Dimethylphthalate | 70 U |
| 208-96-8 | Acenaphthylene | 70 U |
| 99-09-2 | 3-Nitroaniline | 420 U |
| 83-32-9 | Acenaphthene | 70 U |
| 51-28-5 | 2,4-Dinitrophenol | 700 U |
| 100-02-7 | 4-Nitrophenol | 350 U |
| 132-64-9 | Dibenzofuran | 70 U |
| 606-20-2 | 2,6-Dinitrotoluene | 350 U |



ANALYTICAL
RESOURCES
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ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2

Lab Sample ID: J442D

LIMS ID: 95-294

Matrix: Soil

Data Release Authorized: ✓

Reported: 01/17/95

Sample No: MW-3-20

Analytical
Chemists &
Consultants

QC Report No: J442-Landau Associates, Incorporated

Project: 275002.23

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Paxton Sales

Date Sampled: 01/06/95

Date Received: 01/11/95

Date extracted: 01/13/95

Date analyzed: 01/16/95

Instrument: FINN8

GPC Cleanup: NO

Sample Amount: 28.6 g-dry-wt

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: 5.7%

pH: 6.7

| CAS Number | Analyte | ug/kg |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 350 U |
| 84-66-2 | Diethylphthalate | 70 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 70 U |
| 86-73-7 | Fluorene | 70 U |
| 100-01-6 | 4-Nitroaniline | 350 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 700 U |
| 86-30-6 | N-Nitrosodiphenylamine | 70 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 70 U |
| 118-74-1 | Hexachlorobenzene | 70 U |
| 87-86-5 | Pentachlorophenol | 350 U |
| 85-01-8 | Phenanthrene | 70 U |
| 86-74-8 | Carbazole | 70 U |
| 120-12-7 | Anthracene | 70 U |
| 84-74-2 | Di-n-Butylphthalate | 70 U |
| 206-44-0 | Fluoranthene | 70 U |
| 129-00-0 | Pyrene | 70 U |
| 85-68-7 | Butylbenzylphthalate | 70 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 350 U |
| 56-55-3 | Benzo(a)anthracene | 70 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 70 U |
| 218-01-9 | Chrysene | 70 U |
| 117-84-0 | Di-n-Octyl phthalate | 70 U |
| 205-99-2 | Benzo(b)fluoranthene | 70 U |
| 207-08-9 | Benzo(k)fluoranthene | 70 U |
| 50-32-8 | Benzo(a)pyrene | 70 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 70 U |
| 53-70-3 | Dibenz(a,h)anthracene | 70 U |
| 191-24-2 | Benzo(g,h,i)perylene | 70 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 69.8% | d5-Phenol | 60.5% |
| 2-Fluorobiphenyl | 75.2% | 2-Fluorophenol | 55.6% |
| d14-p-Terphenyl | 90.2% | 2,4,6-Tribromophenol | 68.2% |
| d4-1,2-Dichlorobenzene | 71.3% | d4-2-Chlorophenol | 64.7% |



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(206) 621-7523 (FAX)

ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Lab Sample ID: J442SB
LIMS ID: 95-294
Matrix: Soil

QC Report No: J442-Landau Associates, Incorporated
Project: 275002.23
Paxton Sales

Data Release Authorized: *MH*
Reported: 01/17/95

| LABORATORY CONTROL SAMPLE CONSTITUENT | SPIKE VALUE | SPIKE ADDED | % RECOVERY |
|--|----------------|----------------|---------------|
| Phenol | 1770 | 2500 | 70.8% |
| 2-Chlorophenol | 1620 | 2500 | 64.8% |
| 1,4-Dichlorobenzene | 1200 | 1670 | 72.0% |
| N-Nitroso-Di-N-Propylamine | 1250 | 1670 | 75.0% |
| 1,2,4-Trichlorobenzene | 1330 | 1670 | 79.8% |
| 4-Chloro-3-methylphenol | 2270 | 2500 | 90.8% |
| Acenaphthene | 1090 | 1670 | 65.4% |
| 4-Nitrophenol | 1970 | 2500 | 78.8% |
| 2,4-Dinitrotoluene | 1360 | 1670 | 81.6% |
| Pentachlorophenol | 1480 | 2500 | 59.2% |
| Pyrene | 1580 | 1670 | 94.8% |

Spike Blank Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 65.0% | d5-Phenol | 62.7% |
| 2-Fluorobiphenyl | 69.7% | 2-Fluorophenol | 55.1% |
| d14-p-Terphenyl | 85.6% | 2,4,6-Tribromophenol | 56.8% |
| d4-1,2-Dichlorobenzene | 75.7% | d4-2-Chlorophenol | 64.3% |

Values reported in ug/kg



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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Final Report
Laboratory Analysis of Conventional Parameters

Lab Sample ID: J442C
LIMS ID: 95-293
Matrix: Soil
Data Release Authorized: *MJG*
Reported: 01/27/95 *1-27*

QC Report No: J442-Landau Associates, Incorporated
Project: 275002.23
Paxton Sales
Date Sampled: 01/06/95
Date Received: 01/11/95

| <u>Analyte</u> | <u>Analysis Date</u> | <u>Method</u> | <u>Dilution Factor</u> | <u>RL</u> | <u>Units</u> | <u>Result</u> |
|----------------|----------------------|-------------------------|------------------------|-----------|--------------|---------------|
| Total Solids | 01/11/95 | EPA 160.3 SM 2540 B | | 0.01 | Percent | 92.9 |
| Total Cyanide | 01/18/95 | EPA 335.2 SM4500CN-C | | 0.20 | mg/kg | < 0.20 U |

RL Analytical reporting limit
U Undetected at reported detection limit

Report for J442 received 01/11/95



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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

QA Report - Replicate Analysis

Matrix: Soil

QC Report No: J442-Landau Associates, Incorporated

Project: 275002.23

Paxton Sales

Date Received: 01/11/95

Data Release Authorized: *Mof*
)-27

**DUPLICATE ANALYSIS RESULTS
CONVENTIONALS**

| Constituent | Units | Sample Value | Duplicate Value | RPD |
|---|--------------|---------------------|------------------------|------------|
| ARI ID: 95-293, J442 C Client Sample ID: MW-3-19.8 | | | | |
| Total Solids | Percent | 92.9 | 93.2 | 0.3% |
| Total Cyanide | mg/kg | < 0.20 U | < 0.20 U | NA |



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QA Report - Matrix Spike Analysis

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Matrix: Soil QC Report No: J442-Landau Associates, Incorporated
 Project: 275002.23
 Paxton Sales
 Date Received: 01/11/95

Data Release Authorized: *MOP*
1-27

MATRIX SPIKE QA/QC REPORT
CONVENTIONALS

| <u>Constituent</u> | <u>Units</u> | <u>Sample Value</u> | <u>Spike Value</u> | <u>Spike Added</u> | <u>Recovery</u> |
|--|--------------|---------------------|--------------------|--------------------|-----------------|
| ARI ID: 95-293, J442 C Client Sample ID: MW-3-19.8 | | | | | |
| Total Cyanide | mg/kg | < 0.2 | 95.9 | 103 | 93.1% |



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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

QC Report No: J442-Landau Associates, Incorporated (206) 621-7523 (FAX)

QA Report - Method Blank Analysis

Matrix: Soil

Project: 275002.23
Paxton Sales
Date Received: 01/11/95

Data Release Authorized: *Mof*
1-27

METHOD BLANK RESULTS
CONVENTIONALS

| <u>Analysis Date</u> | <u>Constituent</u> | <u>Units</u> | <u>Result</u> |
|--------------------------|--------------------|--------------|---------------|
| Method Blank 01/11/95 | Total Solids | mg residue | < 1 U |
| Method Blank 01/18/95 | Total Cyanide | mg/L | < 0.004 U |



**ANALYTICAL
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QA Report - Standard Reference Material Analysis

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

QC Report No: J442-Landau Associates, Incorporated
Project: 275002.23
Paxton Sales

Data Release Authorized: *MOP*
1-27

Date Received: 01/11/95

**STANDARD REFERENCE MATERIAL ANALYSIS
CONVENTIONALS**

| Constituent | Units | Value | True Value | Recovery |
|---------------------|----------|-------|------------|----------|
| SPEX QC Lot#6-153AS | | | | |
| Total Cyanide | mg/L | 0.193 | 0.200 | 96.5% |
| Date analyzed: | 01/18/95 | | | |

EXPLANATION OF INORGANIC DATA REPORT CODES

The columns labeled 'PREP', 'C', and 'M' contain important information about your analyses. The codes are defined below.

PREPARATION CODES

These 3-letter codes describe methods used to prepare samples for analysis:

| | |
|-----|--|
| AEN | USEPA, Metals in air filters |
| AHM | ARI, Mercury in air filters |
| AHN | ARI, Metals in air filters |
| ANN | NIOSH 7300, Metals in air filters |
| CAN | AOAC (1984) 25.024, Metals in earthenware |
| DE6 | EPA 600/4-79-020 218.5, Cr(VI) in water |
| DMM | DMN followed by TMM, Dissolved mercury |
| DMN | Filtered through .45u filter, Dissolved metals |
| EW6 | EWN followed by DE6 |
| EWM | EWN followed by TMM |
| VN | USEPA SW-846 1310, EP Toxicity |
| AP | ARI, Metals in tissue (HNO ₃ /HClO ₄) |
| FPP | PSEP, Metals in tissue (HNO ₃ /HClO ₄) |
| FRM | Journal, Mercury in tissue |
| FRN | Journal, Metals in tissue (HNO ₃ /H ₂ O ₂) |
| KRN | ARI, Concentration by coprecipitation |
| LEM | USEPA 1311, TCLP followed by TMM |
| LEN | USEPA 1311, TCLP Extraction |
| THM | ARI, Mercury in miscellaneous materials |
| MHN | ARI, Metals in miscellaneous materials |
| OAM | ARI, Mercury in oil, grease or tar |
| OAN | ARI, Metals in oil, grease or tar |
| PHM | ARI, Mercury in wipes |
| PHN | ARI, Metals in wipes |
| RCC | USEPA CLP, Water digestion, HCl matrix |
| RCN | USEPA CLP, Water digestion, HNO ₃ matrix |
| REC | EPA 600/4-79-020 4.1.4, HCl matrix |
| I | EPA 600/4-79-020 200.7 and 9.4 |
| EN | EPA 600/4-79-020 4.1.4, HNO ₃ matrix |

| | |
|-----|--|
| RMA | EPA 600/4-79-020 206.2 |
| RWC | USEPA SW-846 3005 |
| SCC | USEPA CLP, Soil digestion, HCl matrix |
| SCM | USEPA CLP, Mercury in soil |
| SCN | USEPA CLP, Soil digestion, HNO ₃ matrix |
| SEM | EPA 600/4-79-020 245.5, Mercury in soil |
| SHF | ARI, Metals in soil, HF digestion |
| SMN | Agronomy, Metals in soil, Water extract |
| SMM | SMN followed by DMM, Dissolved mercury |
| SPM | USEPA 1312, SPLP extraction followed by TMM |
| SPN | USEPA 1312, SPLP Extraction |
| SSS | Standard Methods 302C, Ti in soil |
| SW6 | USEPA SW-846 3060, Cr(VI) in soil |
| SWC | USEPA SW-846 3050, HCl matrix |
| SWN | USEPA SW-846 3050, HNO ₃ matrix |
| SZF | PSEP/PSDDA, Microwave, Total acid digestion |
| TEC | EPA 600/4-79-020 4.1.3, HCl matrix |
| TEG | EPA 600/4-79-020 272.1, Silver in water |
| TEI | EPA 600/4-79-020 200.7 and 9.3 |
| TEN | EPA 600/4-79-020 4.1.3, HNO ₃ matrix |
| THG | ARI, Silver in photographic solutions |
| TMM | EPA 600/4-79-020 245.1, Mercury in water |
| TSC | Standard Methods 302C, Sb/Sn in water |
| TSN | Standard Methods 302D |
| TSS | Standard Methods 302E, Ti in water |
| TWC | USEPA SW-846 3010, HCl matrix |
| TWG | USEPA SW-846 7760, Silver in water |
| TWN | USEPA SW-846 3020, HNO ₃ matrix |
| WMN | EPA 600/4-79-020, Preserved, undigested water |

CONCENTRATION CODES

These codes are used to qualify reported concentrations:

U No analyte was detected. The reported value is the lower limit of detection.

METHOD CODES

These codes signify the instrumental technique used for analysis:

| | |
|-----|---|
| CVA | Cold Vapor Atomic Absorption Spectrophotometry |
| FA | Flame Atomic Absorption Spectrophotometry |
| GFA | Graphite Furnace Atomic Absorption Spectrophotometry |
| ICP | Inductively Coupled Plasma Atomic Emission Spectrometry |



ANALYTICAL
RESOURCES
INCORPORATED

ARI job number: J442

ARI Sample number: B

Client: Landau Associates

Contact: Bill Evans

Matrix: Soil

% Solids: 93.39

ID number: MW-3-14.8

Project: 275002.23

Description:

Sampled: 01/06/95

Received: 01/11/95

Released by: B.C.

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|----------------|---|------|------|-----|----------|
| 7440-36-0 | Antimony | 0.1 mg/kg-dry | U | 0.1 | SWC | GFA | 01/17/95 |
| 7440-38-2 | Arsenic | 3.5 mg/kg-dry | | 0.5 | SWN | GFA | 01/20/95 |
| 7440-41-7 | Beryllium | 0.2 mg/kg-dry | | 0.1 | SWC | ICP | 01/13/95 |
| 7440-43-9 | Cadmium | 0.2 mg/kg-dry | | 0.2 | SWC | ICP | 01/13/95 |
| 7440-47-3 | Chromium | 14.7 mg/kg-dry | | 0.5 | SWC | ICP | 01/13/95 |
| 7440-50-8 | Copper | 27.8 mg/kg-dry | | 0.2 | SWC | ICP | 01/13/95 |
| 7439-92-1 | Lead | 3.6 mg/kg-dry | | 0.2 | SWN | GFA | 01/18/95 |
| 7439-97-6 | Mercury | 0.04 mg/kg-dry | U | 0.04 | SCM | CVA | 01/13/95 |
| 7440-02-0 | Nickel | 11 mg/kg-dry | | 1 | SWC | ICP | 01/13/95 |
| 7782-49-2 | Selenium | 0.2 mg/kg-dry | | 0.1 | SWN | GFA | 01/24/95 |
| 7440-22-4 | Silver | 0.3 mg/kg-dry | U | 0.3 | SWC | ICP | 01/13/95 |
| 7440-28-0 | Thallium | 0.2 mg/kg-dry | | 0.1 | SWN | GFA | 01/19/95 |
| 7440-66-6 | Zinc | 49.1 mg/kg-dry | | 0.4 | SWC | ICP | 01/13/95 |



**ANALYTICAL
RESOURCES
INCORPORATED**

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

ARI job number: J442

ARI Sample number: MB

Client: Landau Associates

Contact: Bill Evans

Matrix: Soil

% Solids: 100.0

ID number:

Project: 275002.23

Description: Method Blank

Sampled: / /

Received: / /

Released by: [Signature]

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|----------------|---|------|------|-----|----------|
| 7440-36-0 | Antimony | 0.1 mg/kg-dry | U | 0.1 | SWC | GFA | 01/17/95 |
| 7440-38-2 | Arsenic | 0.1 mg/kg-dry | U | 0.1 | SWN | GFA | 01/17/95 |
| 7440-41-7 | Beryllium | 0.1 mg/kg-dry | U | 0.1 | SWC | ICP | 01/13/95 |
| 7440-43-9 | Cadmium | 0.2 mg/kg-dry | U | 0.2 | SWC | ICP | 01/13/95 |
| 7440-47-3 | Chromium | 0.5 mg/kg-dry | U | 0.5 | SWC | ICP | 01/13/95 |
| 7440-50-8 | Copper | 0.2 mg/kg-dry | U | 0.2 | SWC | ICP | 01/13/95 |
| 7439-92-1 | Lead | 0.1 mg/kg-dry | U | 0.1 | SWN | GFA | 01/18/95 |
| 7439-97-6 | Mercury | 0.05 mg/kg-dry | U | 0.05 | SCM | CVA | 01/13/95 |
| 7440-02-0 | Nickel | 1 mg/kg-dry | U | 1 | SWC | ICP | 01/13/95 |
| 7782-49-2 | Selenium | 0.1 mg/kg-dry | U | 0.1 | SWN | GFA | 01/18/95 |
| 7440-22-4 | Silver | 0.3 mg/kg-dry | U | 0.3 | SWC | ICP | 01/13/95 |
| 7440-28-0 | Thallium | 0.1 mg/kg-dry | U | 0.1 | SWN | GFA | 01/19/95 |
| 7440-66-6 | Zinc | 0.6 mg/kg-dry | | 0.4 | SWC | ICP | 01/13/95 |



**ANALYTICAL
RESOURCES
INCORPORATED**

ARI job number: J442
ARI Sample number: MBSPK
Client: Landau Associates
Contact: Bill Evans
Matrix: Soil

% Solids: 100.0

ID number:
Project: 275002.23
Description: Method Blank Spike
Sampled: / /
Received: / /

Released by: [Signature]

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Consultants

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Seattle, WA 98109-5187
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(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|----------------|---|------|------|-----|----------|
| 7440-36-0 | Antimony | 10.3 mg/kg-dry | | 0.5 | SWC | GFA | 01/17/95 |
| 7440-38-2 | Arsenic | 10.6 mg/kg-dry | | 0.5 | SWN | GFA | 01/17/95 |
| 7440-41-7 | Beryllium | 4.7 mg/kg-dry | | 0.1 | SWC | ICP | 01/13/95 |
| 7440-43-9 | Cadmium | 9.7 mg/kg-dry | | 0.2 | SWC | ICP | 01/13/95 |
| 7440-47-3 | Chromium | 25.0 mg/kg-dry | | 0.5 | SWC | ICP | 01/13/95 |
| 7440-50-8 | Copper | 10.0 mg/kg-dry | | 0.2 | SWC | ICP | 01/13/95 |
| 7439-92-1 | Lead | 10.8 mg/kg-dry | | 0.5 | SWN | GFA | 01/18/95 |
| 7439-97-6 | Mercury | 0.45 mg/kg-dry | | 0.05 | SCM | CVA | 01/13/95 |
| 7440-02-0 | Nickel | 50 mg/kg-dry | | 1 | SWC | ICP | 01/13/95 |
| 7782-49-2 | Selenium | 9.7 mg/kg-dry | | 0.5 | SWN | GFA | 01/18/95 |
| 7440-22-4 | Silver | 24.0 mg/kg-dry | | 0.3 | SWC | ICP | 01/13/95 |
| 7440-28-0 | Thallium | 11.4 mg/kg-dry | | 0.5 | SWN | GFA | 01/19/95 |
| 7440-66-6 | Zinc | 48.5 mg/kg-dry | | 0.4 | SWC | ICP | 01/13/95 |



**ANALYTICAL
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INCORPORATED**

Analytical
Chemists &
Consultants

Matrix Spike Quality Control Report

Client: Landau Associates
Client's sample ID:
ARI sample ID: J442 MBSPK
Units: mg/kg-dry

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

| Analyte | Meth | Sample | Matrix Spike | Spike Added | %R | Control Limit | Q |
|-----------|------|--------|--------------|-------------|-------|---------------|---|
| Antimony | GFA | 0 | 10.3 | 10.0 | 103.0 | 75-125% | |
| Arsenic | GFA | 0 | 10.6 | 10.0 | 106.0 | 75-125% | |
| Beryllium | ICP | 0 | 4.7 | 5.0 | 94.0 | 75-125% | |
| Cadmium | ICP | 0 | 9.7 | 10.0 | 97.0 | 75-125% | |
| Chromium | ICP | 0 | 25.0 | 25.0 | 100.0 | 75-125% | |
| Copper | ICP | 0 | 10.0 | 10.0 | 100.0 | 75-125% | |
| Lead | GFA | 0 | 10.8 | 10.0 | 108.0 | 75-125% | |
| Mercury | CVA | 0 | 0.45 | 0.50 | 90.0 | 75-125% | |
| Nickel | ICP | 0 | 50 | 50 | 100.0 | 75-125% | |
| Selenium | GFA | 0 | 9.7 | 10.0 | 97.0 | 75-125% | |
| Silver | ICP | 0 | 24.0 | 25.0 | 96.0 | 75-125% | |
| Thallium | GFA | 0 | 11.4 | 10.0 | 114.0 | 75-125% | |
| Zinc | ICP | 0.6 | 48.5 | 50.0 | 95.8 | 75-125% | |

%R = Percent Recovery

'Q' codes: 'N' = control limit not met
'H' = %R not applicable, sample concentration too high
'S' = Analyte not spiked



Analytical Resources, Incorporated

Analytical Chemists and Consultants

275002

cctomum

February 6, 1995

RECEIVED

Mr. Bill Evans
Landau Associates, Inc.
3600 Port of Tacoma Road
Suite 501
Tacoma, WA 98424

FEB - 8 1995
LANDAU ASSOCIATES, INC.
TACOMA

RE: Project No. 275002 (Yakima) Paxton Sales / ARI Job No. J477

Dear Mr. Evans:

Please find enclosed original results and Chain-of-Custody Record (COC NO. 2509) for the above-referenced project. Analytical Resources, Incorporated (ARI), accepted the following six soil samples on January 16, 1995:

| | | |
|------------|-----------|-----------|
| MW-2-20.3 | MW-2-20.5 | MW-2-20.2 |
| MW-2-120.2 | MW-2-20.8 | MW-2-20.4 |

ARI received the samples intact. Prior to sample analysis the pesticide analysis was omitted from the COC and the laboratory added semivolatile organic compounds, according to the December 21, 1994, Subcontractor Agreement, Supplement 275002. The laboratory analyzed the samples for volatile organic compounds by EPA method 8260, semivolatile organics by EPA method 8270, total petroleum hydrocarbons by WDOE method WTPH-HCID, metals following the EPA method 6010/7000 series, and cyanide by EPA method 335.2.

Sample analysis was routine and no analytical complications were noted for this sample delivery group.

As always, a copy of this report and all associated raw data will remain on file with ARI. If you have any questions or require additional information, please feel free to contact me at your convenience. If I am unavailable, you can leave a message on my voice mail and I will return your call as soon as possible.

Sincerely,

ANALYTICAL RESOURCES, INC.

Bryan D. Anderson

Bryan D. Anderson
Project Manager
(206)340-2866, ext. 116

enclosures
cc: file J477

BDA/dn

Chain-of-Custody Record

Date 1/16/95
Page 1 of 1

Project DAXTON SALES Job No. 275002

Client Veljinkanje, Moore & SHORE
V-Vis
VIA

Sampler's Name CHRIS KIMMEL

EXPLANATION OF INORGANIC DATA REPORT CODES

The columns labeled 'PREP', 'C', and 'M' contain important information about your analyses. The codes are defined below.

PREPARATION CODES

These 3-letter codes describe methods used to prepare samples for analysis:

| | |
|-----|--|
| AEN | USEPA, Metals in air filters |
| AHM | ARI, Mercury in air filters |
| AHN | ARI, Metals in air filters |
| ANN | NIOSH 7300, Metals in air filters |
| CAN | AOAC (1984) 25.024, Metals in earthenware |
| DE6 | EPA 600/4-79-020 218.5, Cr(VI) in water |
| DMM | DMN followed by TMM, Dissolved mercury |
| DMN | Filtered through .45u filter, Dissolved metals |
| EW6 | EWN followed by DE6 |
| EWM | EWN followed by TMM |
| I_N | USEPA SW-846 1310, EP Toxicity |
| THP | ARI, Metals in tissue (HNO ₃ /HClO ₄) |
| PP | PSEP, Metals in tissue (HNO ₃ /HClO ₄) |
| FRM | Journal, Mercury in tissue |
| FRN | Journal, Metals in tissue (HNO ₃ /H ₂ O ₂) |
| KRN | ARI, Concentration by coprecipitation |
| LEM | USEPA 1311, TCLP followed by TMM |
| LEN | USEPA 1311, TCLP Extraction |
| HM | ARI, Mercury in miscellaneous materials |
| HN | ARI, Metals in miscellaneous materials |
| OAM | ARI, Mercury in oil, grease or tar |
| OAN | ARI, Metals in oil, grease or tar |
| HM | ARI, Mercury in wipes |
| PHN | ARI, Metals in wipes |
| RCC | USEPA CLP, Water digestion, HCl matrix |
| RCN | USEPA CLP, Water digestion, HNO ₃ matrix |
| REC | EPA 600/4-79-020 4.1.4, HCl matrix |
| REI | EPA 600/4-79-020 200.7 and 9.4 |
| REN | EPA 600/4-79-020 4.1.4, HNO ₃ matrix |

| | |
|-----|--|
| RMA | EPA 600/4-79-020 206.2 |
| RWC | USEPA SW-846 3005 |
| SCC | USEPA CLP, Soil digestion, HCl matrix |
| SCM | USEPA CLP, Mercury in soil |
| SCN | USEPA CLP, Soil digestion, HNO ₃ matrix |
| SEM | EPA 600/4-79-020 245.5, Mercury in soil |
| SHF | ARI, Metals in soil, HF digestion |
| SMN | Agronomy, Metals in soil, Water extract |
| SMM | SMN followed by DMM, Dissolved mercury |
| SPM | USEPA 1312, SPLP extraction followed by TMM |
| SPN | USEPA 1312, SPLP Extraction |
| SSS | Standard Methods 302C, Ti in soil |
| SW6 | USEPA SW-846 3060, Cr(VI) in soil |
| SWC | USEPA SW-846 3050, HCl matrix |
| SWN | USEPA SW-846 3050, HNO ₃ matrix |
| SZF | PSEP/PSDDA, Microwave, Total acid digestion |
| TEC | EPA 600/4-79-020 4.1.3, HCl matrix |
| TEG | EPA 600/4-79-020 272.1, Silver in water |
| TEI | EPA 600/4-79-020 200.7 and 9.3 |
| TEN | EPA 600/4-79-020 4.1.3, HNO ₃ matrix |
| THG | ARI, Silver in photographic solutions |
| TMM | EPA 600/4-79-020 245.1, Mercury in water |
| TSC | Standard Methods 302C, Sb/Sn in water |
| TSN | Standard Methods 302D |
| TSS | Standard Methods 302E, Ti in water |
| TWC | USEPA SW-846 3010, HCl matrix |
| TWG | USEPA SW-846 7760, Silver in water |
| TWN | USEPA SW-846 3020, HNO ₃ matrix |
| WMN | EPA 600/4-79-020, Preserved, undigested water |

CONCENTRATION CODES

These codes are used to qualify reported concentrations:

J No analyte was detected. The reported value is the lower limit of detection.

METHOD CODES

These codes signify the instrumental technique used for analysis:

| | |
|-----|---|
| CVA | Cold Vapor Atomic Absorption Spectrophotometry |
| FAA | Flame Atomic Absorption Spectrophotometry |
| GFA | Graphite Furnace Atomic Absorption Spectrophotometry |
| CP | Inductively Coupled Plasma Atomic Emission Spectrometry |



ANALYTICAL
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ARI job number: J477
ARI Sample number: A
Client: Landau Associates
Contact: Bill Evans
Matrix: Soil

% Solids: 89.31

ID number: MW-2-20.3
Project: 275002
Description:
Sampled: 01/14/95
Received: 01/16/95

Released by: 

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|----------------|---|------|------|-----|----------|
| 7440-36-0 | Antimony | 0.1 mg/kg-dry | U | 0.1 | SWC | GFA | 01/24/95 |
| 7440-38-2 | Arsenic | 3.1 mg/kg-dry | | 0.5 | SWN | GFA | 01/26/95 |
| 7440-41-7 | Beryllium | 0.3 mg/kg-dry | | 0.1 | SWC | ICP | 01/19/95 |
| 7440-43-9 | Cadmium | 0.2 mg/kg-dry | U | 0.2 | SWC | ICP | 01/19/95 |
| 7440-47-3 | Chromium | 24.6 mg/kg-dry | | 0.6 | SWC | ICP | 01/19/95 |
| 7440-50-8 | Copper | 21.9 mg/kg-dry | | 0.2 | SWC | ICP | 01/19/95 |
| 7439-92-1 | Lead | 3.9 mg/kg-dry | | 0.2 | SWN | GFA | 01/20/95 |
| 7439-97-6 | Mercury | 0.04 mg/kg-dry | U | 0.04 | SCM | CVA | 01/19/95 |
| 7440-02-0 | Nickel | 16 mg/kg-dry | | 1 | SWC | ICP | 01/19/95 |
| 7782-49-2 | Selenium | 0.1 mg/kg-dry | | 0.1 | SWN | GFA | 01/24/95 |
| 7440-22-4 | Silver | 0.3 mg/kg-dry | U | 0.3 | SWC | ICP | 01/19/95 |
| 7440-28-0 | Thallium | 0.5 mg/kg-dry | U | 0.5 | SWN | GFA | 01/24/95 |
| 7440-66-6 | Zinc | 53.2 mg/kg-dry | | 0.4 | SWC | ICP | 01/19/95 |



ANALYTICAL
RESOURCES
INCORPORATED

ARI job number: J477

ARI Sample number: MB

Client: Landau Associates

Contact: Bill Evans

Matrix: Soil

% Solids: 100.0

ID number:

Project: 275002

Description: Method Blank

Sampled: / /

Received: / /

Released by: DJK

Analytical
Chemists &
Consultants

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Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|----------------|---|------|------|-----|----------|
| 7440-36-0 | Antimony | 0.1 mg/kg-dry | U | 0.1 | SWC | GFA | 01/24/95 |
| 7440-38-2 | Arsenic | 0.1 mg/kg-dry | U | 0.1 | SWN | GFA | 01/23/95 |
| 7440-41-7 | Beryllium | 0.1 mg/kg-dry | U | 0.1 | SWC | ICP | 01/19/95 |
| 7440-43-9 | Cadmium | 0.2 mg/kg-dry | U | 0.2 | SWC | ICP | 01/19/95 |
| 7440-47-3 | Chromium | 0.5 mg/kg-dry | U | 0.5 | SWC | ICP | 01/19/95 |
| 7440-50-8 | Copper | 0.2 mg/kg-dry | U | 0.2 | SWC | ICP | 01/19/95 |
| 7439-92-1 | Lead | 0.1 mg/kg-dry | U | 0.1 | SWN | GFA | 01/20/95 |
| 7439-97-6 | Mercury | 0.05 mg/kg-dry | U | 0.05 | SCM | CVA | 01/19/95 |
| 7440-02-0 | Nickel | 1 mg/kg-dry | U | 1 | SWC | ICP | 01/19/95 |
| 7782-49-2 | Selenium | 0.1 mg/kg-dry | U | 0.1 | SWN | GFA | 01/24/95 |
| 7440-22-4 | Silver | 0.3 mg/kg-dry | U | 0.3 | SWC | ICP | 01/19/95 |
| 7440-28-0 | Thallium | 0.1 mg/kg-dry | U | 0.1 | SWN | GFA | 01/24/95 |
| 7440-66-6 | Zinc | 0.4 mg/kg-dry | | 0.4 | SWC | ICP | 01/19/95 |



**ANALYTICAL
RESOURCES
INCORPORATED**

ARI job number: J477
ARI Sample number: MBSPK
Client: Landau Associates
Contact: Bill Evans
Matrix: Soil
% Solids: 100.0

ID number:
Project: 275002
Description: Method Blank Spike
Sampled: / /
Received: / /
Released by: *[Signature]*

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Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|----------------|---|------|------|-----|----------|
| 7440-36-0 | Antimony | 10.7 mg/kg-dry | | 0.5 | SWC | GFA | 01/24/95 |
| 7440-38-2 | Arsenic | 10.4 mg/kg-dry | | 0.5 | SWN | GFA | 01/23/95 |
| 7440-41-7 | Beryllium | 4.9 mg/kg-dry | | 0.1 | SWC | ICP | 01/19/95 |
| 7440-43-9 | Cadmium | 9.9 mg/kg-dry | | 0.2 | SWC | ICP | 01/19/95 |
| 7440-47-3 | Chromium | 25.5 mg/kg-dry | | 0.5 | SWC | ICP | 01/19/95 |
| 7440-50-8 | Copper | 10.0 mg/kg-dry | | 0.2 | SWC | ICP | 01/19/95 |
| 7439-92-1 | Lead | 9.5 mg/kg-dry | | 0.5 | SWN | GFA | 01/20/95 |
| 7439-97-6 | Mercury | 0.44 mg/kg-dry | | 0.05 | SCM | CVA | 01/19/95 |
| 7440-02-0 | Nickel | 50 mg/kg-dry | | 1 | SWC | ICP | 01/19/95 |
| 7782-49-2 | Selenium | 10.6 mg/kg-dry | | 0.5 | SWN | GFA | 01/24/95 |
| 7440-22-4 | Silver | 24.6 mg/kg-dry | | 0.3 | SWC | ICP | 01/19/95 |
| 7440-28-0 | Thallium | 10.1 mg/kg-dry | | 0.5 | SWN | GFA | 01/24/95 |
| 7440-66-6 | Zinc | 50.3 mg/kg-dry | | 0.4 | SWC | ICP | 01/19/95 |



ANALYTICAL
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Matrix Spike Quality Control Report

Client: Landau Associates
Client's sample ID:
ARI sample ID: J477 MBSPK
Units: mg/kg-dry

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

| Analyte | Meth | Sample | Matrix Spike | Spike Added | %R | Control Limit | Q |
|-----------|------|--------|--------------|-------------|-------|---------------|---|
| Antimony | GFA | 0 | 10.7 | 10.0 | 107.0 | 75-125% | |
| Arsenic | GFA | 0 | 10.4 | 10.0 | 104.0 | 75-125% | |
| Beryllium | ICP | 0 | 4.9 | 5.0 | 98.0 | 75-125% | |
| Cadmium | ICP | 0 | 9.9 | 10.0 | 99.0 | 75-125% | |
| Chromium | ICP | 0 | 25.5 | 25.0 | 102.0 | 75-125% | |
| Copper | ICP | 0 | 10.0 | 10.0 | 100.0 | 75-125% | |
| Lead | GFA | 0 | 9.5 | 10.0 | 95.0 | 75-125% | |
| Mercury | CVA | 0 | 0.44 | 0.50 | 88.0 | 75-125% | |
| Nickel | ICP | 0 | 51 | 50 | 102.0 | 75-125% | |
| Selenium | GFA | 0 | 10.6 | 10.0 | 106.0 | 75-125% | |
| Silver | ICP | 0 | 24.6 | 25.0 | 98.4 | 75-125% | |
| Thallium | GFA | 0 | 10.1 | 10.0 | 101.0 | 75-125% | |
| Zinc | ICP | 0.4 | 50.3 | 50.0 | 99.8 | 75-125% | |

%R = Percent Recovery

'Q' codes: 'N' = control limit not met
 'H' = %R not applicable, sample concentration too high
 'S' = Analyte not spiked



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Final Report
Laboratory Analysis of Conventional Parameters

Lab Sample ID: J477B
LIMS ID: 95-450
Matrix: Soil
Data Release Authorized: *mof*
Reported: 01/30/95 *1-38*

QC Report No: J477-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/14/95
Date Received: 01/16/95

| Analyte | Analysis Date | Method | Dilution Factor | RL | Units | Result |
|---------------|---------------|-------------------------|-----------------|----|---------|----------|
| Total Solids | 01/18/95 | EPA 160.3 SM 2540 B | 0.01 | | Percent | 88.9 |
| Total Cyanide | 01/24/95 | EPA 335.2 SM4500CN-C | 0.11 | | mg/kg | < 0.11 U |

RL Analytical reporting limit
U Undetected at reported detection limit

Report for J477 received 01/16/95



**ANALYTICAL
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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

QA Report - Replicate Analysis

Matrix: Soil

QC Report No: J477-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Received: 01/16/95

Data Release Authorized: *Mof*
1-30

**DUPLICATE ANALYSIS RESULTS
CONVENTIONALS**

| Constituent | Units | Sample Value | Duplicate Value | RPD |
|---|--------------|---------------------|------------------------|------------|
| ARI ID: 95-450, J477 B Client Sample ID: MW-2-20.5 | | | | |
| Total Solids | Percent | 88.9 | 86.7 | 2.5% |
| Total Cyanide | mg/kg | < 0.11 U | < 0.11 U | NA |



**ANALYTICAL
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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

QA Report - Matrix Spike Analysis

Matrix: Soil

QC Report No: J477-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/16/95

Data Release Authorized: *Mop*
1.36

**MATRIX SPIKE QA/QC REPORT
CONVENTIONALS**

| <u>Constituent</u> | <u>Units</u> | <u>Sample Value</u> | <u>Spike Value</u> | <u>Spike Added</u> | <u>Recovery</u> |
|---|--------------|---------------------|--------------------|--------------------|-----------------|
| ARI ID: 95-450, J477 B Client Sample ID: MW-2-20.5 | | | | | |
| Total Cyanide | mg/kg | < 0.11 | 43.7 | 48.6 | 90.0% |



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Seattle, WA 98109-5187
(206) 621-6490

~~(206) 621-7523 (FAX)~~

QA Report - Method Blank Analysis

Matrix: Soil

QC Report No: J477-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Received: 01/16/95

Data Release Authorized: *Mof*
1-30

METHOD BLANK RESULTS
CONVENTIONALS

| <u>Analysis Date</u> | <u>Constituent</u> | <u>Units</u> | <u>Result</u> |
|--------------------------|--------------------|--------------|---------------|
| Method Blank 01/18/95 | Total Solids | mg residue | < 1 U |
| Method Blank 01/24/95 | Total Cyanide | mg/L | < 0.004 U |



**ANALYTICAL
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QA Report - Standard Reference Material Analysis

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

QC Report No: J477-Landau Associates, Incorporated
Project: 275002
Paxton Sales

Data Release Authorized: *Map*
1-30

Date Received: 01/16/95

**STANDARD REFERENCE MATERIAL ANALYSIS
CONVENTIONALS**

| Constituent | Units | Value | True Value | Recovery |
|---------------------|----------|-------|------------|----------|
| SPEX QC Lot#6-153AS | | | | |
| Total Cyanide | mg/L | 0.185 | 0.200 | 92.5% |
| Date analyzed: | 01/24/95 | | | |



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ORGANIC COMPOUND DATA REPORTING QUALIFIERS

- U** Indicates the compound was analyzed for, but not detected at the given detection limit.
- J** Indicates an estimated value when the result is less than the calculated detection limit.
- D** Indicates the surrogate/spike(s) was not detected, due to dilution of extract.
- NR** Indicates the surrogate recovery cannot be reported due to matrix interference.
- E** Indicates a value above the linear range of the detector.
Sample dilution required.
- S** Indicates no value reported due to saturation of the detector.
Dilution required.
- Y** Indicates a raised detection limit due to matrix interferences.
- NA** Indicates compound was not analyzed.
- M** Indicates an estimated value of analyte found and confirmed by analyst but with low spectral match.
- B** Indicates compound was found in the associated method blank.

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



Sample No: MW-2-20.2

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Lab Sample ID: J477C

QC Report No: J477-Landau Associates, Incorporated

LIMS ID: 95-451

Project: 275002

Matrix: Soil

Paxton Sales

Data Release Authorized: *BMB*

Date Sampled: 01/14/95

Reported: 02/01/95

Date Received: 01/16/95

Instrument: FINN5

Sample Amount: 4.65 g dry Wt

Date Analyzed: 01/26/95

Percent Moisture: 10.6%

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.2 U |
| 75-01-4 | Vinyl Chloride | 2.2 U |
| 75-00-3 | Chloroethane | 2.2 U |
| 75-09-2 | Methylene Chloride | 2.2 U |
| 75-35-4 | 1,1-Dichloroethene | 1.1 U |
| 75-34-3 | 1,1-Dichloroethane | 1.1 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.1 U |
| 67-66-3 | Chloroform | 1.1 U |
| 107-06-2 | 1,2-Dichloroethane | 1.1 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.1 U |
| 56-23-5 | Carbon Tetrachloride | 1.1 U |
| 75-27-4 | Bromodichloromethane | 1.1 U |
| 78-87-5 | 1,2-Dichloropropane | 1.1 U |
| 79-01-6 | Trichloroethene | 1.1 U |
| 124-48-1 | Dibromochloromethane | 1.1 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.1 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.1 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.4 U |
| 75-25-2 | Bromoform | 1.1 U |
| 127-18-4 | Tetrachloroethene | 1.1 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.1 U |
| 108-90-7 | Chlorobenzene | 1.1 U |
| 75-69-4 | Trichlorofluoromethane | 2.2 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.1 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.1 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.1 U |
| 74-96-4 | Bromoethane | 2.2 U |
| 74-95-3 | Dibromomethane | 1.1 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.1 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.1 U |
| 108-86-1 | Bromobenzene | 1.1 U |
| 95-49-8 | 2-Chlorotoluene | 1.1 U |
| 106-43-4 | 4-Chlorotoluene | 1.1 U |

Volatile Surrogate Recovery

| | |
|------------------------|------|
| d4-1,2-Dichloroethane | 107% |
| d8-Toluene | 107% |
| Bromofluorobenzene | 101% |
| d4-1,2-Dichlorobenzene | 103% |

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



Sample No: MW-2-120.2

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Lab Sample ID: J477D

QC Report No: J477-Landau Associates, Incorporated

LIMS ID: 95-452

Project: 275002

Matrix: Soil

Paxton Sales

Data Release Authorized: *bnd*

Date Sampled: 01/14/95

Reported: 02/01/95

Date Received: 01/16/95

Instrument: FINN5

Sample Amount: 4.52 g dry Wt

Date Analyzed: 01/26/95

Percent Moisture: 10.1%

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.2 U |
| 75-01-4 | Vinyl Chloride | 2.2 U |
| 75-00-3 | Chloroethane | 2.2 U |
| 75-09-2 | Methylene Chloride | 2.2 U |
| 75-35-4 | 1,1-Dichloroethene | 1.1 U |
| 75-34-3 | 1,1-Dichloroethane | 1.1 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.1 U |
| 67-66-3 | Chloroform | 1.1 U |
| 107-06-2 | 1,2-Dichloroethane | 1.1 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.1 U |
| 56-23-5 | Carbon Tetrachloride | 1.1 U |
| 75-27-4 | Bromodichloromethane | 1.1 U |
| 78-87-5 | 1,2-Dichloropropane | 1.1 U |
| 79-01-6 | Trichloroethene | 1.1 U |
| 124-48-1 | Dibromochloromethane | 1.1 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.1 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.1 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.5 U |
| 75-25-2 | Bromoform | 1.1 U |
| 127-18-4 | Tetrachloroethene | 1.1 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.1 U |
| 108-90-7 | Chlorobenzene | 1.1 U |
| 75-69-4 | Trichlorofluoromethane | 2.2 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.1 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.1 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.1 U |
| 74-96-4 | Bromoethane | 2.2 U |
| 74-95-3 | Dibromomethane | 1.1 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.1 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.1 U |
| 108-86-1 | Bromobenzene | 1.1 U |
| 95-49-8 | 2-Chlorotoluene | 1.1 U |
| 106-43-4 | 4-Chlorotoluene | 1.1 U |

Volatile Surrogate Recovery

| | |
|------------------------|------|
| d4-1,2-Dichloroethane | 101% |
| d8-Toluene | 110% |
| Bromofluorobenzene | 101% |
| d4-1,2-Dichlorobenzene | 101% |

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



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| | |
|-------------------------------------|--|
| Lab Sample ID: 012695MB | QC Report No: J477-Landau Associates, Incorporated |
| LIMS ID: 95-451 | Project: 275002 |
| Matrix: Soil | Paxton Sales |
| Data Release Authorized: <i>EWB</i> | Date Sampled: NA |
| Reported: 02/01/95 | Date Received: NA |
| Instrument: FINN5 | Sample Amount: 5.00 g dry Wt Equiv |
| Date Analyzed: 01/26/95 | Percent Moisture: NA |

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.0 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.0 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-83-9 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|------|
| d4-1,2-Dichloroethane | 113% |
| d8-Toluene | 102% |
| Bromofluorobenzene | 101% |
| d4-1,2-Dichlorobenzene | 106% |



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ORGANICS ANALYSIS DATA SHEET
Volatile by GC/MS
Page 1 of 1

Lab Sample ID: J477SB
LIMS ID: 95-451
Matrix: Soil
Data Release Authorized: *BWB*
Reported: 02/01/95
Date Analyzed: 01/26/95

QC Report No: J477-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/16/95

| LABORATORY CONTROL SAMPLE CONSTITUENT | SPIKE VALUE | SPIKE AMT | % RECOVERY |
|--|----------------|--------------|---------------|
| Chloromethane | 57.2 | 50.0 | 114% |
| Vinyl Chloride | 63.1 | 50.0 | 126% |
| Chloroethane | 52.7 | 50.0 | 105% |
| Methylene Chloride | 50.5 | 50.0 | 101% |
| 1,1-Dichloroethene | 55.2 | 50.0 | 110% |
| 1,1-Dichloroethane | 50.8 | 50.0 | 102% |
| trans-1,2-Dichloroethene | 51.2 | 50.0 | 102% |
| Chloroform | 50.5 | 50.0 | 101% |
| 1,2-Dichloroethane | 53.0 | 50.0 | 106% |
| 1,1,1-Trichloroethane | 50.2 | 50.0 | 100% |
| Carbon Tetrachloride | 53.3 | 50.0 | 107% |
| Bromodichloromethane | 51.0 | 50.0 | 102% |
| 1,2-Dichloropropane | 52.8 | 50.0 | 106% |
| Trichloroethene | 51.4 | 50.0 | 103% |
| Dibromochloromethane | 49.2 | 50.0 | 98.4% |
| 1,1,2-Trichloroethane | 50.7 | 50.0 | 101% |
| trans-1,3-Dichloropropene | 50.9 | 50.0 | 102% |
| 2-Chloroethylvinylether | 54.2 | 50.0 | 108% |
| Bromoform | 53.3 | 50.0 | 107% |
| Tetrachloroethene | 49.5 | 50.0 | 99.0% |
| 1,1,2,2-Tetrachloroethane | 52.2 | 50.0 | 104% |
| Chlorobenzene | 49.7 | 50.0 | 99.4% |
| Trichlorofluoromethane | 49.3 | 50.0 | 98.6% |
| 1,2-Dichlorobenzene | 51.0 | 50.0 | 102% |
| 1,3-Dichlorobenzene | 51.8 | 50.0 | 104% |
| 1,4-Dichlorobenzene | 48.1 | 50.0 | 96.2% |
| Bromoethane | 57.2 | 50.0 | 114% |
| Dibromomethane | 45.0 | 50.0 | 90.0% |
| 1,1,1,2-Tetrachloroethane | 48.7 | 50.0 | 97.4% |
| 1,2,3-Trichloropropane | 51.0 | 50.0 | 102% |
| Bromobenzene | 50.3 | 50.0 | 101% |
| 2-Chlorotoluene | 47.8 | 50.0 | 95.6% |
| 4-Chlorotoluene | 53.7 | 50.0 | 107% |

Spike Blank Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 107.% |
| d8-Toluene | 102.% |
| Bromofluorobenzene | 109.% |
| d4-1,2-Dichlorobenzene | 109.% |

Reported in ug/kg-dry-Wt



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TOTAL PETROLEUM HYDROCARBONS
WA HCID Method by GC/FID

Matrix: Soil

Project: 275002

Paxton Sales

Data Release Authorized:
Reported: 02/06/95

Date Received: 01/16/95

| Lab ID | Client Sample ID | Date Analyzed | Dilution Factor | Gas Range | Diesel Range | Oil Range | Surrogate Recovery |
|---------------|------------------|---------------|-----------------|-----------|--------------|-----------|--------------------|
| 95-453-0116MB | Method Blank | 01/16/95 | 1:1 | 20 U | 25 U | 50 U | 92.5% |
| 95-453-J477E | MW-2-20.8 | 01/18/95 | 1:1 | 20 U | 25 U | 50 U | 106% |

Values reported in ppm (mg/kg) on a dry weight basis.

Surrogate is Methyl Arachidate.

Gas value based on total peaks in the range from Toluene to C12.

Diesel value based on the total peaks in the range from C12 to C24.

Oil value based on the total peaks in the range from C24 to C32.

Data Qualifiers

- U Compound not detected at the given detection limit.
- X Value detected above linear range of instrument. Dilution required.
- J Indicates an estimated value below the calculated detection limit.
- S No value reported due to saturation of the detector. Dilution required.
- D Indicates the surrogate was not detected because of dilution of the extract.
- C Indicates a probable value which cannot be confirmed due to matrix interference.
- NR Indicates no recovery due to matrix interference and/or dilution.



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TOTAL PETROLEUM HYDROCARBONS
WA HCID Method by GC/FID

Lab Sample ID: J477SB
LIMS ID: 95-453
Matrix: Soil

QC Report No: J477-Landau Associates, Incorporated
Project: 275002
Paxton Sales

Data Release Authorized:
Reported: 01/23/95

Date extracted: 01/16/95

LABORATORY CONTROL SAMPLE RECOVERY REPORT

| CONSTITUENT | SPIKE VALUE | SPIKE ADDED | % RECOVERY |
|---------------------------|----------------|----------------|---------------|
| LABORATORY CONTROL SAMPLE | | | |
| Diesel Range | 539 | 500 | 108% |

HCID Surrogate Recovery

| | | |
|-------------|------------------|------|
| Spike Blank | Methylarachidate | 104% |
|-------------|------------------|------|

Values reported in parts per million (mg/kg)

HCID SPIKE CONTROL LIMITS

| | |
|------------------|---------|
| Percent Recovery | 50-150% |
| Duplicate RPD | <50% |

Advisory QA Limits



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ORGANIC COMPOUND DATA REPORTING QUALIFIERS

- U** Indicates the compound was analyzed for, but not detected at the given detection limit.
- J** Indicates an estimated value when the result is less than the calculated detection limit.
- D** Indicates the surrogate/spike(s) was not detected, due to dilution of extract.
- NR** Indicates the surrogate recovery cannot be reported due to matrix interference.
- E** Indicates a value above the linear range of the detector.
Sample dilution required.
- S** Indicates no value reported due to saturation of the detector.
Dilution required.
- Y** Indicates a raised detection limit due to matrix interferences.
- NA** Indicates compound was not analyzed.
- M** Indicates an estimated value of analyte found and confirmed by analyst but with low spectral match.
- B** Indicates compound was found in the associated method blank.



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ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 1 of 2

Lab Sample ID: J477F

LIMS ID: 95-454

Matrix: Soil

Data Release Authorized: MHA

Reported: 01/19/95

Sample No: MW-2-20.4

QC Report No: J477-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Sampled: 01/14/95

Date Received: 01/16/95

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Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN4

GPC Cleanup: NO

Sample Amount: 27.4 g-dry-wt

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: 8.9%

pH: 7.9

| CAS Number | Analyte | ug/kg |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 150 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 150 U |
| 95-57-8 | 2-Chlorophenol | 73 U |
| 541-73-1 | 1,3-Dichlorobenzene | 73 U |
| 106-46-7 | 1,4-Dichlorobenzene | 73 U |
| 100-51-6 | Benzyl Alcohol | 370 U |
| 95-50-1 | 1,2-Dichlorobenzene | 73 U |
| 95-48-7 | 2-Methylphenol | 150 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 73 U |
| 106-44-5 | 4-Methylphenol | 73 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 150 U |
| 67-72-1 | Hexachloroethane | 150 U |
| 98-95-3 | Nitrobenzene | 73 U |
| 78-59-1 | Isophorone | 73 U |
| 88-75-5 | 2-Nitrophenol | 370 U |
| 105-67-9 | 2,4-Dimethylphenol | 220 U |
| 65-85-0 | Benzoic Acid | 730 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 73 U |
| 120-83-2 | 2,4-Dichlorophenol | 220 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 73 U |
| 91-20-3 | Naphthalene | 220 U |
| 106-47-8 | 4-Chloroaniline | 150 U |
| 87-68-3 | Hexachlorobutadiene | 150 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 150 U |
| 91-57-6 | 2-Methylnaphthalene | 73 U |
| 77-47-4 | Hexachlorocyclopentadiene | 370 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 370 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 370 U |
| 91-58-7 | 2-Chloronaphthalene | 73 U |
| 88-74-4 | 2-Nitroaniline | 370 U |
| 131-11-3 | Dimethylphthalate | 73 U |
| 208-96-8 | Acenaphthylene | 73 U |
| 99-09-2 | 3-Nitroaniline | 440 U |
| 83-32-9 | Acenaphthene | 73 U |
| 51-28-5 | 2,4-Dinitrophenol | 730 U |
| 100-02-7 | 4-Nitrophenol | 370 U |
| 132-64-9 | Dibenzofuran | 73 U |
| 606-20-2 | 2,6-Dinitrotoluene | 370 U |



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ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2

Lab Sample ID: J477F

LIMS ID: 95-454

Matrix: Soil

Data Release Authorized: M+J

Reported: 01/19/95

Sample No: MW-2-20.4

Analytical
Chemists &
Consultants

QC Report No: J477-Landau Associates, Incorporated

Project: 275002

Paxton Sales

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Date Sampled: 01/14/95

Date Received: 01/16/95

Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN4

GPC Cleanup: NO

Sample Amount: 27.4 g-dry-wt

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: 8.9%

pH: 7.9

| CAS Number | Analyte | ug/kg |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 370 U |
| 84-66-2 | Diethylphthalate | 73 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 73 U |
| 86-73-7 | Fluorene | 73 U |
| 100-01-6 | 4-Nitroaniline | 370 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 730 U |
| 86-30-6 | N-Nitrosodiphenylamine | 73 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 73 U |
| 118-74-1 | Hexachlorobenzene | 73 U |
| 87-86-5 | Pentachlorophenol | 370 U |
| 85-01-8 | Phenanthrene | 73 U |
| 86-74-8 | Carbazole | 73 U |
| 120-12-7 | Anthracene | 73 U |
| 84-74-2 | Di-n-Butylphthalate | 73 U |
| 206-44-0 | Fluoranthene | 73 U |
| 129-00-0 | Pyrene | 73 U |
| 85-68-7 | Butylbenzylphthalate | 73 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 370 U |
| 56-55-3 | Benzo(a)anthracene | 73 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 73 U |
| 218-01-9 | Chrysene | 73 U |
| 117-84-0 | Di-n-Octyl phthalate | 73 U |
| 205-99-2 | Benzo(b)fluoranthene | 73 U |
| 207-08-9 | Benzo(k)fluoranthene | 73 U |
| 50-32-8 | Benzo(a)pyrene | 73 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 73 U |
| 53-70-3 | Dibenz(a,h)anthracene | 73 U |
| 191-24-2 | Benzo(g,h,i)perylene | 73 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 85.0% | d5-Phenol | 75.3% |
| 2-Fluorobiphenyl | 79.0% | 2-Fluorophenol | 76.0% |
| d14-p-Terphenyl | 87.4% | 2,4,6-Tribromophenol | 82.0% |
| d4-1,2-Dichlorobenzene | 83.6% | d4-2-Chlorophenol | 79.1% |



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Semivolatiles by GC/MS

Page 1 of 2

Lab Sample ID: J477MB

LIMS ID: 95-454

Matrix: Soil

Data Release Authorized: M A

Reported: 01/19/95

Sample No: Method Blank

Analytical
Chemists &
Consultants

QC Report No: J477-Landau Associates, Incorporated

Project: 275002

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Paxton Sales

Date Sampled: NA

(206) 621-6490

Date Received: NA

(206) 621-7523 (FAX)

Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN2

GPC Cleanup: NO

Sample Amount: 30.0 g-dry-wt Equiv

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: NA

pH: NA

| CAS Number | Analyte | ug/kg |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 130 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 130 U |
| 95-57-8 | 2-Chlorophenol | 67 U |
| 541-73-1 | 1,3-Dichlorobenzene | 67 U |
| 106-46-7 | 1,4-Dichlorobenzene | 67 U |
| 100-51-6 | Benzyl Alcohol | 330 U |
| 95-50-1 | 1,2-Dichlorobenzene | 67 U |
| 95-48-7 | 2-Methylphenol | 130 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 67 U |
| 106-44-5 | 4-Methylphenol | 67 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 130 U |
| 67-72-1 | Hexachloroethane | 130 U |
| 98-95-3 | Nitrobenzene | 67 U |
| 78-59-1 | Isophorone | 67 U |
| 88-75-5 | 2-Nitrophenol | 330 U |
| 105-67-9 | 2,4-Dimethylphenol | 200 U |
| 65-85-0 | Benzoic Acid | 670 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 67 U |
| 120-83-2 | 2,4-Dichlorophenol | 200 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 67 U |
| 91-20-3 | Naphthalene | 67 U |
| 106-47-8 | 4-Chloroaniline | 200 U |
| 87-68-3 | Hexachlorobutadiene | 130 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 130 U |
| 91-57-6 | 2-Methylnaphthalene | 67 U |
| 77-47-4 | Hexachlorocyclopentadiene | 330 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 330 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 330 U |
| 91-58-7 | 2-Chloronaphthalene | 67 U |
| 88-74-4 | 2-Nitroaniline | 330 U |
| 131-11-3 | Dimethylphthalate | 67 U |
| 208-96-8 | Acenaphthylene | 67 U |
| 99-09-2 | 3-Nitroaniline | 400 U |
| 83-32-9 | Acenaphthene | 67 U |
| 51-28-5 | 2,4-Dinitrophenol | 670 U |
| 100-02-7 | 4-Nitrophenol | 330 U |
| 132-64-9 | Dibenzofuran | 67 U |
| 606-20-2 | 2,6-Dinitrotoluene | 330 U |



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2

Lab Sample ID: J477MB

LIMS ID: 95-454

Matrix: Soil

Data Release Authorized: *MH*

Reported: 01/19/95

Sample No: Method Blank

Analytical
Chemists &
Consultants

QC Report No: J477-Landau Associates, Incorporated

Project: 275002

333 Ninth Ave. North
Seattle, WA 98109-5187

Paxton Sales

(206) 621-6490

Date Sampled: NA

(206) 621-7523 (FAX)

Date Received: NA

Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN2

GPC Cleanup: NO

Sample Amount: 30.0 g-dry-wt Equiv

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: NA

pH: NA

| CAS Number | Analyte | ug/kg |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 330 U |
| 84-66-2 | Diethylphthalate | 67 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 67 U |
| 86-73-7 | Fluorene | 67 U |
| 100-01-6 | 4-Nitroaniline | 330 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 670 U |
| 86-30-6 | N-Nitrosodiphenylamine | 67 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 67 U |
| 118-74-1 | Hexachlorobenzene | 67 U |
| 87-86-5 | Pentachlorophenol | 330 U |
| 85-01-8 | Phenanthrene | 67 U |
| 86-74-8 | Carbazole | 67 U |
| 120-12-7 | Anthracene | 67 U |
| 84-74-2 | Di-n-Butylphthalate | 67 U |
| 206-44-0 | Fluoranthene | 67 U |
| 129-00-0 | Pyrene | 67 U |
| 85-68-7 | Butylbenzylphthalate | 67 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 330 U |
| 56-55-3 | Benzo(a)anthracene | 67 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 67 U |
| 218-01-9 | Chrysene | 67 U |
| 117-84-0 | Di-n-Octyl phthalate | 67 U |
| 205-99-2 | Benzo(b)fluoranthene | 67 U |
| 207-08-9 | Benzo(k)fluoranthene | 67 U |
| 50-32-8 | Benzo(a)pyrene | 67 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 67 U |
| 53-70-3 | Dibenz(a,h)anthracene | 67 U |
| 191-24-2 | Benzo(g,h,i)perylene | 67 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 69.7% | d5-Phenol | 70.2% |
| 2-Fluorobiphenyl | 82.2% | 2-Fluorophenol | 82.3% |
| d14-p-Terphenyl | 94.7% | 2,4,6-Tribromophenol | 46.3% |
| d4-1,2-Dichlorobenzene | 79.2% | d4-2-Chlorophenol | 75.7% |



ANALYTICAL
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INCORPORATED

ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Analytical
Chemists &
Consultants

Lab Sample ID: J477SB
LIMS ID: 95-454
Matrix: Soil

QC Report No: J477-Landau Associates, Incorporated
Project: 275002
Paxton Sales

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Data Release Authorized: M4
Reported: 01/19/95

| LABORATORY CONTROL SAMPLE CONSTITUENT | SPIKE VALUE | SPIKE ADDED | % RECOVERY |
|--|----------------|----------------|---------------|
| Phenol | 2220 | 2500 | 88.8% |
| 2-Chlorophenol | 1980 | 2500 | 79.2% |
| 1,4-Dichlorobenzene | 1350 | 1670 | 81.0% |
| N-Nitroso-Di-N-Propylamine | 1430 | 1670 | 85.8% |
| 1,2,4-Trichlorobenzene | 1430 | 1670 | 85.8% |
| 4-Chloro-3-methylphenol | 2320 | 2500 | 92.8% |
| Acenaphthene | 1470 | 1670 | 88.2% |
| 4-Nitrophenol | 1790 | 2500 | 71.6% |
| 2,4-Dinitrotoluene | 1640 | 1670 | 98.4% |
| Pentachlorophenol | 1290 | 2500 | 51.6% |
| Pyrene | 1740 | 1670 | 104% |

Spike Blank Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 69.6% | d5-Phenol | 77.6% |
| 2-Fluorobiphenyl | 82.9% | 2-Fluorophenol | 82.0% |
| d14-p-Terphenyl | 93.2% | 2,4,6-Tribromophenol | 66.0% |
| d4-1,2-Dichlorobenzene | 74.3% | d4-2-Chlorophenol | 77.8% |

Values reported in ug/kg



Analytical Resources, Incorporated
Analytical Chemists and Consultants

275002

February 2, 1995

CC TO MUM

Mr. Bill Evans
Landau Associates, Inc.
3600 Port of Tacoma Road
Suite 501
Tacoma, WA 98424

RECEIVED

FEB - 8 1995

LANDAU ASSOCIATES, INC.
TACOMA

RE: Project No. 275002 (Yakima) Paxton Sales / ARI Job No. J468

Dear Mr. Evans:

Please find enclosed original results and Chain-of-Custody Record (COC NO. 2509) for the above-referenced project. Analytical Resources, Incorporated (ARI), accepted the fourteen soil samples on January 13, 1995. ARI received the samples intact with no discrepancies on the chain-of-custody documentation.

The laboratory analyzed the samples for volatile organic compounds by EPA method 8260, semivolatile organics by EPA method 8270, total petroleum hydrocarbons by WDOE method WTPH-HCID, metals following the EPA method 6010/7000 series, and cyanide by EPA method 335.2.

Sample analysis was routine and no analytical complications were noted for this sample delivery group.

As always, a copy of this report and all associated raw data will remain on file with ARI. If you have any questions or require additional information, please feel free to contact me at your convenience. If I am unavailable, you can leave a message on my voice mail and I will return your call as soon as possible.

Sincerely,

ANALYTICAL RESOURCES, INC.

Bryan D. Anderson

Bryan D. Anderson
Project Manager
(206)340-2866, ext. 116

enclosures
cc: file J468

BDA/dn

Chain-of-Custody Record

No. 2514

Date 1/12/95
Page 1 of 1

Project PAXTON SALES Job No. 275002

Client PAXTON

Project Location YAKIMA WA.

Sampler's Name CHRISTINE KIMMEL

Testing Parameters

| Sample No. | Date | Time | MATRIX | Location | No. of Containers | Observations/Comments |
|--|--|---|---|---|---|---|
| MWN-1-5.2' | 1/11/95 | 501L | MWN-1 | | 1 | X X |
| MWN-1-5.3' | | | | | | X X |
| MWN-1-5.1' | | | | | | X X |
| MWN-1-7.5 | | | | | | X X |
| MWN-1-10.0 | | | | | | X X |
| MWN-1-15.2 | | | | | | X X |
| MWN-1-17.6 | | | | | | X X |
| MWN-1-20.2 | | | | | | X X X X |
| MWN-1-22.3 | | | | | | X X X X |
| MWN-1-22.4 | | | | | | X X X X |
| MWN-1-25.0 | 1/12/95 | | | | | X |
| MWN-1-27.5 | | | | | | |
| MWN-1-30.0 | | | | | | |
| MWN-1-30.1 | | | | | | |
| Special Shipment/Handling or Storage Requirements | | | | | | Method of Shipment UPS GROUND |
| Relinquished by | Received by | Relinquished by | Received by | Relinquished by | Received by | Received by |
| Christine Kimmel Printed Name LANDAU ASSOCIATES Company Date 1/12/95 Time 1400 | Signature Christine Kimmel Printed Name LANDAU ASSOCIATES Company Date 1/12/95 Time 12:00 | Signature Christine Kimmel Printed Name B. A. T. Company Date 1/13/95 Time 12:00 | Signature Christine Kimmel Printed Name B. A. T. Company Date 1/13/95 Time 12:00 | Signature Christine Kimmel Printed Name B. A. T. Company Date 1/13/95 Time 12:00 | Signature Christine Kimmel Printed Name B. A. T. Company Date 1/13/95 Time 12:00 | Signature Christine Kimmel Printed Name B. A. T. Company Date 1/13/95 Time 12:00 |



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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

(206) 621-7523 (FAX)

TOTAL PETROLEUM HYDROCARBONS
WA HCID Method by GC/FID

Matrix: Soil

Project: 275002

Paxton Sales

Data Release Authorized:

Date Received: 01/13/95

Reported: 02/06/95

Allen M. Newson

| Lab ID | Client Sample ID | Date Analyzed | Dilution Factor | Gas Range | Diesel Range | Oil Range | Surrogate Recovery |
|---------------|---------------------|------------------|--------------------|--------------|-----------------|--------------|-----------------------|
| 95-411-0116MB | Method Blank | 01/16/95 | 1:1 | 20 U | 25 U | 50 U | 92.5% |
| 95-411-J468A | MW-1-5.2' | 01/17/95 | 1:1 | 20 U | 25 U | 50 U | 93.4% |
| 95-420-J468J | MW-1-22.4 | 01/17/95 | 1:1 | 20 U | 25 U | 50 U | 112% |
| 95-421-J468K | MW-1-25.0 | 01/17/95 | 1:1 | 20 U | 25 U | 50 U | 100% |
| 95-422-J468L | MW-1-27.5 | 01/17/95 | 1:1 | 20 U | 25 U | 50 U | 98.7% |

Values reported in ppm (mg/kg) on a dry weight basis.

Surrogate is Methyl Arachidate.

Gas value based on total peaks in the range from Toluene to C12.

Diesel value based on the total peaks in the range from C12 to C24.

Oil value based on the total peaks in the range from C24 to C32.

Data Qualifiers

- U Compound not detected at the given detection limit.
- X Value detected above linear range of instrument. Dilution required.
- J Indicates an estimated value below the calculated detection limit.
- S No value reported due to saturation of the detector. Dilution required.
- D Indicates the surrogate was not detected because of dilution of the extract.
- C Indicates a probable value which cannot be confirmed due to matrix interference.
- NR Indicates no recovery due to matrix interference and/or dilution.



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Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

TOTAL PETROLEUM HYDROCARBONS
WA HCID Method by GC/FID

Lab Sample ID: J468SB
LIMS ID: 95-411
Matrix: Soil

QC Report No: J468-Landau Associates, Incorporated
Project: 275002
Paxton Sales

Data Release Authorized:
Reported: 01/23/95

Date extracted: 01/16/95

LABORATORY CONTROL SAMPLE RECOVERY REPORT

| CONSTITUENT | SPIKE VALUE | SPIKE ADDED | % RECOVERY |
|---------------------------|----------------|----------------|---------------|
| LABORATORY CONTROL SAMPLE | | | |
| Diesel Range | 539 | 500 | 108% |

HCID Surrogate Recovery

| | | |
|-------------|------------------|------|
| Spike Blank | Methylarachidate | 104% |
|-------------|------------------|------|

Values reported in parts per million (mg/kg)

HCID SPIKE CONTROL LIMITS

Percent Recovery 50-150%
Duplicate RPD <50%

Advisory QA Limits



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Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

ORGANIC COMPOUND DATA REPORTING QUALIFIERS

- U** Indicates the compound was analyzed for, but not detected at the given detection limit.
- J** Indicates an estimated value when the result is less than the calculated detection limit.
- D** Indicates the surrogate/spike(s) was not detected, due to dilution of extract.
- NR** Indicates the surrogate recovery cannot be reported due to matrix interference.
- E** Indicates a value above the linear range of the detector.
Sample dilution required.
- S** Indicates no value reported due to saturation of the detector.
Dilution required.
- Y** Indicates a raised detection limit due to matrix interferences.
- NA** Indicates compound was not analyzed.
- M** Indicates an estimated value of analyte found and confirmed by analyst but with low spectral match.
- B** Indicates compound was found in the associated method blank.

ORGANICS ANALYSIS DATA SHEET
Volatile by Purge & Trap GC/MS
Page 1 of 1



ANALYTICAL
RESOURCES
INCORPORATED

Sample No: MW-1-5.1'

Lab Sample ID: J468C

LIMS ID: 95-413

Matrix: Soil

Data Release Authorized: *OK*

Reported: 01/31/95

QC Report No: J468-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Sampled: 01/11/95

Date Received: 01/13/95

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Instrument: FINN5

Sample Amount: 4.63 g dry Wt

Date Analyzed: 01/25/95

Percent Moisture: 7.7%

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.2 U |
| 75-01-4 | Vinyl Chloride | 2.2 U |
| 75-00-3 | Chloroethane | 2.2 U |
| 75-09-2 | Methylene Chloride | 2.4 B |
| 75-35-4 | 1,1-Dichloroethene | 1.1 U |
| 75-34-3 | 1,1-Dichloroethane | 1.1 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.1 U |
| 67-66-3 | Chloroform | 1.1 U |
| 107-06-2 | 1,2-Dichloroethane | 1.1 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.1 U |
| 56-23-5 | Carbon Tetrachloride | 1.1 U |
| 75-27-4 | Bromodichloromethane | 1.1 U |
| 78-87-5 | 1,2-Dichloropropane | 1.1 U |
| 79-01-6 | Trichloroethene | 1.1 U |
| 124-48-1 | Dibromochloromethane | 1.1 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.1 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.1 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.4 U |
| 75-25-2 | Bromoform | 1.1 U |
| 127-18-4 | Tetrachloroethene | 1.1 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.1 U |
| 108-90-7 | Chlorobenzene | 1.1 U |
| 75-69-4 | Trichlorofluoromethane | 2.2 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.1 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.1 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.1 U |
| 74-96-4 | Bromoethane | 2.2 U |
| 74-95-3 | Dibromomethane | 1.1 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.1 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.1 U |
| 108-86-1 | Bromobenzene | 1.1 U |
| 95-49-8 | 2-Chlorotoluene | 1.1 U |
| 106-43-4 | 4-Chlorotoluene | 1.1 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 99.5% |
| d8-Toluene | 99.0% |
| Bromofluorobenzene | 99.2% |
| d4-1,2-Dichlorobenzene | 99.7% |

ORGANICS ANALYSIS DATA SHEET
Volatile by Purge & Trap GC/MS
Page 1 of 1



Sample No: MW-1-7.5

ANALYTICAL
RESOURCES
INCORPORATED

| | |
|-------------------------------------|--|
| Lab Sample ID: J468D | QC Report No: J468-Landau Associates, Incorporated |
| LIMS ID: 95-414 | Project: 275002 |
| Matrix: Soil | Paxton Sales |
| Data Release Authorized: <i>BSB</i> | Date Sampled: 01/11/95 |
| Reported: 01/31/95 | Date Received: 01/13/95 |

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Instrument: FINN5 Sample Amount: 4.67 g dry Wt
Date Analyzed: 01/25/95 Percent Moisture: 8.4%

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.1 U |
| 75-01-4 | Vinyl Chloride | 2.1 U |
| 75-00-3 | Chloroethane | 2.1 U |
| 75-09-2 | Methylene Chloride | 2.2 B |
| 75-35-4 | 1,1-Dichloroethene | 1.1 U |
| 75-34-3 | 1,1-Dichloroethane | 1.1 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.1 U |
| 67-66-3 | Chloroform | 1.1 U |
| 107-06-2 | 1,2-Dichloroethane | 1.1 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.1 U |
| 56-23-5 | Carbon Tetrachloride | 1.1 U |
| 75-27-4 | Bromodichloromethane | 1.1 U |
| 78-87-5 | 1,2-Dichloropropane | 1.1 U |
| 79-01-6 | Trichloroethene | 1.1 U |
| 124-48-1 | Dibromochloromethane | 1.1 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.1 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.1 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.4 U |
| 75-25-2 | Bromoform | 1.1 U |
| 127-18-4 | Tetrachloroethene | 1.1 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.1 U |
| 108-90-7 | Chlorobenzene | 1.1 U |
| 75-69-4 | Trichlorofluoromethane | 2.1 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.1 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.1 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.1 U |
| 74-96-4 | Bromoethane | 2.1 U |
| 74-95-3 | Dibromomethane | 1.1 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.1 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.1 U |
| 108-86-1 | Bromobenzene | 1.1 U |
| 95-49-8 | 2-Chlorotoluene | 1.1 U |
| 106-43-4 | 4-Chlorotoluene | 1.1 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 96.6% |
| d8-Toluene | 97.8% |
| Bromofluorobenzene | 98.2% |
| d4-1,2-Dichlorobenzene | 102% |

ORGANICS ANALYSIS DATA SHEET
Volatile by Purge & Trap GC/MS

Page 1 of 1



Sample No: MW-1-22.3

**ANALYTICAL
RESOURCES
INCORPORATED**

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J468I

LIMS ID: 95-419

Matrix: Soil

Data Release Authorized: *ON*

Reported: 01/31/95

QC Report No: J468-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Sampled: 01/11/95

Date Received: 01/13/95

Instrument: FINN5

Date Analyzed: 01/25/95

Sample Amount: 4.48 g dry Wt

Percent Moisture: 12.6%

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.2 U |
| 75-01-4 | Vinyl Chloride | 2.2 U |
| 75-00-3 | Chloroethane | 2.2 U |
| 75-09-2 | Methylene Chloride | 2.2 U |
| 75-35-4 | 1,1-Dichloroethene | 1.1 U |
| 75-34-3 | 1,1-Dichloroethane | 1.1 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.1 U |
| 67-66-3 | Chloroform | 1.1 U |
| 107-06-2 | 1,2-Dichloroethane | 1.1 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.1 U |
| 56-23-5 | Carbon Tetrachloride | 1.1 U |
| 75-27-4 | Bromodichloromethane | 1.1 U |
| 78-87-5 | 1,2-Dichloropropane | 1.1 U |
| 79-01-6 | Trichloroethene | 1.1 U |
| 124-48-1 | Dibromochloromethane | 1.1 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.1 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.1 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.6 U |
| 75-25-2 | Bromoform | 1.1 U |
| 127-18-4 | Tetrachloroethene | 1.1 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.1 U |
| 108-90-7 | Chlorobenzene | 1.1 U |
| 75-69-4 | Trichlorofluoromethane | 2.2 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.1 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.1 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.1 U |
| 74-96-4 | Bromoethane | 2.2 U |
| 74-95-3 | Dibromomethane | 1.1 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.1 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.1 U |
| 108-86-1 | Bromobenzene | 1.1 U |
| 95-49-8 | 2-Chlorotoluene | 1.1 U |
| 106-43-4 | 4-Chlorotoluene | 1.1 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 92.4% |
| d8-Toluene | 92.3% |
| Bromofluorobenzene | 88.1% |
| d4-1,2-Dichlorobenzene | 83.5% |

ORGANICS ANALYSIS DATA SHEET
Volatile by Purge & Trap GC/MS

Page 1 of 1



**ANALYTICAL
RESOURCES
INCORPORATED**

Sample No: MW-1-30.0

Lab Sample ID: J468M

LIMS ID: 95-423

Matrix: Soil

Data Release Authorized: *BNS*

Reported: 01/31/95

QC Report No: J468-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Sampled: 01/12/95

Date Received: 01/13/95

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Instrument: FINN5

Sample Amount: 4.80 g dry Wt

Date Analyzed: 01/25/95

Percent Moisture: 6.9%

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.1 U |
| 75-01-4 | Vinyl Chloride | 2.1 U |
| 75-00-3 | Chloroethane | 2.1 U |
| 75-09-2 | Methylene Chloride | 2.1 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.2 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.0 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.1 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-96-4 | Bromoethane | 2.1 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 96.6% |
| d8-Toluene | 97.2% |
| Bromofluorobenzene | 91.7% |
| d4-1,2-Dichlorobenzene | 87.4% |

ORGANICS ANALYSIS DATA SHEET
Volatile by Purge & Trap GC/MS



Page 1 of 1

**ANALYTICAL
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Analytical
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Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

| | |
|-------------------------------------|--|
| Lab Sample ID: 012595MB | QC Report No: J468-Landau Associates, Incorporated |
| LIMS ID: 95-413 | Project: 275002 |
| Matrix: Soil | Paxton Sales |
| Data Release Authorized: <i>BNB</i> | Date Sampled: NA |
| Reported: 01/31/95 | Date Received: NA |
| Instrument: FINN5 | Sample Amount: 5.00 g dry Wt Equiv |
| Date Analyzed: 01/25/95 | Percent Moisture: NA |

| CAS Number | Analyte | ug/kg |
|------------|----------------------------|--------------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.3 |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.0 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.5 |
| 541-73-1 | 1,3-Dichlorobenzene | 1.6 |
| 106-46-7 | 1,4-Dichlorobenzene | 1.6 M |
| 74-83-9 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.2 |
| 95-49-8 | 2-Chlorotoluene | 1.1 |
| 106-43-4 | 4-Chlorotoluene | 1.4 |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 101% |
| d8-Toluene | 94.4% |
| Bromofluorobenzene | 91.4% |
| d4-1,2-Dichlorobenzene | 90.2% |

ORGANICS ANALYSIS DATA SHEET
Volatile by Purge & Trap GC/MS
Page 1 of 1



Sample No: Method Blank

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Analytical
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Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: 012595MB

QC Report No: J468-Landau Associates, Incorporated

LIMS ID: 95-419

Project: 275002

Matrix: Soil

Paxton Sales

Data Release Authorized: *BNB*

Date Sampled: NA

Reported: 01/31/95

Date Received: NA

Instrument: FINN5

Sample Amount: 5.00 g dry Wt Equiv

Date Analyzed: 01/25/95

Percent Moisture: NA

| CAS Number | Analyte | ug/kg |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.0 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.0 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-83-9 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 87.6% |
| d8-Toluene | 93.9% |
| Bromofluorobenzene | 91.2% |
| d4-1,2-Dichlorobenzene | 87.0% |

ORGANICS ANALYSIS DATA SHEET
Volatiles by GC/MS
Page 1 of 1



ANALYTICAL
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Analytical
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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J468SB
LIMS ID: 95-419
Matrix: Soil
Data Release Authorized: *BMS*
Reported: 01/31/95
Date Analyzed: 01/25/95

QC Report No: J468-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/13/95

| LABORATORY CONTROL SAMPLE CONSTITUENT | SPIKE VALUE | SPIKE AMT | % RECOVERY |
|--|----------------|--------------|---------------|
| Chloromethane | 52.3 | 50.0 | 105% |
| Vinyl Chloride | 48.2 | 50.0 | 96.4% |
| Chloroethane | 46.1 | 50.0 | 92.2% |
| Methylene Chloride | 55.0 | 50.0 | 110% |
| 1,1-Dichloroethene | 55.3 | 50.0 | 111% |
| 1,1-Dichloroethane | 51.3 | 50.0 | 103% |
| trans-1,2-Dichloroethene | 54.7 | 50.0 | 109% |
| Chloroform | 51.0 | 50.0 | 102% |
| 1,2-Dichloroethane | 51.3 | 50.0 | 103% |
| 1,1,1-Trichloroethane | 51.3 | 50.0 | 103% |
| Carbon Tetrachloride | 49.2 | 50.0 | 98.4% |
| Bromodichloromethane | 48.8 | 50.0 | 97.6% |
| 1,2-Dichloropropane | 49.5 | 50.0 | 99.0% |
| Trichloroethene | 48.0 | 50.0 | 96.0% |
| Dibromochloromethane | 47.9 | 50.0 | 95.8% |
| 1,1,2-Trichloroethane | 45.4 | 50.0 | 90.8% |
| trans-1,3-Dichloropropene | 46.1 | 50.0 | 92.2% |
| 2-Chloroethylvinylether | 53.0 | 50.0 | 106% |
| Bromoform | 46.2 | 50.0 | 92.4% |
| Tetrachloroethene | 47.3 | 50.0 | 94.6% |
| 1,1,2,2-Tetrachloroethane | 51.8 | 50.0 | 104% |
| Chlorobenzene | 47.9 | 50.0 | 95.8% |
| Trichlorofluoromethane | 43.7 | 50.0 | 87.4% |
| 1,2-Dichlorobenzene | 47.3 | 50.0 | 94.6% |
| 1,3-Dichlorobenzene | 46.2 | 50.0 | 92.4% |
| 1,4-Dichlorobenzene | 47.7 | 50.0 | 95.4% |
| Bromoethane | 57.5 | 50.0 | 115% |
| Dibromomethane | 48.3 | 50.0 | 96.6% |
| 1,1,1,2-Tetrachloroethane | 47.8 | 50.0 | 95.6% |
| 1,2,3-Trichloropropane | 52.6 | 50.0 | 105% |
| Bromobenzene | 46.0 | 50.0 | 92.0% |
| 2-Chlorotoluene | 43.9 | 50.0 | 87.8% |
| 4-Chlorotoluene | 51.3 | 50.0 | 103% |

| <u>Spike Blank Surrogate Recovery</u> | |
|---------------------------------------|-------|
| d4-1,2-Dichloroethane | 98.0% |
| d8-Toluene | 89.4% |
| Bromofluorobenzene | 97.0% |
| d4-1,2-Dichlorobenzene | 94.0% |

Reported in ug/kg-dry-Wt



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Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

ORGANIC COMPOUND DATA REPORTING QUALIFIERS

- U** Indicates the compound was analyzed for, but not detected at the given detection limit.
- J** Indicates an estimated value when the result is less than the calculated detection limit.
- D** Indicates the surrogate/spike(s) was not detected, due to dilution of extract.
- NR** Indicates the surrogate recovery cannot be reported due to matrix interference.
- E** Indicates a value above the linear range of the detector.
Sample dilution required.
- S** Indicates no value reported due to saturation of the detector.
Dilution required.
- Y** Indicates a raised detection limit due to matrix interferences.
- NA** Indicates compound was not analyzed.
- M** Indicates an estimated value of analyte found and confirmed by analyst but with low spectral match.
- B** Indicates compound was found in the associated method blank.



ANALYTICAL
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ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 1 of 2

Lab Sample ID: J468G

LIMS ID: 95-417

Matrix: Soil

Data Release Authorized: MTH

Reported: 01/19/95

Sample No: MW-1-17.6

Analytical
Chemists &
Consultants

QC Report No: J468-Landau Associates, Incorporated

Project: 275002

Paxton Sales

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Date Sampled: 01/11/95

Date Received: 01/13/95

Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN4

GPC Cleanup: NO

Sample Amount: 27.1 g-dry-wt

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: 9.9%

pH: 8.2

| CAS Number | Analyte | ug/kg |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 150 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 150 U |
| 95-57-8 | 2-Chlorophenol | 74 U |
| 541-73-1 | 1,3-Dichlorobenzene | 74 U |
| 106-46-7 | 1,4-Dichlorobenzene | 74 U |
| 100-51-6 | Benzyl Alcohol | 370 U |
| 95-50-1 | 1,2-Dichlorobenzene | 74 U |
| 95-48-7 | 2-Methylphenol | 150 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 74 U |
| 106-44-5 | 4-Methylphenol | 74 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 150 U |
| 67-72-1 | Hexachloroethane | 150 U |
| 98-95-3 | Nitrobenzene | 74 U |
| 78-59-1 | Isophorone | 74 U |
| 88-75-5 | 2-Nitrophenol | 370 U |
| 105-67-9 | 2,4-Dimethylphenol | 220 U |
| 65-85-0 | Benzoic Acid | 740 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 74 U |
| 120-83-2 | 2,4-Dichlorophenol | 220 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 74 U |
| 91-20-3 | Naphthalene | 74 U |
| 106-47-8 | 4-Chloroaniline | 220 U |
| 87-68-3 | Hexachlorobutadiene | 150 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 150 U |
| 91-57-6 | 2-Methylnaphthalene | 74 U |
| 77-47-4 | Hexachlorocyclopentadiene | 370 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 370 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 370 U |
| 91-58-7 | 2-Chloronaphthalene | 74 U |
| 88-74-4 | 2-Nitroaniline | 370 U |
| 131-11-3 | Dimethylphthalate | 74 U |
| 208-96-8 | Acenaphthylene | 74 U |
| 99-09-2 | 3-Nitroaniline | 440 U |
| 83-32-9 | Acenaphthene | 74 U |
| 51-28-5 | 2,4-Dinitrophenol | 740 U |
| 100-02-7 | 4-Nitrophenol | 370 U |
| 132-64-9 | Dibenzofuran | 74 U |
| 606-20-2 | 2,6-Dinitrotoluene | 370 U |



ANALYTICAL
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Final Report
Laboratory Analysis of Conventional Parameters

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547A

LIMS ID: 95-824

Matrix: Water

Data Release Authorized: *Mop*

Reported: 02/10/95 *2-18*

QC Report No: J547-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Sampled: 01/23/95

Date Received: 01/24/95

| Analyte | Analysis Date | Method | RL | Units | Result |
|---------------|---------------|-----------|-------|-------|--------|
| Total Cyanide | 02/09/95 | EPA 335.2 | 0.004 | mg/L | 0.020 |

RL Analytical reporting limit

U Undetected at reported detection limit

Report for J547 received 01/24/95



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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

**Final Report
Laboratory Analysis of Conventional Parameters**

Lab Sample ID: J547B

LIMS ID: 95-825

Matrix: Water

Data Release Authorized: *Mof*
Reported: 02/10/95 *2-10*

Sample No: MW-1A-123
QC Report No: J547-Landau Associates, Incorporated (206) 621-7523 (FAX)
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

| Analyte | Analysis Date | Method | RL | Units | Result |
|----------------|----------------------|---------------|-----------|--------------|---------------|
| Total Cyanide | 02/09/95 | EPA 335.2 | 0.004 | mg/L | 0.031 |

RL Analytical reporting limit

U Undetected at reported detection limit

Report for J547 received 01/24/95



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(206) 621-6490

(206) 621-7523 (FAX)

**Final Report
Laboratory Analysis of Conventional Parameters**

Lab Sample ID: J547C
LIMS ID: 95-826
Matrix: Water
Data Release Authorized: *Moy*
Reported: 02/10/95
2-10

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

| <u>Analyte</u> | <u>Analysis Date</u> | <u>Method</u> | <u>RL</u> | <u>Units</u> | <u>Result</u> |
|----------------|----------------------|---------------|-----------|--------------|---------------|
| Total Cyanide | 02/09/95 | EPA 335.2 | 0.004 | mg/L | < 0.005 U |

RL Analytical reporting limit

U Undetected at reported detection limit

Report for J547 received 01/24/95



**ANALYTICAL
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333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

(206) 621-7523 (FAX)

**Final Report
Laboratory Analysis of Conventional Parameters**

Lab Sample ID: J547D
LIMS ID: 95-827

Matrix: Water

Data Release Authorized: *Mof*
Reported: 02/10/95
2/10

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

| <u>Analyte</u> | <u>Analysis Date</u> | <u>Method</u> | <u>RL</u> | <u>Units</u> | <u>Result</u> |
|----------------|----------------------|---------------|-----------|--------------|---------------|
| Total Cyanide | 02/09/95 | EPA 335.2 | 0.004 | mg/L | < 0.005 U |

RL Analytical reporting limit

U Undetected at reported detection limit

Report for J547 received 01/24/95



**ANALYTICAL
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INCORPORATED**

Analytical
Chemists &
Consultants

QA Report - Replicate Analysis

Matrix: Water

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/24/95

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Data Release Authorized: *Mof*
2-10

**DUPLICATE ANALYSIS RESULTS
CONVENTIONALS**

| Constituent | Units | Sample Value | Duplicate Value | RPD |
|--|--------------|---------------------|------------------------|------------|
| ARI ID: 95-826, J547 C Client Sample ID: MW-3-123 | | | | |
| Total Cyanide | mg/L | < 0.005 U | < 0.005 U | NA |



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(206) 621-7523 (FAX)

QA Report - Matrix Spike Analysis

Matrix: Water

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/24/95

Data Release Authorized: *Mof*
2/70

**MATRIX SPIKE QA/QC REPORT
CONVENTIONALS**

| <u>Constituent</u> | <u>Units</u> | <u>Sample Value</u> | <u>Spike Value</u> | <u>Spike Added</u> | <u>Recovery</u> |
|--|--------------|---------------------|--------------------|--------------------|-----------------|
| ARI ID: 95-827, J547 D Client Sample ID: MW-2-123 | | | | | |
| Total Cyanide | mg/L | < 0.00 | 0.208 | 0.189 | 110% |

ORGANICS ANALYSIS DATA SHEET
Volatile by Purge & Trap GC/MS

Page 1 of 1



Sample No: MW-2-123

ANALYTICAL
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Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547D

LIMS ID: 95-827

Matrix: Water

Data Release Authorized: *DR*

Reported: 02/08/95

QC Report No: J547-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Sampled: 01/23/95

Date Received: 01/24/95

Instrument: FINN3

Sample Amount: 5.00 mL

Date Analyzed: 01/31/95

| CAS Number | Analyte | ug/L |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.0 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.7 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-96-4 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 106% |
| d8-Toluene | 100% |
| Bromofluorobenzene | 103% |
| d4-1,2-Dichlorobenzene | 98.1% |



**ANALYTICAL
RESOURCES
INCORPORATED**

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

4206 621-7523 (FAX)

QA Report - Method Blank Analysis

Matrix: Water

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/24/95

Data Release Authorized: *Mop*
2-16

**METHOD BLANK RESULTS
CONVENTIONALS**

| <u>Analysis Date</u> | <u>Constituent</u> | <u>Units</u> | <u>Result</u> |
|----------------------|--------------------|--------------|---------------|
| 02/09/95 | Total Cyanide | mg/L | < 0.004 U |



**ANALYTICAL
RESOURCES
INCORPORATED**

Analytical
Chemists &
Consultants

QA Report - Standard Reference Material Analysis

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

(206) 621-7523 (FAX)

QC Report No: J547-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Data Release Authorized: *MSP*

Date Received: 01/24/95

2-10

**STANDARD REFERENCE MATERIAL ANALYSIS
CONVENTIONALS**

| Constituent | Units | Value | True Value | Recovery |
|---------------------|----------|-------|------------|----------|
| SPEX QC Lot#6-153AS | | | | |
| Total Cyanide | mg/L | 0.183 | 0.200 | 91.5% |
| Date analyzed: | 02/09/95 | | | |

EXPLANATION OF INORGANIC DATA REPORT CODES

The columns labeled 'PREP', 'C', and 'M' contain important information about your analyses. The codes are defined below.

PREPARATION CODES

These 3-letter codes describe methods used to prepare samples for analysis:

| | |
|-----|--|
| AEN | USEPA, Metals in air filters |
| AHM | ARI, Mercury in air filters |
| AHN | ARI, Metals in air filters |
| ANN | NIOSH 7300, Metals in air filters |
| CAN | AOAC (1984) 25.024, Metals in earthenware |
| DE6 | EPA 600/4-79-020 218.5, Cr(VI) in water |
| DMM | DMN followed by TMM, Dissolved mercury |
| DMN | Filtered through .45u filter, Dissolved metals |
| EW6 | EWN followed by DE6 |
| EWM | EWN followed by TMM |
| F N | USEPA SW-846 1310, EP Toxicity |
| FHP | ARI, Metals in tissue (HNO ₃ /HClO ₄) |
| FPP | PSEP, Metals in tissue (HNO ₃ /HClO ₄) |
| FRM | Journal, Mercury in tissue |
| FRN | Journal, Metals in tissue (HNO ₃ /H ₂ O ₂) |
| KRN | ARI, Concentration by coprecipitation |
| LEM | USEPA 1311, TCLP followed by TMM |
| LEN | USEPA 1311, TCLP Extraction |
| HJM | ARI, Mercury in miscellaneous materials |
| JHN | ARI, Metals in miscellaneous materials |
| OAM | ARI, Mercury in oil, grease or tar |
| OAN | ARI, Metals in oil, grease or tar |
| PHM | ARI, Mercury in wipes |
| PHN | ARI, Metals in wipes |
| RCC | USEPA CLP, Water digestion, HCl matrix |
| RCN | USEPA CLP, Water digestion, HNO ₃ matrix |
| REC | EPA 600/4-79-020 4.1.4, HCl matrix |
| I | EPA 600/4-79-020 200.7 and 9.4 |
| REN | EPA 600/4-79-020 4.1.4, HNO ₃ matrix |

| | |
|-----|--|
| PMA | EPA 600/4-79-020 206.2 |
| RWC | USEPA SW-846 3005 |
| SCC | USEPA CLP, Soil digestion, HCl matrix |
| SCM | USEPA CLP, Mercury in soil |
| SCN | USEPA CLP, Soil digestion, HNO ₃ matrix |
| SEM | EPA 600/4-79-020 245.5, Mercury in soil |
| SHF | ARI, Metals in soil, HF digestion |
| SMN | Agronomy, Metals in soil, Water extract |
| SMM | SMN followed by DMM, Dissolved mercury |
| SPM | USEPA 1312, SPLP extraction followed by TMM |
| SPN | USEPA 1312, SPLP Extraction |
| SSS | Standard Methods 302C, Ti in soil |
| SW6 | USEPA SW-846 3060, Cr(VI) in soil |
| SWC | USEPA SW-846 3050, HCl matrix |
| SWN | USEPA SW-846 3050, HNO ₃ matrix |
| SZF | PSEP/PSDDA, Microwave, Total acid digestion |
| TEC | EPA 600/4-79-020 4.1.3, HCl matrix |
| TEG | EPA 600/4-79-020 272.1, Silver in water |
| TEI | EPA 600/4-79-020 200.7 and 9.3 |
| TEN | EPA 600/4-79-020 4.1.3, HNO ₃ matrix |
| THG | ARI, Silver in photographic solutions |
| TMM | EPA 600/4-79-020 245.1, Mercury in water |
| TSC | Standard Methods 302C, Sb/Sn in water |
| TSN | Standard Methods 302D |
| TSS | Standard Methods 302E, Ti in water |
| TWC | USEPA SW-846 3010, HCl matrix |
| TWG | USEPA SW-846 7760, Silver in water |
| TWN | USEPA SW-846 3020, HNO ₃ matrix |
| WMN | EPA 600/4-79-020, Preserved, undigested water |

CONCENTRATION CODES

These codes are used to qualify reported concentrations:

J No analyte was detected. The reported value is the lower limit of detection.

METHOD CODES

These codes signify the instrumental technique used for analysis:

| | |
|-----|---|
| CVA | Cold Vapor Atomic Absorption Spectrophotometry |
| A | Flame Atomic Absorption Spectrophotometry |
| GFA | Graphite Furnace Atomic Absorption Spectrophotometry |
| CP | Inductively Coupled Plasma Atomic Emission Spectrometry |



ANALYTICAL
RESOURCES
INCORPORATED

ARI job number: J547

ID number: MW-1-123

Analytical
Chemists &
Consultants

ARI Sample number: A

Project: 275002

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Client: Landau Associates

Description:

Contact: Bill Evans

Sampled: 01/23/95

Matrix: Water

Received: 01/24/95

% Solids: 0.00

Released by: *[Signature]*

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|---------------|---|--------|------|-----|----------|
| 7440-36-0 | Antimony | 0.001 mg/L | U | 0.001 | RWC | GFA | 01/31/95 |
| 7440-38-2 | Arsenic | 0.024 mg/L | | 0.001 | RMA | GFA | 02/07/95 |
| 7440-41-7 | Beryllium | 0.003 mg/L | | 0.001 | TWC | ICP | 01/31/95 |
| 7440-43-9 | Cadmium | 0.002 mg/L | U | 0.002 | TWC | ICP | 01/31/95 |
| 7440-47-3 | Chromium | 0.106 mg/L | | 0.005 | TWC | ICP | 01/31/95 |
| 7440-50-8 | Copper | 0.210 mg/L | | 0.002 | TWC | ICP | 01/31/95 |
| 7439-92-1 | Lead | 0.038 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7439-97-6 | Mercury | 0.0005 mg/L | | 0.0001 | TMM | CVA | 01/27/95 |
| 7440-02-0 | Nickel | 0.09 mg/L | | 0.01 | TWC | ICP | 01/31/95 |
| 7782-49-2 | Selenium | 0.005 mg/L | U | 0.005 | RMA | GFA | 02/02/95 |
| 7440-22-4 | Silver | 0.003 mg/L | U | 0.003 | TWC | ICP | 01/31/95 |
| 7440-28-0 | Thallium | 0.002 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7440-66-6 | Zinc | 0.343 mg/L | | 0.004 | TWC | ICP | 01/31/95 |



ANALYTICAL
RESOURCES
INCORPORATED

ARI job number: J547
ARI Sample number: ASPK
Client: Landau Associates
Contact: Bill Evans
Matrix: Water
% Solids: 0.00

ID number: MW-1-123
Project: 275002
Description: Matrix Spike
Sampled: 01/23/95
Received: 01/24/95
Released by: _____

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|---------------|---|--------|------|-----|----------|
| 7440-36-0 | Antimony | 0.036 mg/L | | 0.005 | RWC | GFA | 01/31/95 |
| 7440-38-2 | Arsenic | 0.114 mg/L | | 0.005 | RMA | GFA | 02/07/95 |
| 7440-41-7 | Beryllium | 0.050 mg/L | | 0.001 | TWC | ICP | 01/31/95 |
| 7440-43-9 | Cadmium | 0.099 mg/L | | 0.002 | TWC | ICP | 01/31/95 |
| 7440-47-3 | Chromium | 0.353 mg/L | | 0.005 | TWC | ICP | 01/31/95 |
| 7440-50-8 | Copper | 0.307 mg/L | | 0.002 | TWC | ICP | 01/31/95 |
| 7439-92-1 | Lead | 0.140 mg/L | | 0.005 | TWN | GFA | 01/27/95 |
| 7439-97-6 | Mercury | 0.0015 mg/L | | 0.0001 | TMM | CVA | 01/27/95 |
| 7440-02-0 | Nickel | 0.58 mg/L | | 0.01 | TWC | ICP | 01/31/95 |
| 7782-49-2 | Selenium | 0.09 mg/L | | 0.01 | RMA | GFA | 02/02/95 |
| 7440-22-4 | Silver | 0.234 mg/L | | 0.003 | TWC | ICP | 01/31/95 |
| 7440-28-0 | Thallium | 0.100 mg/L | | 0.005 | TWN | GFA | 01/27/95 |
| 7440-66-6 | Zinc | 0.812 mg/L | | 0.004 | TWC | ICP | 01/31/95 |



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2

Lab Sample ID: J468G

LIMS ID: 95-417

Matrix: Soil

Data Release Authorized: MJA

Reported: 01/19/95

Sample No: MW-1-17.6

Analytical
Chemists &
Consultants

QC Report No: J468-Landau Associates, Incorporated

Project: 275002

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Paxton Sales

Date Sampled: 01/11/95

Date Received: 01/13/95

Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN4

GPC Cleanup: NO

Sample Amount: 27.1 g-dry-wt

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: 9.9%

pH: 8.2

| CAS Number | Analyte | ug/kg |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 370 U |
| 84-66-2 | Diethylphthalate | 74 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 74 U |
| 86-73-7 | Fluorene | 74 U |
| 100-01-6 | 4-Nitroaniline | 370 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 740 U |
| 86-30-6 | N-Nitrosodiphenylamine | 74 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 74 U |
| 118-74-1 | Hexachlorobenzene | 74 U |
| 87-86-5 | Pentachlorophenol | 370 U |
| 85-01-8 | Phenanthrene | 74 U |
| 86-74-8 | Carbazole | 74 U |
| 120-12-7 | Anthracene | 74 U |
| 84-74-2 | Di-n-Butylphthalate | 74 U |
| 206-44-0 | Fluoranthene | 74 U |
| 129-00-0 | Pyrene | 74 U |
| 85-68-7 | Butylbenzylphthalate | 74 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 370 U |
| 56-55-3 | Benzo(a)anthracene | 74 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 74 U |
| 218-01-9 | Chrysene | 74 U |
| 117-84-0 | Di-n-Octyl phthalate | 74 U |
| 205-99-2 | Benzo(b)fluoranthene | 74 U |
| 207-08-9 | Benzo(k)fluoranthene | 74 U |
| 50-32-8 | Benzo(a)pyrene | 74 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 74 U |
| 53-70-3 | Dibenz(a,h)anthracene | 74 U |
| 191-24-2 | Benzo(g,h,i)perylene | 74 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 85.4% | d5-Phenol | 77.7% |
| 2-Fluorobiphenyl | 87.2% | 2-Fluorophenol | 72.8% |
| d14-p-Terphenyl | 106% | 2,4,6-Tribromophenol | 75.8% |
| d4-1,2-Dichlorobenzene | 89.7% | d4-2-Chlorophenol | 78.6% |



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 1 of 2

Lab Sample ID: J468H

LIMS ID: 95-418

Matrix: Soil

Data Release Authorized: *MTH*

Reported: 01/19/95

Sample No: MW-1-20.2

Analytical
Chemists &
Consultants

QC Report No: J468-Landau Associates, Incorporated

Project: 275002

Paxton Sales

333 Ninth Ave. North
Seattle, WA 98109-5187

Date Sampled: 01/11/95

(206) 621-6490

Date Received: 01/13/95

(206) 621-7523 (FAX)

Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN4

GPC Cleanup: NO

Sample Amount: 26.2 g-dry-wt

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: 12.8%

pH: 8.0

| CAS Number | Analyte | ug/kg |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 150 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 150 U |
| 95-57-8 | 2-Chlorophenol | 76 U |
| 541-73-1 | 1,3-Dichlorobenzene | 76 U |
| 106-46-7 | 1,4-Dichlorobenzene | 76 U |
| 100-51-6 | Benzyl Alcohol | 380 U |
| 95-50-1 | 1,2-Dichlorobenzene | 76 U |
| 95-48-7 | 2-Methylphenol | 150 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 76 U |
| 106-44-5 | 4-Methylphenol | 76 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 150 U |
| 67-72-1 | Hexachloroethane | 150 U |
| 98-95-3 | Nitrobenzene | 76 U |
| 78-59-1 | Isophorone | 76 U |
| 88-75-5 | 2-Nitrophenol | 380 U |
| 105-67-9 | 2,4-Dimethylphenol | 230 U |
| 65-85-0 | Benzoic Acid | 760 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 76 U |
| 120-83-2 | 2,4-Dichlorophenol | 230 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 76 U |
| 91-20-3 | Naphthalene | 76 U |
| 106-47-8 | 4-Chloroaniline | 230 U |
| 87-68-3 | Hexachlorobutadiene | 150 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 150 U |
| 91-57-6 | 2-Methylnaphthalene | 76 U |
| 77-47-4 | Hexachlorocyclopentadiene | 380 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 380 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 380 U |
| 91-58-7 | 2-Chloronaphthalene | 76 U |
| 88-74-4 | 2-Nitroaniline | 380 U |
| 131-11-3 | Dimethylphthalate | 76 U |
| 208-96-8 | Acenaphthylene | 76 U |
| 99-09-2 | 3-Nitroaniline | 460 U |
| 83-32-9 | Acenaphthene | 76 U |
| 51-28-5 | 2,4-Dinitrophenol | 760 U |
| 100-02-7 | 4-Nitrophenol | 380 U |
| 132-64-9 | Dibenzofuran | 76 U |
| 606-20-2 | 2,6-Dinitrotoluene | 380 U |



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2

Lab Sample ID: J468H

LIMS ID: 95-418

Matrix: Soil

Data Release Authorized: MJA

Reported: 01/19/95

Sample No: MW-1-20.2

Analytical
Chemists &
Consultants

QC Report No: J468-Landau Associates, Incorporated

Project: 275002

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Paxton Sales

Date Sampled: 01/11/95

Date Received: 01/13/95

Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN4

GPC Cleanup: NO

Sample Amount: 26.2 g-dry-wt

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: 12.8%

pH: 8.0

| CAS Number | Analyte | ug/kg |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 380 U |
| 84-66-2 | Diethylphthalate | 76 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 76 U |
| 86-73-7 | Fluorene | 76 U |
| 100-01-6 | 4-Nitroaniline | 380 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 760 U |
| 86-30-6 | N-Nitrosodiphenylamine | 76 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 76 U |
| 118-74-1 | Hexachlorobenzene | 76 U |
| 87-86-5 | Pentachlorophenol | 380 U |
| 85-01-8 | Phenanthrene | 76 U |
| 86-74-8 | Carbazole | 76 U |
| 120-12-7 | Anthracene | 76 U |
| 84-74-2 | Di-n-Butylphthalate | 76 U |
| 206-44-0 | Fluoranthene | 76 U |
| 129-00-0 | Pyrene | 76 U |
| 85-68-7 | Butylbenzylphthalate | 76 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 380 U |
| 56-55-3 | Benzo(a)anthracene | 76 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 76 U |
| 218-01-9 | Chrysene | 76 U |
| 117-84-0 | Di-n-Octyl phthalate | 76 U |
| 205-99-2 | Benzo(b)fluoranthene | 76 U |
| 207-08-9 | Benzo(k)fluoranthene | 76 U |
| 50-32-8 | Benzo(a)pyrene | 76 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 76 U |
| 53-70-3 | Dibenz(a,h)anthracene | 76 U |
| 191-24-2 | Benzo(g,h,i)perylene | 76 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 84.2% | d5-Phenol | 76.3% |
| 2-Fluorobiphenyl | 78.2% | 2-Fluorophenol | 71.1% |
| d14-p-Terphenyl | 86.8% | 2,4,6-Tribromophenol | 73.0% |
| d4-1,2-Dichlorobenzene | 82.7% | d4-2-Chlorophenol | 79.0% |



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 1 of 2

Lab Sample ID: J468MB

LIMS ID: 95-417

Matrix: Soil

Data Release Authorized: MA

Reported: 01/19/95

Sample No: Method Blank

Analytical
Chemists &
Consultants

QC Report No: J468-Landau Associates, Incorporated

Project: 275002

333 Ninth Ave. North
Seattle, WA 98109-5187

Paxton Sales

(206) 621-6490

(206) 621-7523 (FAX)

Date Sampled: NA

Date Received: NA

Date extracted: 01/17/95

Date analyzed: 01/18/95

Instrument: FINN2

GPC Cleanup: NO

Sample Amount: 30.0 g-dry-wt Equiv

Final Extract Volume: 2.0 mL

Dilution Factor: 1:1

Percent Moisture: NA

pH: NA

| CAS Number | Analyte | ug/kg |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 130 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 130 U |
| 95-57-8 | 2-Chlorophenol | 67 U |
| 541-73-1 | 1,3-Dichlorobenzene | 67 U |
| 106-46-7 | 1,4-Dichlorobenzene | 67 U |
| 100-51-6 | Benzyl Alcohol | 330 U |
| 95-50-1 | 1,2-Dichlorobenzene | 67 U |
| 95-48-7 | 2-Methylphenol | 130 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 67 U |
| 106-44-5 | 4-Methylphenol | 67 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 130 U |
| 67-72-1 | Hexachloroethane | 130 U |
| 98-95-3 | Nitrobenzene | 67 U |
| 78-59-1 | Isophorone | 67 U |
| 88-75-5 | 2-Nitrophenol | 330 U |
| 105-67-9 | 2,4-Dimethylphenol | 200 U |
| 65-85-0 | Benzoic Acid | 670 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 67 U |
| 120-83-2 | 2,4-Dichlorophenol | 200 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 67 U |
| 91-20-3 | Naphthalene | 67 U |
| 106-47-8 | 4-Chloroaniline | 200 U |
| 87-68-3 | Hexachlorobutadiene | 130 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 130 U |
| 91-57-6 | 2-Methylnaphthalene | 67 U |
| 77-47-4 | Hexachlorocyclopentadiene | 330 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 330 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 330 U |
| 91-58-7 | 2-Chloronaphthalene | 67 U |
| 88-74-4 | 2-Nitroaniline | 330 U |
| 131-11-3 | Dimethylphthalate | 67 U |
| 208-96-8 | Acenaphthylene | 67 U |
| 99-09-2 | 3-Nitroaniline | 400 U |
| 83-32-9 | Acenaphthene | 67 U |
| 51-28-5 | 2,4-Dinitrophenol | 670 U |
| 100-02-7 | 4-Nitrophenol | 330 U |
| 132-64-9 | Dibenzofuran | 67 U |
| 606-20-2 | 2,6-Dinitrotoluene | 330 U |



ANALYTICAL
RESOURCES
INCORPORATED

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

621-7523 (FAX)

QA Report - Method Blank Analysis

Matrix: Soil

QC Report No: J468-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/13/95

Data Release Authorized: *MWP*
1-30

METHOD BLANK RESULTS
CONVENTIONALS

| <u>Analysis Date</u> | <u>Constituent</u> | <u>Units</u> | <u>Result</u> |
|--------------------------|--------------------|--------------|---------------|
| Method Blank 01/16/95 | Total Solids | mg residue | < 1 U |
| Method Blank 01/18/95 | Total Cyanide | mg/L | < 0.004 U |



ANALYTICAL
RESOURCES
INCORPORATED

Analytical
Chemists &
Consultants

QA Report - Standard Reference Material Analysis

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490

406-621-7523 (FAX)

QC Report No: J468-Landau Associates, Incorporated
Project: 275002

Paxton Sales

Data Release Authorized: *Mof*
1-30

Date Received: 01/13/95

**STANDARD REFERENCE MATERIAL ANALYSIS
CONVENTIONALS**

| Constituent | Units | Value | True Value | Recovery |
|---------------------|----------|-------|------------|----------|
| SPEX QC Lot#6-153AS | | | | |
| Total Cyanide | mg/L | 0.193 | 0.200 | 96.5% |
| Date analyzed: | 01/18/95 | | | |

EXPLANATION OF INORGANIC DATA REPORT CODES

The columns labeled 'PREP', 'C', and 'M' contain important information about your analyses. The codes are defined below.

PREPARATION CODES

These 3-letter codes describe methods used to prepare samples for analysis:

| | | | |
|-------------------|--|-----|--|
| AEN | USEPA, Metals in air filters | RMA | EPA 600/4-79-020 206.2 |
| AHM | ARI, Mercury in air filters | RWC | USEPA SW-846 3005 |
| AHN | ARI, Metals in air filters | SCC | USEPA CLP, Soil digestion, HCl matrix |
| ANN | NIOSH 7300, Metals in air filters | SCM | USEPA CLP, Mercury in soil |
| CAN | AOAC (1984) 25.024, Metals in earthenware | SCN | USEPA CLP, Soil digestion, HNO ₃ matrix |
| DE6 | EPA 600/4-79-020 218.5, Cr(VI) in water | SEM | EPA 600/4-79-020 245.5, Mercury in soil |
| DMM | DMN followed by TMM, Dissolved mercury | SHF | ARI, Metals in soil, HF digestion |
| DMN | Filtered through .45u filter, Dissolved metals | SMN | Agronomy, Metals in soil, Water extract |
| EW6 | EWN followed by DE6 | SMM | SMN followed by DMM, Dissolved mercury |
| EWM | EWN followed by TMM | SPM | USEPA 1312, SPLP extraction followed by TMM |
| F ^{EW} N | USEPA SW-846 1310, EP Toxicity | SPN | USEPA 1312, SPLP Extraction |
| F ^{EW} T | ARI, Metals in tissue (HNO ₃ /HClO ₄) | SSS | Standard Methods 302C, Ti in soil |
| FPP | PSEP, Metals in tissue (HNO ₃ /HClO ₄) | SW6 | USEPA SW-846 3060, Cr(VI) in soil |
| FRM | Journal, Mercury in tissue | SWC | USEPA SW-846 3050, HCl matrix |
| FRN | Journal, Metals in tissue (HNO ₃ /H ₂ O ₂) | SWN | USEPA SW-846 3050, HNO ₃ matrix |
| KRN | ARI, Concentration by coprecipitation | SZF | PSEP/PSDDA, Microwave, Total acid digestion |
| LEM | USEPA 1311, TCLP followed by TMM | TEC | EPA 600/4-79-020 4.1.3, HCl matrix |
| LEN | USEPA 1311, TCLP Extraction | TEG | EPA 600/4-79-020 272.1, Silver in water |
| MHM | ARI, Mercury in miscellaneous materials | TEI | EPA 600/4-79-020 200.7 and 9.3 |
| HN | ARI, Metals in miscellaneous materials | TEN | EPA 600/4-79-020 4.1.3, HNO ₃ matrix |
| OAM | ARI, Mercury in oil, grease or tar | THG | ARI, Silver in photographic solutions |
| OAN | ARI, Metals in oil, grease or tar | TMM | EPA 600/4-79-020 245.1, Mercury in water |
| PHM | ARI, Mercury in wipes | TSC | Standard Methods 302C, Sb/Sn in water |
| PHN | ARI, Metals in wipes | TSN | Standard Methods 302D |
| RCC | USEPA CLP, Water digestion, HCl matrix | TSS | Standard Methods 302E, Ti in water |
| RCN | USEPA CLP, Water digestion, HNO ₃ matrix | TWC | USEPA SW-846 3010, HCl matrix |
| REC | EPA 600/4-79-020 4.1.4, HCl matrix | TWG | USEPA SW-846 7760, Silver in water |
| RCW | EPA 600/4-79-020 200.7 and 9.4 | TWN | USEPA SW-846 3020, HNO ₃ matrix |
| RKN | EPA 600/4-79-020 4.1.4, HNO ₃ matrix | WMN | EPA 600/4-79-020, Preserved, undigested water |

CONCENTRATION CODES

These codes are used to qualify reported concentrations:

U No analyte was detected. The reported value is the lower limit of detection.

METHOD CODES

These codes signify the instrumental technique used for analysis:

| | |
|-----|---|
| CVA | Cold Vapor Atomic Absorption Spectrophotometry |
| FLA | Flame Atomic Absorption Spectrophotometry |
| GA | Graphite Furnace Atomic Absorption Spectrophotometry |
| ICP | Inductively Coupled Plasma Atomic Emission Spectrometry |



Analytical Resources, Incorporated
Analytical Chemists and Consultants

cc to MCM
mp 2/14

RECEIVED

FEB 14 1995

LANDAU ASSOCIATES, INC.
TACOMA

February 13, 1995

Mr. Bill Evans
Landau Associates, Inc.
3600 Port of Tacoma Road
Suite 501
Tacoma, WA 98424

RE: Project No. 275002 (Yakima) Paxton Sales / ARI Job No. J547

Dear Mr. Evans:

Please find enclosed original results and Chain-of-Custody Record (COC NO. 2508) for the above-referenced project. Analytical Resources, Incorporated (ARI), accepted the following four water samples on January 24, 1995:

MW-1-123

MW-1A-123

MW-3-123

MW-2-123

Two VOC containers associated with sample **MW-1-123** were received broken. The remaining samples were received intact and the laboratory was able to complete the requested analyses without incident. The laboratory analyzed the samples for volatile organic compounds by EPA method 8260 with a modified 8010 compound list, semivolatile organics by EPA method 8270, total petroleum hydrocarbons by WDOE method WTPH-D, metals following the EPA method 6010/7000 series, and cyanide by EPA method 335.2.

As always, a copy of this report and all associated raw data will remain on file with ARI. If you have any questions or require additional information, please feel free to contact me at your convenience. If I am unavailable, you can leave a message on my voice mail and I will return your call as soon as possible.

Sincerely,

ANALYTICAL RESOURCES, INC.

Bryan D. Anderson

Bryan D. Anderson
Project Manager
(206)340-2866, ext. 116

enclosures
cc: file J547

BDA/bda



**ANALYTICAL
RESOURCES
INCORPORATED**

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

ORGANIC COMPOUND DATA REPORTING QUALIFIERS

- U** Indicates the compound was analyzed for, but not detected at the given detection limit.
- J** Indicates an estimated value when the result is less than the calculated detection limit.
- D** Indicates the surrogate/spike(s) was not detected, due to dilution of extract.
- NR** Indicates the surrogate recovery cannot be reported due to matrix interference.
- E** Indicates a value above the linear range of the detector.
Sample dilution required.
- S** Indicates no value reported due to saturation of the detector.
Dilution required.
- Y** Indicates a raised detection limit due to matrix interferences.
- NA** Indicates compound was not analyzed.
- M** Indicates an estimated value of analyte found and confirmed by analyst but with low spectral match.
- B** Indicates compound was found in the associated method blank.

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



ANALYTICAL
RESOURCES
INCORPORATED

Sample No: MW-1-123

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

| | |
|------------------------------------|--|
| Lab Sample ID: J547A | QC Report No: J547-Landau Associates, Incorporated |
| LIMS ID: 95-824 | Project: 275002 |
| Matrix: Water | Paxton Sales |
| Data Release Authorized: <i>MM</i> | Date Sampled: 01/23/95 |
| Reported: 02/08/95 | Date Received: 01/24/95 |
| Instrument: FINN3 | Sample Amount: 5.00 mL |
| Date Analyzed: 01/31/95 | |

| CAS Number | Analyte | ug/L |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.0 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 2.3 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-96-4 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 107% |
| d8-Toluene | 97.9% |
| Bromofluorobenzene | 102% |
| d4-1,2-Dichlorobenzene | 104% |

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



ANALYTICAL
RESOURCES
INCORPORATED

Sample No: MW-1A-123

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547B QC Report No: J547-Landau Associates, Incorporated
LIMS ID: 95-825 Project: 275002
Matrix: Water Paxton Sales
Data Release Authorized: *NY* Date Sampled: 01/23/95
Reported: 02/08/95 Date Received: 01/24/95

Instrument: FINN3 Sample Amount: 5.00 mL
Date Analyzed: 01/31/95

| CAS Number | Analyte | ug/L |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.0 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 2.2 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-96-4 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 106% |
| d8-Toluene | 98.5% |
| Bromofluorobenzene | 96.2% |
| d4-1,2-Dichlorobenzene | 104% |

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



ANALYTICAL
RESOURCES
INCORPORATED

Sample No: MW-3-123

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547C
LIMS ID: 95-826
Matrix: Water
Data Release Authorized: *NOP*
Reported: 02/08/95

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

Instrument: FINN3
Date Analyzed: 01/31/95

Sample Amount: 5.00 mL

| CAS Number | Analyte | ug/L |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.0 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 3.1 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-96-4 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 110% |
| d8-Toluene | 98.6% |
| Bromofluorobenzene | 102% |
| d4-1,2-Dichlorobenzene | 102% |

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



ANALYTICAL
RESOURCES
INCORPORATED

Sample No: Trip Blank

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

| | |
|------------------------------------|--|
| Lab Sample ID: J547E | QC Report No: J547-Landau Associates, Incorporated |
| LIMS ID: 95-828 | Project: 275002 |
| Matrix: Water | Paxton Sales |
| Data Release Authorized: <i>DK</i> | Date Sampled: 01/20/95 |
| Reported: 02/08/95 | Date Received: 01/24/95 |
| Instrument: FINN3 | Sample Amount: 5.00 mL |
| Date Analyzed: 01/31/95 | |

| CAS Number | Analyte | ug/L |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.0 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.0 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-96-4 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 105% |
| d8-Toluene | 96.5% |
| Bromofluorobenzene | 108% |
| d4-1,2-Dichlorobenzene | 98.2% |

ORGANICS ANALYSIS DATA SHEET
Volatiles by Purge & Trap GC/MS
Page 1 of 1



ANALYTICAL
RESOURCES
INCORPORATED

Sample No: Method Blank

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

| | |
|-------------------------------------|--|
| Lab Sample ID: 013195MB | QC Report No: J547-Landau Associates, Incorporated |
| LIMS ID: 95-824 | Project: 275002 |
| Matrix: Water | Paxton Sales |
| Data Release Authorized: <i>OSR</i> | Date Sampled: NA |
| Reported: 02/08/95 | Date Received: NA |
| Instrument: FINN3 | Sample Amount: 5.00 mL |
| Date Analyzed: 01/31/95 | |

| CAS Number | Analyte | ug/L |
|------------|---------------------------|-------|
| 74-87-3 | Chloromethane | 2.0 U |
| 75-01-4 | Vinyl Chloride | 2.0 U |
| 75-00-3 | Chloroethane | 2.0 U |
| 75-09-2 | Methylene Chloride | 2.0 U |
| 75-35-4 | 1,1-Dichloroethene | 1.0 U |
| 75-34-3 | 1,1-Dichloroethane | 1.0 U |
| 156-60-5 | trans-1,2-Dichloroethene | 1.0 U |
| 67-66-3 | Chloroform | 1.0 U |
| 107-06-2 | 1,2-Dichloroethane | 1.0 U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.0 U |
| 56-23-5 | Carbon Tetrachloride | 1.0 U |
| 75-27-4 | Bromodichloromethane | 1.0 U |
| 78-87-5 | 1,2-Dichloropropane | 1.0 U |
| 79-01-6 | Trichloroethene | 1.0 U |
| 124-48-1 | Dibromochloromethane | 1.0 U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.0 U |
| 10061-02-6 | trans-1,3-Dichloropropene | 1.0 U |
| 110-75-8 | 2-Chloroethylvinylether | 5.0 U |
| 75-25-2 | Bromoform | 1.0 U |
| 127-18-4 | Tetrachloroethene | 1.0 U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.0 U |
| 108-90-7 | Chlorobenzene | 1.0 U |
| 75-69-4 | Trichlorofluoromethane | 2.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 74-83-9 | Bromoethane | 2.0 U |
| 74-95-3 | Dibromomethane | 1.0 U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.0 U |
| 96-18-4 | 1,2,3-Trichloropropane | 1.0 U |
| 108-86-1 | Bromobenzene | 1.0 U |
| 95-49-8 | 2-Chlorotoluene | 1.0 U |
| 106-43-4 | 4-Chlorotoluene | 1.0 U |

Volatile Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 117% |
| d8-Toluene | 101% |
| Bromofluorobenzene | 99.9% |
| d4-1,2-Dichlorobenzene | 103% |

ORGANICS ANALYSIS DATA SHEET
Volatile s by GC/MS
Page 1 of 1



ANALYTICAL
RESOURCES
INCORPORATED

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547SB
LIMS ID: 95-824
Matrix: Water
Data Release Authorized: *MH*
Reported: 02/08/95
Date Analyzed: 01/31/95

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/24/95

| LABORATORY CONTROL SAMPLE CONSTITUENT | SPIKE VALUE | SPIKE AMT | % RECOVERY |
|--|----------------|--------------|---------------|
| Chloromethane | 40.8 | 50.0 | 81.6% |
| Vinyl Chloride | 42.7 | 50.0 | 85.4% |
| Chloroethane | 37.1 | 50.0 | 74.2% |
| Methylene Chloride | 49.6 | 50.0 | 99.2% |
| 1,1-Dichloroethene | 46.8 | 50.0 | 93.6% |
| 1,1-Dichloroethane | 50.5 | 50.0 | 101% |
| trans-1,2-Dichloroethene | 50.5 | 50.0 | 101% |
| Chloroform | 46.0 | 50.0 | 92.0% |
| 1,2-Dichloroethane | 46.1 | 50.0 | 92.2% |
| 1,1,1-Trichloroethane | 51.6 | 50.0 | 103% |
| Carbon Tetrachloride | 47.3 | 50.0 | 94.6% |
| Bromodichloromethane | 48.7 | 50.0 | 97.4% |
| 1,2-Dichloropropane | 49.1 | 50.0 | 98.2% |
| Trichloroethene | 48.1 | 50.0 | 96.2% |
| Dibromochloromethane | 48.0 | 50.0 | 96.0% |
| 1,1,2-Trichloroethane | 46.5 | 50.0 | 93.0% |
| trans-1,3-Dichloropropene | 45.5 | 50.0 | 91.0% |
| 2-Chloroethylvinylether | 51.9 | 50.0 | 104% |
| Bromoform | 45.6 | 50.0 | 91.2% |
| Tetrachloroethene | 47.1 | 50.0 | 94.2% |
| 1,1,2,2-Tetrachloroethane | 48.0 | 50.0 | 96.0% |
| Chlorobenzene | 48.6 | 50.0 | 97.2% |
| Trichlorofluoromethane | 34.4 | 50.0 | 68.8% |
| 1,2-Dichlorobenzene | 48.0 | 50.0 | 96.0% |
| 1,3-Dichlorobenzene | 47.5 | 50.0 | 95.0% |
| 1,4-Dichlorobenzene | 47.5 | 50.0 | 95.0% |
| Bromoethane | 51.8 | 50.0 | 104% |
| Dibromomethane | 47.7 | 50.0 | 95.4% |
| 1,1,1,2-Tetrachloroethane | 47.2 | 50.0 | 94.4% |
| 1,2,3-Trichloropropane | 50.2 | 50.0 | 100% |
| Bromobenzene | 45.3 | 50.0 | 90.6% |
| 2-Chlorotoluene | 45.6 | 50.0 | 91.2% |
| 4-Chlorotoluene | 48.5 | 50.0 | 97.0% |

Spike Blank Surrogate Recovery

| | |
|------------------------|-------|
| d4-1,2-Dichloroethane | 106% |
| d8-Toluene | 98.1% |
| Bromofluorobenzene | 98.8% |
| d4-1,2-Dichlorobenzene | 102% |

Reported in ug/L



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TOTAL DIESEL RANGE HYDROCARBONS
WA TPHd Range C12 to C24 by GC/FID

Matrix: Water

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Received: 01/24/95
Reported: 02/01/95

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Cathy M. Neumeier

| Lab ID | Client Sample ID | Date Extracted | Date Analyzed | Dilution Factor | Diesel Hydrocarbons | Range ID | *Diesel Surrogate Recovery | |
|--------------|------------------|----------------|---------------|-----------------|---------------------|----------|----------------------------|------|
| J547-0128MB | Method Blank | 01/25/95 | 01/28/95 | 1:1 | 0.25 | U | -- | 126% |
| 95-824-J547A | MW-1-123 | 01/25/95 | 01/28/95 | 1:1 | 0.25 | U | -- | 123% |
| 95-825-J547B | MW-1A-123 | 01/25/95 | 01/28/95 | 1:1 | 0.25 | U | -- | 123% |
| 95-826-J547C | MW-3-123 | 01/25/95 | 01/28/95 | 1:1 | 0.25 | U | -- | 121% |
| 95-827-J547D | MW-2-123 | 01/25/95 | 01/28/95 | 1:1 | 0.25 | U | -- | 139% |

Surrogate is Methyl Arachidate.

* ID indicates, in the opinion of the analyst, the petroleum product with the best pattern match. 'NO' indicates that there was not a good match for any of the requested products. Values reported in ppm (mg/L) on an weight-as-received basis.
Diesel quantitation on total peaks in the range from C12 to C24.

Data Qualifiers

- U Compound not detected at the given detection limit.
- X Value detected above linear range of instrument. Dilution required.
- J Indicates an estimated value below the calculated detection limit.
- S No value reported due to saturation of the detector. Dilution required.
- D Indicates the surrogate was not detected because of dilution of the extract.
- C Indicates a probable value which cannot be confirmed due to matrix interference.
- NR Indicates no recovery due to matrix interference and/or dilution.



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TOTAL DIESEL RANGE HYDROCARBONS
WA TPHd Range C12 to C24 by GC/FID

Lab Sample ID: J547SB
LIMS ID: 95-824
Matrix: Water

QC Report No: J547-Landau Associates, Inc.
Project: 275002
Paxton Sales

Data Release Authorized:
Reported: 02/01/95

LABORATORY CONTROL SAMPLE RECOVERY REPORT

| CONSTITUENT | SPIKE FOUND | SPIKE ADDED | % RECOVERY |
|---------------------------|----------------|----------------|---------------|
| Diesel Range Hydrocarbons | 2.66 | 2.50 | 106% |

TPHd Surrogate Recovery

Methylarachidate 139%

Values reported in parts per million (mg/L)



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Matrix Spike Quality Control Report

Client: Landau Associates
Client's sample ID: MW-1-123
ARI sample ID: J547 ASPK
Units: mg/L

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

| Analyte | Meth | Sample | Matrix Spike | Spike Added | %R | Control Limit | Q |
|-----------|------|--------|--------------|-------------|-------|---------------|---|
| Antimony | GFA | 0 | 0.036 | 0.100 | 36.0 | 75-125% | N |
| Arsenic | GFA | 0.024 | 0.114 | 0.100 | 90.0 | 75-125% | |
| Beryllium | ICP | 0.003 | 0.050 | 0.050 | 94.0 | 75-125% | |
| Cadmium | ICP | 0 | 0.100 | 0.100 | 100.0 | 75-125% | |
| Chromium | ICP | 0.106 | 0.353 | 0.250 | 98.8 | 75-125% | |
| Copper | ICP | 0.210 | 0.307 | 0.100 | 97.0 | 75-125% | |
| Lead | GFA | 0.038 | 0.140 | 0.100 | 102.0 | 75-125% | |
| Mercury | CVA | 0.0005 | 0.0015 | 0.0010 | 100.0 | 75-125% | |
| Nickel | ICP | 0.09 | 0.58 | 0.50 | 98.0 | 75-125% | |
| Selenium | GFA | 0 | 0.089 | 0.100 | 89.0 | 75-125% | |
| Silver | ICP | 0 | 0.234 | 0.250 | 93.6 | 75-125% | |
| Thallium | GFA | 0.002 | 0.100 | 0.100 | 98.0 | 75-125% | |
| Zinc | ICP | 0.343 | 0.812 | 0.500 | 93.8 | 75-125% | |

%R = Percent Recovery

'Q' codes: 'N' = control limit not met
'H' = %R not applicable, sample concentration too high
'S' = Analyte not spiked



ANALYTICAL
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ARI job number: J547

ARI Sample number: B

Client: Landau Associates

Contact: Bill Evans

Matrix: Water

% Solids: 0.00

ID number: MW-1A-123

Project: 275002

Description:

Sampled: 01/23/95

Received: 01/24/95

Released by: B/E

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Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|---------------|---|--------|------|-----|----------|
| 7440-36-0 | Antimony | 0.001 mg/L | U | 0.001 | RWC | GFA | 01/31/95 |
| 7440-38-2 | Arsenic | 0.007 mg/L | | 0.001 | RMA | GFA | 02/02/95 |
| 7440-41-7 | Beryllium | 0.002 mg/L | | 0.001 | TWC | ICP | 01/31/95 |
| 7440-43-9 | Cadmium | 0.002 mg/L | U | 0.002 | TWC | ICP | 01/31/95 |
| 7440-47-3 | Chromium | 0.085 mg/L | | 0.005 | TWC | ICP | 01/31/95 |
| 7440-50-8 | Copper | 0.185 mg/L | | 0.002 | TWC | ICP | 01/31/95 |
| 7439-92-1 | Lead | 0.034 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7439-97-6 | Mercury | 0.0005 mg/L | | 0.0001 | TMM | CVA | 01/27/95 |
| 7440-02-0 | Nickel | 0.08 mg/L | | 0.01 | TWC | ICP | 01/31/95 |
| 7782-49-2 | Selenium | 0.005 mg/L | U | 0.005 | RMA | GFA | 02/02/95 |
| 7440-22-4 | Silver | 0.003 mg/L | U | 0.003 | TWC | ICP | 01/31/95 |
| 7440-28-0 | Thallium | 0.002 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7440-66-6 | Zinc | 0.289 mg/L | | 0.004 | TWC | ICP | 01/31/95 |



ANALYTICAL
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ARI job number: J547
ARI Sample number: BDUP

Client: Landau Associates
Contact: Bill Evans
Matrix: Water

% Solids: 0.00

ID number: MW-1A-123
Project: 275002

Description: Laboratory Duplicate
Sampled: 01/23/95
Received: 01/24/95

Released by: JK

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(206) 621-6490
(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|---------------|---|--------|------|-----|----------|
| 7440-36-0 | Antimony | 0.001 mg/L | U | 0.001 | RWC | GFA | 01/31/95 |
| 7440-38-2 | Arsenic | 0.007 mg/L | | 0.001 | RMA | GFA | 02/02/95 |
| 7440-41-7 | Beryllium | 0.002 mg/L | | 0.001 | TWC | ICP | 01/31/95 |
| 7440-43-9 | Cadmium | 0.002 mg/L | | 0.002 | TWC | ICP | 01/31/95 |
| 7440-47-3 | Chromium | 0.094 mg/L | | 0.005 | TWC | ICP | 01/31/95 |
| 7440-50-8 | Copper | 0.184 mg/L | | 0.002 | TWC | ICP | 01/31/95 |
| 7439-92-1 | Lead | 0.035 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7439-97-6 | Mercury | 0.0004 mg/L | | 0.0001 | TMM | CVA | 01/27/95 |
| 7440-02-0 | Nickel | 0.08 mg/L | | 0.01 | TWC | ICP | 01/31/95 |
| 7782-49-2 | Selenium | 0.006 mg/L | | 0.005 | RMA | GFA | 02/02/95 |
| 7440-22-4 | Silver | 0.003 mg/L | U | 0.003 | TWC | ICP | 01/31/95 |
| 7440-28-0 | Thallium | 0.001 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7440-66-6 | Zinc | 0.296 mg/L | | 0.004 | TWC | ICP | 01/31/95 |



ANALYTICAL
RESOURCES
INCORPORATED

ARI job number: J547

ARI Sample number: D

Client: Landau Associates

Contact: Bill Evans

Matrix: Water

% Solids: 0.00

ID number: MW-2-123

Project: 275002

Description:

Sampled: 01/23/95

Received: 01/24/95

Released by: PK

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(206) 621-7523 (FAX)

A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|---------------|---|--------|------|-----|----------|
| 7440-36-0 | Antimony | 0.001 mg/L | U | 0.001 | RWC | GFA | 01/31/95 |
| 7440-38-2 | Arsenic | 0.007 mg/L | | 0.001 | RMA | GFA | 01/30/95 |
| 7440-41-7 | Beryllium | 0.003 mg/L | | 0.001 | TWC | ICP | 01/31/95 |
| 7440-43-9 | Cadmium | 0.002 mg/L | U | 0.002 | TWC | ICP | 01/31/95 |
| 7440-47-3 | Chromium | 0.127 mg/L | | 0.005 | TWC | ICP | 01/31/95 |
| 7440-50-8 | Copper | 0.289 mg/L | | 0.002 | TWC | ICP | 01/31/95 |
| 7439-92-1 | Lead | 0.046 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7439-97-6 | Mercury | 0.0005 mg/L | | 0.0001 | TMM | CVA | 01/27/95 |
| 7440-02-0 | Nickel | 0.10 mg/L | | 0.01 | TWC | ICP | 01/31/95 |
| 7782-49-2 | Selenium | 0.007 mg/L | | 0.005 | RMA | GFA | 02/02/95 |
| 7440-22-4 | Silver | 0.003 mg/L | U | 0.003 | TWC | ICP | 01/31/95 |
| 7440-28-0 | Thallium | 0.002 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7440-66-6 | Zinc | 0.403 mg/L | | 0.004 | TWC | ICP | 01/31/95 |



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Matrix Duplicate Quality Control Report

Client: Landau Associates
Client's sample ID: MW-1A-123
ARI sample ID: J547 BDUP
Units: mg/L

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(206) 621-6490
(206) 621-7523 (FAX)

| Analyte | Meth | Original Sample | Matrix Duplicate | RPD | Control Limit | Q |
|-----------|------|-----------------|------------------|------|---------------|---|
| Antimony | GFA | U 0.001 | U 0.001 | 0.0 | ± 0.001 | L |
| Arsenic | GFA | 0.007 | 0.007 | 0.0 | ± 20 % | |
| Beryllium | ICP | 0.002 | 0.002 | 0.0 | ± 0.001 | L |
| Cadmium | ICP | U 0.002 | 0.002 | 0.0 | ± 0.002 | L |
| Chromium | ICP | 0.085 | 0.094 | 10.1 | ± 20 % | |
| Copper | ICP | 0.185 | 0.184 | 0.5 | ± 20 % | |
| Lead | GFA | 0.034 | 0.035 | 2.9 | ± 20 % | |
| Mercury | CVA | 0.0005 | 0.0004 | 22.2 | ± 0.0001 | L |
| Nickel | ICP | 0.08 | 0.08 | 0.0 | ± 20 % | |
| Selenium | GFA | U 0.005 | 0.006 | 18.2 | ± 0.005 | L |
| Silver | ICP | U 0.003 | U 0.003 | 0.0 | ± 0.003 | L |
| Thallium | GFA | 0.002 | 0.001 | 66.7 | ± 0.001 | L |
| Zinc | ICP | 0.289 | 0.296 | 2.4 | ± 20 % | |

RPD = Relative Percent Difference

'Q' codes: '*' = control limit not met
'L' = RPD not valid, alternate limit = ± detection limit



ANALYTICAL
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ARI job number: J547
ARI Sample number: C
Client: Landau Associates
Contact: Bill Evans
Matrix: Water
% Solids: 0.00

ID number: MW-3-123
Project: 275002
Description:
Sampled: 01/23/95
Received: 01/24/95
Released by: *[Signature]*

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A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|---------------|---|--------|------|-----|----------|
| 7440-36-0 | Antimony | 0.001 mg/L | U | 0.001 | RWC | GFA | 01/31/95 |
| 7440-38-2 | Arsenic | 0.007 mg/L | | 0.001 | RMA | GFA | 01/30/95 |
| 7440-41-7 | Beryllium | 0.001 mg/L | | 0.001 | TWC | ICP | 01/31/95 |
| 7440-43-9 | Cadmium | 0.002 mg/L | U | 0.002 | TWC | ICP | 01/31/95 |
| 7440-47-3 | Chromium | 0.042 mg/L | | 0.005 | TWC | ICP | 01/31/95 |
| 7440-50-8 | Copper | 0.074 mg/L | | 0.002 | TWC | ICP | 01/31/95 |
| 7439-92-1 | Lead | 0.015 mg/L | | 0.001 | TWN | GFA | 01/27/95 |
| 7439-97-6 | Mercury | 0.0001 mg/L | U | 0.0001 | TMM | CVA | 01/27/95 |
| 7440-02-0 | Nickel | 0.04 mg/L | | 0.01 | TWC | ICP | 01/31/95 |
| 7782-49-2 | Selenium | 0.007 mg/L | | 0.005 | RMA | GFA | 02/02/95 |
| 7440-22-4 | Silver | 0.003 mg/L | U | 0.003 | TWC | ICP | 01/31/95 |
| 7440-28-0 | Thallium | 0.001 mg/L | U | 0.001 | TWN | GFA | 01/27/95 |
| 7440-66-6 | Zinc | 0.122 mg/L | | 0.004 | TWC | ICP | 01/31/95 |



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ARI job number: J547

ARI Sample number: MB

Client: Landau Associates

Contact: Bill Evans

Matrix: Water

% Solids: 0.00

ID number:

Project: 275002

Description: Method Blank

Sampled: / /

Received: / /

Released by: *Bru*

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A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|-----------|---------------|---|--------|------|-----|----------|
| 7440-36-0 | Antimony | 0.001 mg/L | U | 0.001 | RWC | GFA | 01/31/95 |
| 7440-38-2 | Arsenic | 0.001 mg/L | U | 0.001 | RMA | GFA | 01/30/95 |
| 7440-41-7 | Beryllium | 0.001 mg/L | U | 0.001 | TWC | ICP | 01/31/95 |
| 7440-43-9 | Cadmium | 0.002 mg/L | U | 0.002 | TWC | ICP | 01/31/95 |
| 7440-47-3 | Chromium | 0.005 mg/L | U | 0.005 | TWC | ICP | 01/31/95 |
| 7440-50-8 | Copper | 0.002 mg/L | U | 0.002 | TWC | ICP | 01/31/95 |
| 7439-92-1 | Lead | 0.001 mg/L | | 0.001 | TWN | GFA | 01/30/95 |
| 7439-97-6 | Mercury | 0.0001 mg/L | U | 0.0001 | TMM | CVA | 01/27/95 |
| 7440-02-0 | Nickel | 0.01 mg/L | U | 0.01 | TWC | ICP | 01/31/95 |
| 7782-49-2 | Selenium | 0.001 mg/L | U | 0.001 | RMA | GFA | 01/31/95 |
| 7440-22-4 | Silver | 0.003 mg/L | U | 0.003 | TWC | ICP | 01/31/95 |
| 7440-28-0 | Thallium | 0.001 mg/L | U | 0.001 | TWN | GFA | 01/27/95 |
| 7440-66-6 | Zinc | 0.004 mg/L | U | 0.004 | TWC | ICP | 01/31/95 |



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ARI job number: J547

ARI Sample number: MB2

Client: Landau Associates

Contact: Bill Evans

Matrix: Water

% Solids: 0.00

ID number:

Project: 275002

Description: Method Blank

Sampled: / /

Received: / /

Released by: JK

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Seattle, WA 98109-5187
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A N A L Y T I C A L R E S U L T S

| CAS Number | Analyte | Concentration | C | LOD | Prep | M | Analyzed |
|------------|---------|---------------|---|-------|------|-----|----------|
| 7440-38-2 | Arsenic | 0.001 mg/L | U | 0.001 | RMA | GFA | 02/07/95 |



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ORGANIC COMPOUND DATA REPORTING QUALIFIERS

- U Indicates the compound was analyzed for, but not detected at the given detection limit.
- J Indicates an estimated value when the result is less than the calculated detection limit.
- D Indicates the surrogate/spike(s) was not detected, due to dilution of extract.
- NR Indicates the surrogate recovery cannot be reported due to matrix interference.
- E Indicates a value above the linear range of the detector.
Sample dilution required.
- S Indicates no value reported due to saturation of the detector.
Dilution required.
- Y Indicates a raised detection limit due to matrix interferences.
- NA Indicates compound was not analyzed.
- M Indicates an estimated value of analyte found and confirmed by analyst but with low spectral match.
- B Indicates compound was found in the associated method blank.

ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Page 1 of 2



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333 Ninth Ave. North
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(206) 621-7523 (FAX)

Lab Sample ID: J547A
LIMS ID: 95-824
Matrix: Water
Data Release Authorized: *OK*
Reported: 02/08/95

Sample No: MW-1-123

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 2.0 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 2.0 U |
| 95-57-8 | 2-Chlorophenol | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 100-51-6 | Benzyl Alcohol | 5.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 95-48-7 | 2-Methylphenol | 2.0 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 1.0 U |
| 106-44-5 | 4-Methylphenol | 1.0 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 2.0 U |
| 67-72-1 | Hexachloroethane | 2.0 U |
| 98-95-3 | Nitrobenzene | 1.0 U |
| 78-59-1 | Isophorone | 1.0 U |
| 88-75-5 | 2-Nitrophenol | 5.0 U |
| 105-67-9 | 2,4-Dimethylphenol | 3.0 U |
| 65-85-0 | Benzoic Acid | 10 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 1.0 U |
| 120-83-2 | 2,4-Dichlorophenol | 3.0 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 U |
| 91-20-3 | Naphthalene | 1.0 U |
| 106-47-8 | 4-Chloroaniline | 3.0 U |
| 87-68-3 | Hexachlorobutadiene | 2.0 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 2.0 U |
| 91-57-6 | 2-Methylnaphthalene | 1.0 U |
| 77-47-4 | Hexachlorocyclopentadiene | 5.0 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 5.0 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 5.0 U |
| 91-58-7 | 2-Chloronaphthalene | 1.0 U |
| 88-74-4 | 2-Nitroaniline | 5.0 U |
| 131-11-3 | Dimethylphthalate | 1.0 U |
| 208-96-8 | Acenaphthylene | 1.0 U |
| 99-09-2 | 3-Nitroaniline | 6.0 U |
| 83-32-9 | Acenaphthene | 1.0 U |
| 51-28-5 | 2,4-Dinitrophenol | 10 U |
| 100-02-7 | 4-Nitrophenol | 5.0 U |
| 132-64-9 | Dibenzofuran | 1.0 U |
| 606-20-2 | 2,6-Dinitrotoluene | 5.0 U |

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2



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(206) 621-7523 (FAX)

Lab Sample ID: J547A
LIMS ID: 95-824
Matrix: Water
Data Release Authorized: *OK*
Reported: 02/08/95

Sample No: MW-1-123

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 5.0 U |
| 84-66-2 | Diethylphthalate | 1.0 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 1.0 U |
| 86-73-7 | Fluorene | 1.0 U |
| 100-01-6 | 4-Nitroaniline | 5.0 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 10 U |
| 86-30-6 | N-Nitrosodiphenylamine | 1.0 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 1.0 U |
| 118-74-1 | Hexachlorobenzene | 1.0 U |
| 87-86-5 | Pentachlorophenol | 5.0 U |
| 85-01-8 | Phenanthrene | 1.0 U |
| 86-74-8 | Carbazole | 1.0 U |
| 120-12-7 | Anthracene | 1.0 U |
| 84-74-2 | Di-n-Butylphthalate | 1.0 U |
| 206-44-0 | Fluoranthene | 1.0 U |
| 129-00-0 | Pyrene | 1.0 U |
| 85-68-7 | Butylbenzylphthalate | 1.0 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 5.0 U |
| 56-55-3 | Benzo(a)anthracene | 1.0 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 1.0 U |
| 218-01-9 | Chrysene | 1.0 U |
| 117-84-0 | Di-n-Octyl phthalate | 1.0 U |
| 205-99-2 | Benzo(b)fluoranthene | 1.0 U |
| 207-08-9 | Benzo(k)fluoranthene | 1.0 U |
| 50-32-8 | Benzo(a)pyrene | 1.0 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 1.0 U |
| 53-70-3 | Dibenz(a,h)anthracene | 1.0 U |
| 191-24-2 | Benzo(g,h,i)perylene | 1.0 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 70.0% | d5-Phenol | 29.9% |
| 2-Fluorobiphenyl | 76.4% | 2-Fluorophenol | 45.3% |
| d14-p-Terphenyl | 78.0% | 2,4,6-Tribromophenol | 56.6% |
| d4-1,2-Dichlorobenzene | 69.2% | d4-2-Chlorophenol | 63.9% |

ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Page 1 of 2



ANALYTICAL
RESOURCES
INCORPORATED

Sample No: MW-1A-123

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547B
LIMS ID: 95-825
Matrix: Water
Data Release Authorized: *DM*
Reported: 02/08/95

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 2.0 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 2.0 U |
| 95-57-8 | 2-Chlorophenol | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 100-51-6 | Benzyl Alcohol | 5.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 95-48-7 | 2-Methylphenol | 2.0 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 1.0 U |
| 106-44-5 | 4-Methylphenol | 1.0 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 2.0 U |
| 67-72-1 | Hexachloroethane | 2.0 U |
| 98-95-3 | Nitrobenzene | 1.0 U |
| 78-59-1 | Isophorone | 1.0 U |
| 88-75-5 | 2-Nitrophenol | 5.0 U |
| 105-67-9 | 2,4-Dimethylphenol | 3.0 U |
| 65-85-0 | Benzoic Acid | 10 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 1.0 U |
| 120-83-2 | 2,4-Dichlorophenol | 3.0 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 U |
| 91-20-3 | Naphthalene | 1.0 U |
| 106-47-8 | 4-Chloroaniline | 3.0 U |
| 87-68-3 | Hexachlorobutadiene | 2.0 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 2.0 U |
| 91-57-6 | 2-Methylnaphthalene | 1.0 U |
| 77-47-4 | Hexachlorocyclopentadiene | 5.0 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 5.0 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 5.0 U |
| 91-58-7 | 2-Chloronaphthalene | 1.0 U |
| 88-74-4 | 2-Nitroaniline | 5.0 U |
| 131-11-3 | Dimethylphthalate | 1.0 U |
| 208-96-8 | Acenaphthylene | 1.0 U |
| 99-09-2 | 3-Nitroaniline | 6.0 U |
| 83-32-9 | Acenaphthene | 1.0 U |
| 51-28-5 | 2,4-Dinitrophenol | 10 U |
| 100-02-7 | 4-Nitrophenol | 5.0 U |
| 132-64-9 | Dibenzofuran | 1.0 U |
| 606-20-2 | 2,6-Dinitrotoluene | 5.0 U |

ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Page 2 of 2



ANALYTICAL
RESOURCES
INCORPORATED

Sample No: MW-1A-123

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547B
LIMS ID: 95-825
Matrix: Water
Data Release Authorized: *BBP*
Reported: 02/08/95

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 5.0 U |
| 84-66-2 | Diethylphthalate | 1.0 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 1.0 U |
| 86-73-7 | Fluorene | 1.0 U |
| 100-01-6 | 4-Nitroaniline | 5.0 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 10 U |
| 86-30-6 | N-Nitrosodiphenylamine | 1.0 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 1.0 U |
| 118-74-1 | Hexachlorobenzene | 1.0 U |
| 87-86-5 | Pentachlorophenol | 5.0 U |
| 85-01-8 | Phenanthrene | 1.0 U |
| 86-74-8 | Carbazole | 1.0 U |
| 120-12-7 | Anthracene | 1.0 U |
| 84-74-2 | Di-n-Butylphthalate | 1.0 U |
| 206-44-0 | Fluoranthene | 1.0 U |
| 129-00-0 | Pyrene | 1.0 U |
| 85-68-7 | Butylbenzylphthalate | 1.0 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 5.0 U |
| 56-55-3 | Benzo(a)anthracene | 1.0 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 1.0 U |
| 218-01-9 | Chrysene | 1.0 U |
| 117-84-0 | Di-n-Octyl phthalate | 1.0 U |
| 205-99-2 | Benzo(b)fluoranthene | 1.0 U |
| 207-08-9 | Benzo(k)fluoranthene | 1.0 U |
| 50-32-8 | Benzo(a)pyrene | 1.0 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 1.0 U |
| 53-70-3 | Dibenz(a,h)anthracene | 1.0 U |
| 191-24-2 | Benzo(g,h,i)perylene | 1.0 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 69.8% | d5-Phenol | 30.7% |
| 2-Fluorobiphenyl | 77.7% | 2-Fluorophenol | 45.8% |
| d14-p-Terphenyl | 80.3% | 2,4,6-Tribromophenol | 60.4% |
| d4-1,2-Dichlorobenzene | 71.9% | d4-2-Chlorophenol | 63.5% |

ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Page 1 of 2



ANALYTICAL
RESOURCES
INCORPORATED

Lab Sample ID: J547C
LIMS ID: 95-826
Matrix: Water
Data Release Authorized: *OKP*
Reported: 02/08/95

Sample No: MW-3-123

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 2.0 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 2.0 U |
| 95-57-8 | 2-Chlorophenol | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 100-51-6 | Benzyl Alcohol | 5.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 95-48-7 | 2-Methylphenol | 2.0 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 1.0 U |
| 106-44-5 | 4-Methylphenol | 1.0 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 2.0 U |
| 67-72-1 | Hexachloroethane | 2.0 U |
| 98-95-3 | Nitrobenzene | 1.0 U |
| 78-59-1 | Isophorone | 1.0 U |
| 88-75-5 | 2-Nitrophenol | 5.0 U |
| 105-67-9 | 2,4-Dimethylphenol | 3.0 U |
| 65-85-0 | Benzoic Acid | 10 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 1.0 U |
| 120-83-2 | 2,4-Dichlorophenol | 3.0 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 U |
| 91-20-3 | Naphthalene | 1.0 U |
| 106-47-8 | 4-Chloroaniline | 3.0 U |
| 87-68-3 | Hexachlorobutadiene | 2.0 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 2.0 U |
| 91-57-6 | 2-Methylnaphthalene | 1.0 U |
| 77-47-4 | Hexachlorocyclopentadiene | 5.0 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 5.0 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 5.0 U |
| 91-58-7 | 2-Chloronaphthalene | 1.0 U |
| 88-74-4 | 2-Nitroaniline | 5.0 U |
| 131-11-3 | Dimethylphthalate | 1.0 U |
| 208-96-8 | Acenaphthylene | 1.0 U |
| 99-09-2 | 3-Nitroaniline | 6.0 U |
| 83-32-9 | Acenaphthene | 1.0 U |
| 51-28-5 | 2,4-Dinitrophenol | 10 U |
| 100-02-7 | 4-Nitrophenol | 5.0 U |
| 132-64-9 | Dibenzofuran | 1.0 U |
| 606-20-2 | 2,6-Dinitrotoluene | 5.0 U |

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2



**ANALYTICAL
RESOURCES
INCORPORATED**

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547C

LIMS ID: 95-826

Matrix: Water

Data Release Authorized: *OBH*

Reported: 02/08/95

Sample No: MW-3-123

QC Report No: J547-Landau Associates, Incorporated

Project: 275002

Paxton Sales

Date Sampled: 01/23/95

Date Received: 01/24/95

Date extracted: 01/27/95

Date analyzed: 01/27/95

Instrument: FINN2

Sample Amount: 500 mL

Final Extract Volume: 0.5 mL

Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 5.0 U |
| 84-66-2 | Diethylphthalate | 1.0 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 1.0 U |
| 86-73-7 | Fluorene | 1.0 U |
| 100-01-6 | 4-Nitroaniline | 5.0 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 10 U |
| 86-30-6 | N-Nitrosodiphenylamine | 1.0 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 1.0 U |
| 118-74-1 | Hexachlorobenzene | 1.0 U |
| 87-86-5 | Pentachlorophenol | 5.0 U |
| 85-01-8 | Phenanthrene | 1.0 U |
| 86-74-8 | Carbazole | 1.0 U |
| 120-12-7 | Anthracene | 1.0 U |
| 84-74-2 | Di-n-Butylphthalate | 1.0 U |
| 206-44-0 | Fluoranthene | 1.0 U |
| 129-00-0 | Pyrene | 1.0 U |
| 85-68-7 | Butylbenzylphthalate | 1.0 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 5.0 U |
| 56-55-3 | Benzo(a)anthracene | 1.0 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 1.0 U |
| 218-01-9 | Chrysene | 1.0 U |
| 117-84-0 | Di-n-Octyl phthalate | 1.0 U |
| 205-99-2 | Benzo(b)fluoranthene | 1.0 U |
| 207-08-9 | Benzo(k)fluoranthene | 1.0 U |
| 50-32-8 | Benzo(a)pyrene | 1.0 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 1.0 U |
| 53-70-3 | Dibenz(a,h)anthracene | 1.0 U |
| 191-24-2 | Benzo(g,h,i)perylene | 1.0 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 67.2% | d5-Phenol | 38.8% |
| 2-Fluorobiphenyl | 77.3% | 2-Fluorophenol | 56.5% |
| d14-p-Terphenyl | 81.0% | 2,4,6-Tribromophenol | 60.4% |
| d4-1,2-Dichlorobenzene | 73.0% | d4-2-Chlorophenol | 72.2% |



ANALYTICAL
RESOURCES
INCORPORATED

ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Page 1 of 2

Sample No: MW-2-123

Analytical
Chemists &
Consultants

Lab Sample ID: J547D
LIMS ID: 95-827
Matrix: Water
Data Release Authorized: *DMH*
Reported: 02/08/95

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 2.0 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 2.0 U |
| 95-57-8 | 2-Chlorophenol | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 100-51-6 | Benzyl Alcohol | 5.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 95-48-7 | 2-Methylphenol | 2.0 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 1.0 U |
| 106-44-5 | 4-Methylphenol | 1.0 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 2.0 U |
| 67-72-1 | Hexachloroethane | 2.0 U |
| 98-95-3 | Nitrobenzene | 1.0 U |
| 78-59-1 | Isophorone | 1.0 U |
| 88-75-5 | 2-Nitrophenol | 5.0 U |
| 105-67-9 | 2,4-Dimethylphenol | 3.0 U |
| 65-85-0 | Benzoic Acid | 10 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 1.0 U |
| 120-83-2 | 2,4-Dichlorophenol | 3.0 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 U |
| 91-20-3 | Naphthalene | 1.0 U |
| 106-47-8 | 4-Chloroaniline | 3.0 U |
| 87-68-3 | Hexachlorobutadiene | 2.0 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 2.0 U |
| 91-57-6 | 2-Methylnaphthalene | 1.0 U |
| 77-47-4 | Hexachlorocyclopentadiene | 5.0 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 5.0 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 5.0 U |
| 91-58-7 | 2-Chloronaphthalene | 1.0 U |
| 88-74-4 | 2-Nitroaniline | 5.0 U |
| 131-11-3 | Dimethylphthalate | 1.0 U |
| 208-96-8 | Acenaphthylene | 1.0 U |
| 99-09-2 | 3-Nitroaniline | 6.0 U |
| 83-32-9 | Acenaphthene | 1.0 U |
| 51-28-5 | 2,4-Dinitrophenol | 10 U |
| 100-02-7 | 4-Nitrophenol | 5.0 U |
| 132-64-9 | Dibenzofuran | 1.0 U |
| 606-20-2 | 2,6-Dinitrotoluene | 5.0 U |

ORGANICS ANALYSIS DATA SHEET

Semivolatiles by GC/MS

Page 2 of 2



**ANALYTICAL
RESOURCES
INCORPORATED**

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547D
LIMS ID: 95-827
Matrix: Water
Data Release Authorized: *MM*
Reported: 02/08/95

Sample No: MW-2-123

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: 01/23/95
Date Received: 01/24/95

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 5.0 U |
| 84-66-2 | Diethylphthalate | 1.0 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 1.0 U |
| 86-73-7 | Fluorene | 1.0 U |
| 100-01-6 | 4-Nitroaniline | 5.0 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 10 U |
| 86-30-6 | N-Nitrosodiphenylamine | 1.0 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 1.0 U |
| 118-74-1 | Hexachlorobenzene | 1.0 U |
| 87-86-5 | Pentachlorophenol | 5.0 U |
| 85-01-8 | Phenanthrene | 1.0 U |
| 86-74-8 | Carbazole | 1.0 U |
| 120-12-7 | Anthracene | 1.0 U |
| 84-74-2 | Di-n-Butylphthalate | 1.0 U |
| 206-44-0 | Fluoranthene | 1.0 U |
| 129-00-0 | Pyrene | 1.0 U |
| 85-68-7 | Butylbenzylphthalate | 1.0 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 5.0 U |
| 56-55-3 | Benzo(a)anthracene | 1.0 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 1.0 U |
| 218-01-9 | Chrysene | 1.0 U |
| 117-84-0 | Di-n-Octyl phthalate | 1.0 U |
| 205-99-2 | Benzo(b)fluoranthene | 1.0 U |
| 207-08-9 | Benzo(k)fluoranthene | 1.0 U |
| 50-32-8 | Benzo(a)pyrene | 1.0 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 1.0 U |
| 53-70-3 | Dibenz(a,h)anthracene | 1.0 U |
| 191-24-2 | Benzo(g,h,i)perylene | 1.0 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 66.2% | d5-Phenol | 29.9% |
| 2-Fluorobiphenyl | 74.4% | 2-Fluorophenol | 43.7% |
| d14-p-Terphenyl | 81.7% | 2,4,6-Tribromophenol | 53.9% |
| d4-1,2-Dichlorobenzene | 71.1% | d4-2-Chlorophenol | 62.6% |

ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Page 1 of 2



ANALYTICAL
RESOURCES
INCORPORATED

Analytical
Chemists &
Consultants

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Lab Sample ID: J547MB
LIMS ID: 95-824
Matrix: Water
Data Release Authorized: *AB*
Reported: 02/08/95

Sample No: Method Blank

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: NA
Date Received: NA

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|------------------------------|-------|
| 108-95-2 | Phenol | 2.0 U |
| 111-44-4 | Bis-(2-Chloroethyl) Ether | 2.0 U |
| 95-57-8 | 2-Chlorophenol | 1.0 U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.0 U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.0 U |
| 100-51-6 | Benzyl Alcohol | 5.0 U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.0 U |
| 95-48-7 | 2-Methylphenol | 2.0 U |
| 108-60-1 | 2,2'-Oxybis(1-Chloropropane) | 1.0 U |
| 106-44-5 | 4-Methylphenol | 1.0 U |
| 621-64-7 | N-Nitroso-Di-N-Propylamine | 2.0 U |
| 67-72-1 | Hexachloroethane | 2.0 U |
| 98-95-3 | Nitrobenzene | 1.0 U |
| 78-59-1 | Isophorone | 1.0 U |
| 88-75-5 | 2-Nitrophenol | 5.0 U |
| 105-67-9 | 2,4-Dimethylphenol | 3.0 U |
| 65-85-0 | Benzoic Acid | 10 U |
| 111-91-1 | bis(2-Chloroethoxy) Methane | 1.0 U |
| 120-83-2 | 2,4-Dichlorophenol | 3.0 U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.0 U |
| 91-20-3 | Naphthalene | 1.0 U |
| 106-47-8 | 4-Chloroaniline | 3.0 U |
| 87-68-3 | Hexachlorobutadiene | 2.0 U |
| 59-50-7 | 4-Chloro-3-methylphenol | 2.0 U |
| 91-57-6 | 2-Methylnaphthalene | 1.0 U |
| 77-47-4 | Hexachlorocyclopentadiene | 5.0 U |
| 88-06-2 | 2,4,6-Trichlorophenol | 5.0 U |
| 95-95-4 | 2,4,5-Trichlorophenol | 5.0 U |
| 91-58-7 | 2-Chloronaphthalene | 1.0 U |
| 88-74-4 | 2-Nitroaniline | 5.0 U |
| 131-11-3 | Dimethylphthalate | 1.0 U |
| 208-96-8 | Acenaphthylene | 1.0 U |
| 99-09-2 | 3-Nitroaniline | 6.0 U |
| 83-32-9 | Acenaphthene | 1.0 U |
| 51-28-5 | 2,4-Dinitrophenol | 10 U |
| 100-02-7 | 4-Nitrophenol | 5.0 U |
| 132-64-9 | Dibenzofuran | 1.0 U |
| 606-20-2 | 2,6-Dinitrotoluene | 5.0 U |



**ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS**

Page 2 of 2

Lab Sample ID: J547MB
LIMS ID: 95-824
Matrix: Water
Data Release Authorized: *DM*
Reported: 02/08/95

Sample No: Method Blank

Analytical
Chemists &
Consultants

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales
Date Sampled: NA
Date Received: NA

333 Ninth Ave. North
Seattle, WA 98109-5187
(206) 621-6490
(206) 621-7523 (FAX)

Date extracted: 01/27/95
Date analyzed: 01/27/95
Instrument: FINN2

Sample Amount: 500 mL
Final Extract Volume: 0.5 mL
Dilution Factor: 1:1

| CAS Number | Analyte | ug/L |
|------------|----------------------------|-------|
| 121-14-2 | 2,4-Dinitrotoluene | 5.0 U |
| 84-66-2 | Diethylphthalate | 1.0 U |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 1.0 U |
| 86-73-7 | Fluorene | 1.0 U |
| 100-01-6 | 4-Nitroaniline | 5.0 U |
| 534-52-1 | 4,6-Dinitro-2-Methylphenol | 10 U |
| 86-30-6 | N-Nitrosodiphenylamine | 1.0 U |
| 101-55-3 | 4-Bromophenyl-phenylether | 1.0 U |
| 118-74-1 | Hexachlorobenzene | 1.0 U |
| 87-86-5 | Pentachlorophenol | 5.0 U |
| 85-01-8 | Phenanthrene | 1.0 U |
| 86-74-8 | Carbazole | 1.0 U |
| 120-12-7 | Anthracene | 1.0 U |
| 84-74-2 | Di-n-Butylphthalate | 1.0 U |
| 206-44-0 | Fluoranthene | 1.0 U |
| 129-00-0 | Pyrene | 1.0 U |
| 85-68-7 | Butylbenzylphthalate | 1.0 U |
| 91-94-1 | 3,3'-Dichlorobenzidine | 5.0 U |
| 56-55-3 | Benzo(a)anthracene | 1.0 U |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 1.0 U |
| 218-01-9 | Chrysene | 1.0 U |
| 117-84-0 | Di-n-Octyl phthalate | 1.0 U |
| 205-99-2 | Benzo(b)fluoranthene | 1.0 U |
| 207-08-9 | Benzo(k)fluoranthene | 1.0 U |
| 50-32-8 | Benzo(a)pyrene | 1.0 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 1.0 U |
| 53-70-3 | Dibenz(a,h)anthracene | 1.0 U |
| 191-24-2 | Benzo(g,h,i)perylene | 1.0 U |

Semivolatiles Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 69.7% | d5-Phenol | 47.5% |
| 2-Fluorobiphenyl | 75.3% | 2-Fluorophenol | 71.9% |
| d14-p-Terphenyl | 76.3% | 2,4,6-Tribromophenol | 65.0% |
| d4-1,2-Dichlorobenzene | 70.7% | d4-2-Chlorophenol | 81.3% |



ANALYTICAL
RESOURCES
INCORPORATED

Analytical
Chemists &
Consultants

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ORGANICS ANALYSIS DATA SHEET
Semivolatiles by GC/MS

Lab Sample ID: J547SB
LIMS ID: 95-824
Matrix: Water

QC Report No: J547-Landau Associates, Incorporated
Project: 275002
Paxton Sales

Data Release Authorized: *MM*
Reported: 02/08/95

| LABORATORY CONTROL SAMPLE CONSTITUENT | SPIKE VALUE | SPIKE ADDED | % RECOVERY |
|--|----------------|----------------|---------------|
| Phenol | 25.0 | 37.5 | 66.7% |
| 2-Chlorophenol | 32.0 | 37.5 | 85.3% |
| 1,4-Dichlorobenzene | 18.1 | 25.0 | 72.4% |
| N-Nitroso-Di-N-Propylamine | 22.4 | 25.0 | 89.6% |
| 1,2,4-Trichlorobenzene | 19.3 | 25.0 | 77.2% |
| 4-Chloro-3-methylphenol | 34.1 | 37.5 | 90.9% |
| Acenaphthene | 20.6 | 25.0 | 82.4% |
| 4-Nitrophenol | 23.3 | 37.5 | 62.1% |
| 2,4-Dinitrotoluene | 24.1 | 25.0 | 96.4% |
| Pentachlorophenol | 35.0 | 37.5 | 93.3% |
| Pyrene | 21.6 | 25.0 | 86.4% |

Spike Blank Surrogate Recovery

| | | | |
|------------------------|-------|----------------------|-------|
| d5-Nitrobenzene | 70.4% | d5-Phenol | 55.1% |
| 2-Fluorobiphenyl | 77.0% | 2-Fluorophenol | 72.4% |
| d14-p-Terphenyl | 76.9% | 2,4,6-Tribromophenol | 74.4% |
| d4-1,2-Dichlorobenzene | 69.0% | d4-2-Chlorophenol | 77.6% |

Values reported in ug/L