

Volume 2 of 2

Olalla Landfill Remedial Investigation/Feasibility Study

Prepared for
**Kitsap County Department of Public Works,
Solid Waste Division**

May 2014

Prepared by
Parametrix



Appendix A

Closure QA Documentation

APPENDIX D

**LANDFILL CLOSURE FINAL COVER
QUALITY ASSURANCE TESTING RESULTS
FROM PARAMETRIX, 1989**

OLALLA LANDFILL CLOSURE, Kitsap County, Washington
Final Cover Quality Assurance Testing Results

| Soil Sample/ Core Number | Test Location | In-place Density (PCF) | Max. Dry Density (PCF) | Percent Compaction (%) | In-Place Moisture Content (%) | Optimum Moisture (%) | Atterberg Limits | Permeability (CM/SEC) | Comments |
|-----------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|
| 1ST BARRIER LIFT | | | | | | | | | |
| S-7 / CORE A | GRID #1 | 124.7 | 125.7 | 99.2 | 10.5 | 9.8 | NP | 1.2E-07 | |
| S-8 / CORE J | GRID #2 | 122.7 | 124.1 | 98.9 | 9.4 | 11.0 | NP | 7.3E-08 | |
| S-9 | GRID #3 | 123.8 | 123.3 | 100+ | 10.4 | 10.2 | NP | | |
| S-10 | GRID #4 | 122.7 | 122.7 | 100.0 | 10.0 | 12.7 | NP | | |
| S-11 | GRID #5 | 125.0 | 117.0 | 100+ | 10.6 | 13.0 | NP | | |
| S-12 / CORE B | GRID #6 | 124.0 | 118.6 | 100+ | 10.6 | 11.8 | NP | 2.8E-07 | |
| S-13 | GRID #7 | 124.4 | 117.0 | 100+ | 10.9 | 14.5 | NP | | |
| S-14 | GRID #8 | 123.5 | 118.9 | 100+ | 11.9 | 13.5 | NP | | |
| S-15 / CORE C | GRID #9 | 124.0 | 120.8 | 100+ | 10.4 | 12.9 | NP | 1.7E-06 | |
| S-16 | GRID #10 | 117.6 | 114.5 | 100+ | 14.9 | 16.6 | NP | | |
| S-17 | GRID #11 | 120.3 | 121.7 | 98.8 | 14.1 | 14.2 | NP | | |
| S-18 / CORE I | GRID #12 | 116.0 | 120.6 | 96.2 | 14.8 | 13.3 | NP | 9.3E-08 | |
| S-19 | GRID #16 | 114.3 | 119.1 | 96.0 | 15.3 | 16.4 | NP | | |
| S-20 | GRID #17 | 113.4 | 119.8 | 94.7 | 15.8 | 15.0 | NP | | |
| S-21 / CORE D | GRID #24 | 120.8 | 122.0 | 99.0 | 14.0 | 9.9 | NP | 4.0E-08 | |
| S-22 / CORE F | GRID #25 | 114.9 | 119.6 | 96.1 | 14.9 | 13.9 | NP | 1.4E-07 | |

OLALLA LANDFILL CLOSURE, Kitsap County, Washington
Final Cover Quality Assurance Testing Results

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|-----------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|
| S-23 | GRID #23 | 115.9 | 123.1 | 94.2 | 10.9 | 12.4 | NP | | |
| S-24 | GRID #18 | 118.3 | 125.8 | 94.0 | 13.6 | 12.8 | NP | | |
| S-25 / CORE E | GRID #15 | 118.3 | 121.8 | 97.1 | 13.4 | 13.1 | NP | 5.5E-07 | |
| S-26 | GRID #14 | 115.3 | 118.0 | 97.7 | 13.3 | 15.2 | NP | | |
| S-27 | GRID #19 | 117.0 | 116.3 | 100+ | 14.6 | 13.7 | NP | | |
| S-28 / CORE G | GRID #22 | 114.2 | 118.3 | 96.5 | 14.9 | 13.7 | NP | 4.3E-08 | |
| S-29 | GRID #26 | 115.0 | 115.7 | 99.4 | 13.7 | 14.6 | NP | | |
| S-30 | GRID #21 | 116.8 | 125.0 | 93.4 | 14.0 | 12.0 | NP | | |
| S-31 / CORE H | GRID #20 | 114.0 | 120.4 | 94.7 | 15.0 | 13.0 | NP | 4.1E-07 | |
| S-32 | GRID #13 | 114.1 | 123.1 | 92.7 | 14.8 | 10.2 | NP | | |

OLALLA LANDFILL CLOSURE, Kitsap County, Washington
Final Cover Quality Assurance Testing Results

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|-----------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|
| 2ND BARRIER LIFT | | | | | | | | | |
| S-33 / CORE P | GRID #21 | 120.3 | 120.8 | 99.6 | 6.0 | 11.4 | NP | 6.6E-07 | |
| S-34 | GRID #20 | 119.0 | 127.6 | 93.3 | 7.8 | 9.8 | NP | | |
| S-35 | GRID #13 | 116.6 | 119.3 | 97.7 | 7.9 | 11.6 | NP | | |
| S-36 / CORE T | GRID #12 | 117.7 | 128.0 | 92.0 | 8.3 | 9.5 | NP | 1.4E-06 | |
| S-37 | GRID #26 | 118.3 | 124.2 | 95.2 | 11.5 | 12.2 | NP | | |
| S-38 | GRID #22 | 120.3 | 118.2 | 100+ | 10.5 | 14.3 | NP | | |
| S-39 / CORE M | GRID #19 | 110.7 | 121.0 | 91.5 | 8.4 | 10.4 | NP | 7.0E-07 | |
| S-40 | GRID #14 | 114.5 | 120.2 | 95.3 | 9.3 | 14.3 | NP | | |
| S-41 / CORE L | GRID #11 | 116.3 | 119.4 | 97.4 | 10.7 | 13.5 | NP | 7.8E-08 | |
| S-42 | GRID #5 | 116.5 | 122.0 | 95.5 | 10.9 | 11.7 | NP | | |
| S-43 / CORE K | GRID #4 | 114.6 | 119.0 | 96.3 | 11.6 | 12.0 | NP | 7.8E-07 | |
| S-44 / CORE O | GRID #3 | 118.7 | 120.9 | 98.2 | 10.0 | 13.2 | NP | 2.8E-07 | |
| S-45 | GRID #6 | 118.9 | 120.6 | 98.6 | 11.6 | 13.2 | NP | | |
| S-46 | GRID #10 | 123.7 | 119.8 | 100+ | 10.5 | 11.1 | NP | | |
| S-47 | GRID #15 | 116.9 | 120.3 | 97.2 | 10.0 | 14.3 | NP | | |

OLALLA LANDFILL CLOSURE, Kitsap County, Washington
Final Cover Quality Assurance Testing Results

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|-----------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|
| S-48 | GRID #18 | 114.5 | 121.8 | 94.0 | 11.5 | 15.4 | NP | | |
| S-49 / CORE N | GRID #23 | 112.9 | 121.5 | 92.9 | 13.6 | 14.1 | NP | 3.1E-07 | |
| S-50 | GRID #25 | 114.0 | 119.3 | 95.6 | 10.0 | 13.4 | NP | | |
| S-51 / CORE Q | GRID #8 | 113.5 | 123.5 | 91.9 | 11.4 | 10.6 | NP | 8.6E-07 | |
| S-52 | GRID #1 | 109.4 | 120.1 | 91.1 | 14.3 | 13.9 | NP | | |
| S-53 | GRID #2 | 109.8 | 121.5 | 90.4 | 12.2 | 11.7 | NP | | |
| S-54 | GRID #7 | 111.7 | 123.4 | 90.5 | 12.4 | 15.0 | NP | | |
| S-55 / CORE R | GRID #9 | 109.7 | 121.6 | 90.2 | 11.5 | 12.0 | NP | 3.0E-07 | |
| S-56 | GRID #16 | 110.8 | 123.0 | 90.1 | 11.7 | 12.4 | NP | | |
| S-57 | GRID #17 | 109.7 | 122.0 | 89.9 | 11.7 | 11.1 | NP | | |
| S-58 / CORE S | GRID #24 | 110.3 | 119.8 | 92.1 | 13.0 | 14.6 | NP | 3.7E-07 | |

OLALLA LANDFILL CLOSURE, Kitsap County, Washington
Final Cover Quality Assurance Testing Results

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|-----------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|
| 3RD BARRIER LIFT | | | | | | | | | |
| S-59 | GRID #8 | 117.1 | 119.3 | 98.2 | 15.6 | 11.7 | NP | | |
| S-60 | GRID #1 | 116.4 | 121.0 | 96.2 | 16.3 | 14.5 | NP | | |
| S-61 | GRID #2 | 114.8 | 123.9 | 92.7 | 16.7 | 14.2 | NP | | |
| S-62 / CORE U | GRID #7 | 116.5 | 123.7 | 94.2 | 16.4 | 12.4 | NP | 6.7E-08 | |
| S-63 | GRID #9 | 118.5 | 125.0 | 94.8 | 15.3 | 14.8 | NP | | |
| S-64 / CORE V | GRID #16 | 111.3 | 122.6 | 90.8 | 15.7 | 15.9 | NP | 1.9E-08 | |
| S-65 / CORE d | GRID #17 | 112.7 | 124.3 | 90.7 | 15.1 | 14.7 | NP | 3.5E-07 | |
| S-66 | GRID #24 | 119.7 | 123.5 | 96.9 | 15.1 | 13.2 | NP | | |
| S-67 | GRID #25 | 112.4 | 123.4 | 91.1 | 17.0 | 13.4 | NP | | |
| S-68 | GRID #23 | 116.6 | 120.8 | 96.5 | 17.1 | 15.3 | NP | | |
| S-69 / CORE W | GRID #18 | 124.4 | 123.9 | 100+ | 15.9 | 12.1 | NP | 4.5E-07 | |
| S-70 | GRID #15 | 115.6 | 123.5 | 93.6 | 16.2 | 14.2 | NP | | |
| S-71 / CORE X | GRID #10 | 117.1 | 121.4 | 96.5 | 14.1 | 13.6 | NP | 1.6E-07 | |
| S-72 | GRID #6 | 115.8 | 122.3 | 94.7 | 15.4 | 13.2 | NP | | |
| S-73 | GRID #3 | 114.7 | 120.3 | 95.3 | 13.8 | 14.0 | NP | | |

OLALLA LANDFILL CLOSURE, Kitsap County, Washington
Final Cover Quality Assurance Testing Results

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|-----------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|---------|
| S-74 | GRID #4 | 112.4 | 121.1 | 92.8 | 13.1 | 13.5 | NP | | | |
| S-75 / CORE Y | GRID #5 | 111.1 | 123.5 | 90.0 | 15.1 | 15.1 | NP | 1.4E-07 | | |
| S-76 | GRID #11 | 110.5 | 121.3 | 91.1 | 15.4 | 13.9 | NP | | | |
| S-77 / CORE Z | GRID #14 | 114.1 | 125.1 | 91.2 | 14.9 | 12.1 | NP | 8.5E-08 | | |
| S-78 | GRID #19 | 110.7 | 120.1 | 92.2 | 16.3 | 13.7 | NP | | | |
| S-79 | GRID #22 | 116.7 | 120.6 | 96.8 | 15.9 | 13.0 | NP | | | |
| S-80 / CORE a | GRID #26 | 118.3 | 121.7 | 97.2 | 16.5 | 14.8 | NP | 5.0E-09 | | |
| S-81 | GRID #21 | 114.3 | 125.1 | 91.4 | 13.3 | 15.1 | NP | | | |
| S-82 | GRID #20 | 114.0 | 123.1 | 92.6 | 13.3 | 14.1 | NP | | | |
| S-83 | GRID #13 | 113.9 | 121.2 | 94.0 | 12.7 | 13.7 | NP | | | |
| S-84 | GRID #12 | 119.0 | 121.8 | 97.7 | 15.3 | 13.4 | NP | | | |
| CORE b | GRID #21/26 | | | | | | | 6.7E-07 | | |
| CORE c | GRID #24/25 | | | | | | | 1.8E-07 | | |
| FINAL COVER PERMEABILITY: | | | | | | | | MEAN DENSITY | ==> | 3.8E-07 |

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|-----------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|
|-----------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|

BANK AND DRAINAGE CONSTRUCTION

1ST BARRIER LIFT

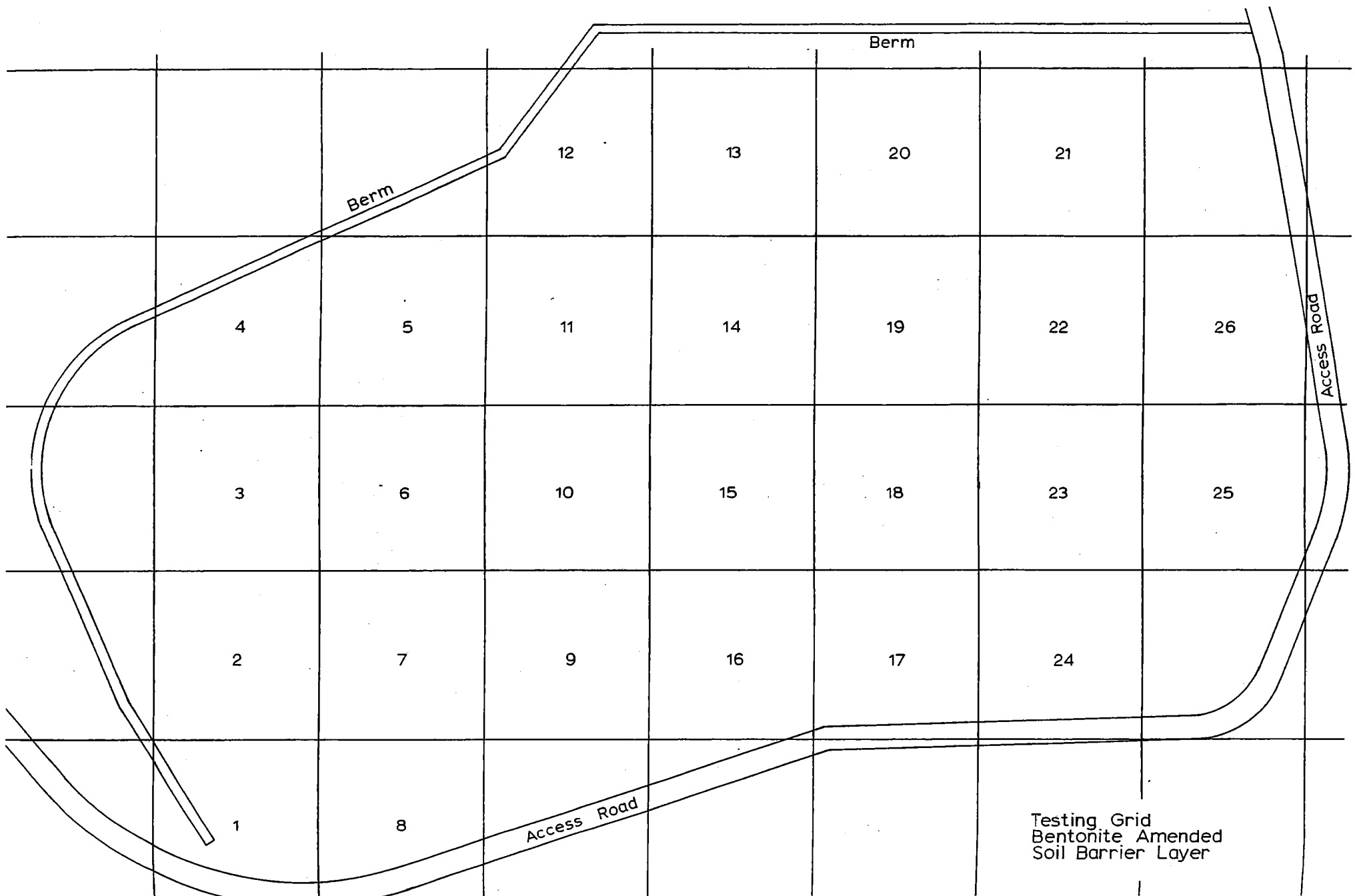
| | | | | | | | | |
|---|--|-------|-------|------|-----|------|-----|---------|
| DRAINAGE DITCH, E. GRID #3 WEST BANK | | 117.9 | 120.5 | 97.8 | 5.1 | 13.1 | NP | |
| DRAINAGE DITCH, S. GRID #8 SOUTH BANK | | 128.0 | 120.5 | 100+ | 4.2 | 13.1 | NP | |
| DRAINAGE DITCH, S. GRID #24 NORTH BANK | | 121.7 | 120.5 | 100+ | 4.2 | 13.1 | NP | |
| DRAINAGE DITCH, W. GRID #21 EAST BANK | | 118.4 | 120.5 | 98.3 | 4.4 | 13.1 | NP | |
| DRAINAGE DITCH, N. GRID #13 NORTH BANK | | 119.3 | 120.5 | 99.0 | 4.6 | 13.3 | NP | |
| DRAINAGE DITCH GRID #5 CORE DDI | | N/A | N/A | N/A | N/A | N/A | N/A | 2.2E-05 |

2ND BARRIER LIFT

| | | | | | | | | |
|-------------------------|--|-------|-------|------|------|------|----|--|
| DRAINAGE DITCH GRID #12 | | 117.3 | 121.5 | 96.5 | 14.6 | 12.6 | NP | |
| DRAINAGE DITCH GRID #4 | | 115.8 | 121.5 | 95.3 | 12.2 | 12.6 | NP | |
| DRAINAGE DITCH GRID #1 | | 118.3 | 121.5 | 97.4 | 11.9 | 12.6 | NP | |
| DRAINAGE DITCH GRID #16 | | 116.9 | 121.5 | 96.2 | 13.4 | 12.6 | NP | |

OLALLA LANDFILL CLOSURE, Kitsap County, Washington
Final Cover Quality Assurance Testing Results

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|------------------------------|------------------|------------------------------|------------------------------|------------------------------|-------------------------------------|----------------------------|---------------------|--------------------------|----------|
| DRAINAGE DITCH | GRID #25 | 114.4 | 121.5 | 94.2 | 14.7 | 12.6 | NP | | |
| DRAINAGE DITCH CORE DDII | GRID #21 | N/A | N/A | N/A | N/A | N/A | N/A | 1.0E-06 | |
| <u>3RD BARRIER LIFT</u> | | | | | | | | | |
| DRAINAGE DITCH | GRID #12 | 114.8 | 122.4 | 93.8 | 15.6 | 13.9 | NP | | |
| DRAINAGE DITCH | GRID #4 | 117.0 | 122.4 | 95.6 | 16.0 | 13.9 | NP | | |
| DRAINAGE DITCH | GRID #2 | 115.4 | 122.4 | 94.3 | 14.7 | 13.9 | NP | | |
| DRAINAGE DITCH | GRID #9 | 119.3 | 122.4 | 97.5 | 15.1 | 13.9 | NP | | |
| DRAINAGE DITCH | GRID #24 | 117.2 | 122.4 | 95.8 | 15.3 | 13.9 | NP | | |
| DRAINAGE DITCH CORE DDIII | GRID #17 | N/A | N/A | N/A | N/A | N/A | N/A | 1.9E-07 | |



Appendix B

Solid Waste Handling Permits

KITSAP PUBLIC HEALTH DISTRICT

2012

SOLID WASTE HANDLING PERMIT

Issued in accordance with the provisions of Chapter 173-304 of the Washington Administrative Code (WAC) and Kitsap County Board of Health Ordinance 2010-1 "Solid Waste Regulations."

PERMIT PERIOD: JANUARY 1, 2012- DECEMBER 31, 2012

SECTION I. PERMITTEE AND ADMINISTRATION INFORMATION

| | |
|---------------------------|--|
| FACILITY TYPE: | Mixed Municipal Solid Waste Landfill (Closed Under Chapter 173-304 WAC in May 1989) |
| ANNUAL PERMIT FEE: | \$109/hr, not to exceed \$10,000/yr during MTCA |
| FACILITY NAME: | Olalla Landfill |
| FACILITY LOCATION: | Approximately 3/4 mile east of Hwy. 16 on Burley-Olalla Rd. |
| FACILITY OWNER: | Kitsap County Department of Public Works |
| CONTACT PERSON: | Ms. Keli McKay-Means |
| TITLE: | Landfill Specialist |
| OWNER ADDRESS: | 614 Division Street, MS-27, Port Orchard, WA 98366 |
| OWNER PHONE: | (360) 337-5665 |

This permit is the property of the Kitsap Public Health District and may be suspended or revoked, after opportunity for hearing, upon violation of any rules and regulations applicable hereto. This permit is not transferable and must be renewed annually. The terms of this permit remain in effect until the date the next permit is issued.



Janet Brower
Program Manager
Solid and Hazardous Waste Program

6/1/2012

Date

File Location: [jb/swwqbcd/shw/common/permits/landfills/Olalla](#)

Date Prepared: January 12, 2012

Date Issued: June 1, 2012

SECTION II. STANDARD PERMIT CONDITIONS

- A. Terms pertaining to solid waste used in this permit shall be as defined per Kitsap County Board of Health (KCBH) Ordinance 2010-1, in Section 100, as amended.
- B. All conditions of this permit shall be binding upon, and the permittee shall be responsible for, all acts and omissions of all contractors and agents of the permittee for the term of the permit and post-closure period.
- C. The permittee shall notify the Kitsap Public Health District (Health District) in writing in advance of any alterations or changes planned in the post-closure plan, when these changes are different or absent from this permit. All alterations or additions to the maintenance and closure of the facility must be included in written plans approved by the Health District before implementation.
- D. Facility maintenance during post-closure shall be conducted in strict compliance with KCBH Ordinance 2010-1, Solid Waste Regulations, (Chapter 173-304 WAC adopted by reference) as amended, and the specific conditions of this permit. In the event of conflict between state and local regulations, the more stringent requirement shall apply.
- E. The Health Officer or his/her duly authorized inspector, or any duly authorized representative of the Washington State Department of Ecology, may, upon giving a minimum 3-day notice prior to entry, enter and inspect any building, structure, property, or portion thereof, at any reasonable time for the purpose of determining compliance with the provisions of this permit, or the provisions of other applicable rules and regulations, or whenever the Health Officer has reason to believe that a violation of this permit has been or is being committed. Entry shall not be unreasonably denied by the permittee, or his/her agent, but may be conditioned on the permittee or an agent of the permittee escorting the inspector, with said escort to be provided immediately upon request. Wet weather inspections conducted during significant rain events will be conducted with prior notice, but not necessarily providing a minimum 3-day notice depending upon conditions.
- F. The permittee shall report all instances of non-compliance (except regulatory exceedances which are outlined in Section IV.E.2.b) in writing within seven (7) days from time of discovery. Instances of non-compliance which may result in potential public health risk and/or environmental damage shall be reported immediately.
- G. This permit or copy thereof shall be located where it can be readily referred to by oversight personnel.
- H. This permit may be amended by the Health District. Amendments shall be made in writing and become specific conditions of the permit. Proposed amendments will be forwarded to the permittee for a fifteen (15) day review and comment period, unless otherwise waived in writing by the Health District. Upon consideration of the permittee's comments, the final

amendment will be issued and become effective. The permittee review and comment period may be waived by the Health District in emergencies where there is an imminent threat to human health or the environment.

SECTION III. FEES

- A. The permittee shall pay an annual permit fee as a condition of permit issuance. As provided for in WAC 173-304-600(2)(e), the Health District may establish reasonable fees for permits and permit renewals. The current annual fee for this permit is shown in Section I. This fee is an estimate of both routine and non-routine activities. The annual permit fee covers the following Health District activities and is based on the time expended by the Health District billed at the approved billing rate in the current fee schedule as provided for in Health District's current Fee Schedule.
1. Reviewing/commenting on three Quarterly Monitoring Reports.
 2. Reviewing/commenting on one Annual Report.
 3. Meet with permit holder annually to discuss the status of the facility and, *to the extent possible, identify proposed Health District activities or requirements related to the facility that would require fees in addition to the Annual Permit Fee.*
 4. Permit drafting/issuance.
 5. Plan and/or design review.
 6. Four site inspections and associated inspection reports.
 7. Administrative costs including routine communications, supervisor review of documents, and billing invoices.
 8. Health District oversight and review required as a result of the Health District's determination that the operation is in non-compliance with this permit and/or applicable regulations;
 9. Health District participation in MTCA review, comment, and meetings through 2012;
 10. Amendments to the existing permit;
 11. Report, plan, or design review associated with any corrective action or to correct non-compliance at the facility; or
 12. Other permit related work beyond those items listed in Section III.A. above.
- B. The annual permit fee will be invoiced semi-annually. Whenever a fee is assessed a Health District billing invoice will be sent to the permittee at the address provided by the permittee. Each invoice will be accompanied by a Detail Report that shows each entry by each staff member and a description of the activity being billed. A copy of the billing invoice with a check or money order for the assessed amount shall be sent by the permittee to the Health District and within thirty (30) days of receipt. Payment and a copy of the invoice shall be sent to:

Kitsap Public Health District
Accounting Office

345 6th Street, Suite 300
Bremerton, WA 98337-1866

- C. The total annual fee for 2012 shall not exceed \$10,000.00, billed at the established Health District hourly rate.

SECTION IV. SPECIFIC PERMIT CONDITIONS

A. Plan of Maintenance and Post-Closure

1. The permittee shall provide post-closure activities to allow for continued facility maintenance and monitoring of air, land, and ground and surface waters, as long as necessary for the facility to stabilize, to protect human health and the environment as determined by the Health District. These post-closure activities shall be conducted in strict compliance with Chapter 173-304 WAC, KCBH Ordinance 2010-1, and the conditions of this permit.
2. Post-closure activities include, but are not limited to, the following: ground water monitoring; surface water monitoring; landfill gas monitoring; and proper operation and maintenance of the facility, facility structures, and facility monitoring systems for their intended use for a period of no less than twenty years from the date of final closure (May 1989).
3. The permittee shall close and maintain the landfill in accordance with the regulations stipulated in Section II.D. of this permit and in conformance with the Health District approved post-closure plan "Olalla Landfill Final Closure Plan," June 1988, as amended, which includes, but is not limited to the actions specified in Section IV.2. above. Closure activities not in conformance with these regulations and this plan shall be considered in violation of this permit.
4. The Health District may determine that the facility post-closure plan is invalid and require the permittee to amend the facility post-closure plan for approval by the Health District.
5. When the Health District determines that a facility post-closure plan amendment is required, the Health District shall, after consultation with the permittee, designate a compliance schedule for submittal of the amendment and its review and approval.
6. The Health District may direct facility post-closure activities to cease, in part or whole, until the post-closure plan amendment has received written approval by the Health District.
7. When post-closure activities are complete, the permittee shall certify to the Health District that the facility is stabilized and state why post-closure activities are no longer necessary. The certification shall be signed by the permittee and a professional engineer registered in the state of Washington.

- B. Minimum Functional Standards for Maintenance.** The permittee shall provide all activities necessary to allow for continued facility maintenance. These activities include, but are not limited to:
1. Stormwater quantity and quality control;
 2. Slope stability, erosion and dust control;
 3. Maintenance of access roads and ditches;
 4. Maintenance of facility structures and systems (i.e., stormwater, landfill cover, gas control, surface and ground water monitoring, etc.);
 5. Control and minimization, or elimination, of threats to human health and environment; and
 6. Unauthorized entry at the facility shall be controlled, at minimum, by means of a lockable gate, barrier, fence, etc., at the property boundary.
- C. Minimum Standards for Performance.** The permittee shall provide all activities necessary to allow for the continued monitoring of air, land, and water as long as is necessary for the facility to stabilize and to protect human health and environment.
1. **Ground Water.** The facility shall not contaminate the ground water underlying the landfill, beyond the point of compliance. The facility shall also not cause exceedances of standards contained in Chapter 173-200 WAC (Water Quality Standards for Ground Waters of the State of Washington) or Chapter 246-290 WAC (Drinking Water Regulations.)
 2. **Air Quality.**
 - a. The facility shall not allow explosive gases to exceed:
 - 1) Twenty-five percent of the lower explosive limit for the gases in facility structures (excluding gas control or recovery system components).
 - 2) The lower explosive limit for the gases at the property boundary or beyond.
 - 3) One hundred parts per million by volume of hydrocarbons (expressed as methane) in off-site structures.
 - b. The facility shall not cause a violation of any ambient air quality standard at the property boundary or emission standard from any emission of landfill gases, combustion, or any other emission associated with a landfill.
 - c. The permittee shall not allow non-burning flares to vent openly to the atmosphere unless it can be shown that little or no landfill gases will be produced or that landfill gases will not support combustion. In such cases installation of vents shall be required as stipulated in Chapter 173-304-460(3)(f)(ii).

3. **Surface Water.** The facility shall not cause exceedances of standards contained in Chapter 173-201A WAC (Water Quality Standards for Surface Waters of the State of Washington) or violate Chapter 90.48 RCW through the discharges of surface runoff, leachate, or any other liquid associated with a landfill.

D. Minimum Environmental Monitoring Requirements.

1. The minimum environmental monitoring requirements specified herein are subject to revision by the Health District based on the Health District's review of reports and plans required through this permit. All environmental monitoring shall be performed in accordance with the approved Sampling and Analysis Plan for the facility, as amended.
2. **Ground Water Monitoring.**
 - a. **Quarterly Monitoring.** Minimum quarterly monitoring required from monitoring wells MW-1, MW-3, MW-6, MW-8, and MW-10. Each individual monitoring well shall be analyzed for:

| | |
|---|---------------------------------|
| Water Level | Temperature |
| Specific Conductance | pH |
| Total Coliform | Sulfate |
| Potassium | Bicarbonate |
| Carbonate | Sodium |
| Chloride | Calcium |
| Total Organic Carbon | Nitrite, Nitrate, & Ammonia (N) |
| Dissolved Arsenic | Chemical Oxygen Demand |
| Dissolved Iron | Dissolved Barium |
| Dissolved Zinc | Dissolved Manganese |
| Volatile Organic Compounds (EPA Method 8260C) | |

Annual Monitoring. A reduced frequency of monitoring for MW-5A and MW-7. A reduced frequency of monitoring consisting of an annual sample collected during the fourth Quarter Sampling event shall be analyzed for:

Water Level
Specific Conductance
Dissolved Arsenic
Dissolved Iron
Dissolved Manganese
Vinyl Chloride
Temperature
pH

- b. Ground water analyses shall be performed by a laboratory accredited by the Washington State Department of Ecology. The methods of analysis shall be in accordance with EPA Publication Number SW-846, "Test Methods for Evaluating Solid Waste - Physical/Chemical Methods". For each test method, the lowest achievable detection limit shall be obtained sufficient to verify compliance with the regulatory standards in Section IV.C.1. of this permit.

3. Surface Water Monitoring.

- a. Annual Monitoring. Minimum annual monitoring required from surface water station SW-2. The annual event shall occur either between January and March or November and December, and shall be analyzed for:

| | | |
|----------------------|----------------|------------------|
| Temperature | pH | Nitrate-Nitrogen |
| Specific Conductance | Fecal Coliform | |

- b. Surface water analysis shall be performed by a laboratory accredited by the Washington State Department of Ecology. The methods of analysis shall be in accordance with 40 CFR Part 136. For each test method, the lowest achievable detection limit shall be obtained sufficient to verify compliance with the standards in Section IV.C.3. of this permit.

- 4. Landfill Gas Monitoring. Minimum quarterly monitoring of landfill gas is required from gas flares GF-1, GF-2, and GF-3. Individual samples shall be monitored for the following:

| | |
|----------------|-----------------------|
| Methane (%LEL) | Gas Pressure |
| Oxygen (%VOL) | Carbon Dioxide (%VOL) |

- E. Reporting Requirements. All reports specified herein must be forwarded to the Health District representative responsible for this permit, and to the Solid Waste Supervisor, Department of Ecology Northwest Regional Office, Waste 2 Resources, 3190 160th Avenue S.E., Bellevue, WA 98008-5452.

1. Annual Report.

- a. Annual report due April 1, following each calendar year.
- b. The permittee shall prepare and submit one copy and one CD of the annual report to the Health District, and one copy and one CD to Ecology. The annual report will serve as the permit application for the next year's permit. The annual report shall cover facility activities during the previous calendar year and must include, at a minimum, the following information:

- 1) Name and address of the facility.
 - 2) Calendar year covered by the report.
 - 3) A summary of all ground water, surface water, and landfill gas Monitoring will be provided in accordance with the Revised Statistical Approach for Olalla Landfill (June 2008) . The summary shall contain the results of applicable statistical analyses, explanations for increasing/decreasing contaminant level trends, confidence intervals and if outliers were discarded, a justification for this action. In addition time series plots for each assessment monitoring well for each parameter that exceeded groundwater quality standards at any time since well installation. All statistically significant increasing/decreasing contaminant level trends and exceedances of ground water, surface water, or air quality standards as described in Section IV shall be reported and explained to the extent possible, and if deemed necessary, a summary of actions taken to verify results. Additionally, if the trends shows either no improvement or increasing contaminate levels or exceedances are detected at or beyond the point of compliance, a plan of action shall be included which describes what activities will be taken to correct the statically significantly elevated contaminants or exceedances, and the expected time frame required to do so. The summary shall also contain a description of the ground water flow rate and direction in the uppermost aquifer. Any changes in the number or location of monitoring systems shall also be noted and explained.
 - 4) A description of the next year's proposed environmental monitoring program, operations, and maintenance activities.
 - 5) A summary of all activities related to the inspection, maintenance, and engineering of the facility and its systems. All field data sheets, inspection and log forms shall be made available to the Health District upon request.
 - 6) An explanation of causes and effects of all instances of non-compliance with the provisions of this permit and a summary of how these issues were or are being corrected.
2. Quarterly Environmental Monitoring Reports.
- a. The permittee shall prepare and submit written quarterly environmental monitoring reports to report the information gathered under the "Minimum Environmental Monitoring Requirements" section of this permit. Gas, surface and ground water monitoring data shall be summarized, evaluated, and compared with all applicable standards, and shall also be submitted in hard copy form. All monitoring data shall be reported in detail on a quarterly basis. Results and

summary of the fourth quarter sampling event shall be combined with the annual report.

- b. Environmental monitoring reports shall be submitted to the Health District and Ecology within sixty (60) days of the date of sample collection. If the report cannot be submitted within sixty (60) days, the Health District will be notified prior to the end of sixty (60) days as to the status of the report and when it will be submitted. In no event shall monitoring reports be submitted more than seventy (70) days after the date of sample collection. All instances of regulatory standard exceedances shall be faxed to the Health District within ten (10) days of the time of discovery.

SECTION V. CORRECTIVE ACTION PLAN

| TASKS | Submittal Date |
|---|-----------------------|
| 2. MTCA PROCESS (see attached schedule) | Ongoing |
| | |
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KITSAP COUNTY HEALTH DISTRICT

2010-2011 SOLID WASTE HANDLING PERMIT

Issued in accordance with the provisions of Chapter 173-304 of the Washington Administrative Code (WAC) and Kitsap County Board of Health Ordinance 2004-2 "Solid Waste Regulations."

PERMIT PERIOD: JANUARY 1, 2010 - DECEMBER 31, 2011

SECTION I. PERMITTEE AND ADMINISTRATION INFORMATION

| | |
|--------------------|--|
| FACILITY TYPE: | Mixed Municipal Solid Waste Landfill (Closed Under Chapter 173-304 WAC in May 1989) |
| ANNUAL PERMIT FEE: | \$109/hr, not to exceed \$10,000 with MTCA |
| FACILITY NAME: | Olalla Landfill |
| FACILITY LOCATION: | Approximately 3/4 mile east of Hwy. 16 on Burley-Olalla Rd. |
| FACILITY OWNER: | Kitsap County Department of Public Works |
| CONTACT PERSON: | Ms. Keli McKay-Means |
| TITLE: | Landfill Specialist |
| OWNER ADDRESS: | 614 Division Street, MS-27, Port Orchard, WA 98366 |
| OWNER PHONE: | (360) 337-5665 |

This permit is the property of the Kitsap County Health District and may be suspended or revoked, after opportunity for hearing, upon violation of any rules and regulations applicable hereto. This permit is not transferable and must be renewed at the end of the permit period.



Janet Brower
Program Manager
Solid and Hazardous Waste Program

4/19/2010
Date

File Location: jb/swwqbcd/shw/common/permits/landfills/Olalla
Date Prepared: April 19, 2009

SECTION II. STANDARD PERMIT CONDITIONS

- A. Terms pertaining to solid waste used in this permit shall be as defined per Kitsap County Board of Health (KCBH) Ordinance 2004-2, in Section 100, as amended.
- B. All conditions of this permit shall be binding upon, and the permittee shall be responsible for, all acts and omissions of all contractors and agents of the permittee for the term of the permit and post-closure period.
- C. The permittee shall notify the Kitsap County Health District (Health District) in writing in advance of any alterations or changes planned in the post-closure plan, when these changes are different or absent from this permit. All alterations or additions to the maintenance and closure of the facility must be included in written plans approved by the Health District before implementation.
- D. Facility maintenance during post-closure shall be conducted in strict compliance with KCBH Ordinance 2004-2, Solid Waste Regulations, (Chapter 173-304 WAC adopted by reference) as amended, and the specific conditions of this permit. In the event of conflict between state and local regulations, the more stringent requirement shall apply.
- E. The Health Officer or his/her duly authorized inspector, or any duly authorized representative of the Washington State Department of Ecology, may, upon giving a minimum 3-day notice prior to entry, enter and inspect any building, structure, property, or portion thereof, at any reasonable time for the purpose of determining compliance with the provisions of this permit, or the provisions of other applicable rules and regulations, or whenever the Health Officer has reason to believe that a violation of this permit has been or is being committed. Entry shall not be unreasonably denied by the permittee, or his/her agent, but may be conditioned on the permittee or an agent of the permittee escorting the inspector, with said escort to be provided immediately upon request. Wet weather inspections conducted during significant rain events will be conducted with prior notice, but not necessarily providing a minimum 3-day notice depending upon conditions.
- F. The permittee shall report all instances of non-compliance (except regulatory exceedances which are outlined in Section IV.E.2.b) in writing within seven (7) days from time of discovery. Instances of non-compliance which may result in potential public health risk and/or environmental damage shall be reported immediately.
- G. This permit or copy thereof shall be located where it can be readily referred to by oversight personnel.
- H. This permit may be amended by the Health District. Amendments shall be made in writing and become specific conditions of the permit. Proposed amendments will be forwarded to the permittee for a fifteen (15) day review and comment period, unless otherwise waived in writing by the Health District. Upon consideration of the permittee's comments, the final

amendment will be issued and become effective. The permittee review and comment period may be waived by the Health District in emergencies where there is an imminent threat to human health or the environment.

SECTION III. FEES

- A. The permittee shall pay an annual permit fee as a condition of permit issuance. As provided for in WAC 173-304-600(2)(e), the Health District may establish reasonable fees for permits and permit renewals. The current annual fee for this permit is shown in Section I. This fee is an estimate of both routine and non-routine activities. The annual permit fee covers the following Health District activities and is based on the time expended by the Health District billed at the approved billing rate in the current fee schedule as provided for in Health District's current Fee Schedule.
1. Reviewing/commenting on three Quarterly Monitoring Reports.
 2. Reviewing/commenting on one Annual Report.
 3. Meet with permit holder annually to discuss the status of the facility and, *to the extent possible, identify proposed Health District activities or requirements related to the facility that would require fees in addition to the Annual Permit Fee.*
 4. Permit drafting/issuance.
 5. Plan and/or design review.
 6. Four site inspections and associated inspection reports.
 7. Administrative costs including routine communications, supervisor review of documents, and billing invoices.
 8. Health District oversight and review required as a result of the Health District's determination that the operation is in non-compliance with this permit and/or applicable regulations;
 9. Health District participation in MTCA review, comment, and meetings through 2011;
 10. Amendments to the existing permit;
 11. Report, plan, or design review associated with any corrective action or to correct non-compliance at the facility; or
 12. Other permit related work beyond those items listed in Section III.A. above.
- B. The annual permit fee will be invoiced semi-annually. Whenever a fee is assessed a Health District billing invoice will be sent to the permittee at the address provided by the permittee. Each invoice will be accompanied by a Detail Report that shows each entry by each staff member and a description of the activity being billed. A copy of the billing invoice with a check or money order for the assessed amount shall be sent by the permittee to the Health District and within thirty (30) days of receipt. Payment and a copy of the invoice shall be sent to:

Kitsap County Health District
Accounting Office
345 6th Street, Suite 300
Bremerton, WA 98337-1866

SECTION IV. SPECIFIC PERMIT CONDITIONS

A. Plan of Maintenance and Post-Closure

1. The permittee shall provide post-closure activities to allow for continued facility maintenance and monitoring of air, land, and ground and surface waters, as long as necessary for the facility to stabilize, to protect human health and the environment as determined by the Health District. These post-closure activities shall be conducted in strict compliance with Chapter 173-304 WAC, KCBH Ordinance 2004-2, and the conditions of this permit.
2. Post-closure activities include, but are not limited to, the following: ground water monitoring; surface water monitoring; landfill gas monitoring; and proper operation and maintenance of the facility, facility structures, and facility monitoring systems for their intended use for a period of no less than twenty years from the date of final closure (May 1989).
3. The permittee shall close and maintain the landfill in accordance with the regulations stipulated in Section II.D. of this permit and in conformance with the Health District approved post-closure plan "Olalla Landfill Final Closure Plan," June 1988, as amended, which includes, but is not limited to the actions specified in Section IV.2. above. Closure activities not in conformance with these regulations and this plan shall be considered in violation of this permit.
4. The Health District may determine that the facility post-closure plan is invalid and require the permittee to amend the facility post-closure plan for approval by the Health District.
5. When the Health District determines that a facility post-closure plan amendment is required, the Health District shall, after consultation with the permittee, designate a compliance schedule for submittal of the amendment and its review and approval.
6. The Health District may direct facility post-closure activities to cease, in part or whole, until the post-closure plan amendment has received written approval by the Health District.
7. When post-closure activities are complete, the permittee shall certify to the Health District that the facility is stabilized and state why post-closure activities are no longer necessary. The certification shall be signed by the permittee and a professional engineer registered in the state of Washington.

- B. Minimum Functional Standards for Maintenance. The permittee shall provide all activities necessary to allow for continued facility maintenance. These activities include, but are not limited to:
1. Stormwater quantity and quality control;
 2. Slope stability, erosion and dust control;
 3. Maintenance of access roads and ditches;
 4. Maintenance of facility structures and systems (i.e., stormwater, landfill cover, gas control, surface and ground water monitoring, etc.);
 5. Control and minimization, or elimination, of threats to human health and environment; and
 6. Unauthorized entry at the facility shall be controlled, at minimum, by means of a lockable gate, barrier, fence, etc., at the property boundary.
- C. Minimum Standards for Performance. The permittee shall provide all activities necessary to allow for the continued monitoring of air, land, and water as long as is necessary for the facility to stabilize and to protect human health and environment.
1. Ground Water. The facility shall not contaminate the ground water underlying the landfill, beyond the point of compliance. The facility shall also not cause exceedances of standards contained in Chapter 173-200 WAC (Water Quality Standards for Ground Waters of the State of Washington) or Chapter 246-290 WAC (Drinking Water Regulations.)
 2. Air Quality.
 - a. The facility shall not allow explosive gases to exceed:
 - 1) Twenty-five percent of the lower explosive limit for the gases in facility structures (excluding gas control or recovery system components).
 - 2) The lower explosive limit for the gases at the property boundary or beyond.
 - 3) One hundred parts per million by volume of hydrocarbons (expressed as methane) in off-site structures.
 - b. The facility shall not cause a violation of any ambient air quality standard at the property boundary or emission standard from any emission of landfill gases, combustion, or any other emission associated with a landfill.
 - c. The permittee shall not allow non-burning flares to vent openly to the atmosphere unless it can be shown that little or no landfill gases will be produced or that landfill gases will not support combustion. In such cases installation of vents shall be required as stipulated in Chapter 173-304-460(3)(f)(ii).
 3. Surface Water. The facility shall not cause exceedances of standards contained in

Chapter 173-201A WAC (Water Quality Standards for Surface Waters of the State of Washington) or violate Chapter 90.48 RCW through the discharges of surface runoff, leachate, or any other liquid associated with a landfill.

D. Minimum Environmental Monitoring Requirements.

1. The minimum environmental monitoring requirements specified herein are subject to revision by the Health District based on the Health District's review of reports and plans required through this permit. All environmental monitoring shall be performed in accordance with the approved Sampling and Analysis Plan for the facility, as amended.

2. Ground Water Monitoring.

a. Quarterly Monitoring. Minimum quarterly monitoring required from monitoring wells MW-1, MW-3, MW-5A, MW-6, and MW-7. Each individual monitoring well shall be analyzed for:

| | |
|---|---------------------------------|
| Water Level | Temperature |
| Specific Conductance | pH |
| Total Coliform | Sulfate |
| Potassium | Bicarbonate |
| Carbonate | Sodium |
| Chloride | Calcium |
| Total Organic Carbon | Nitrite, Nitrate, & Ammonia (N) |
| Dissolved Arsenic | Chemical Oxygen Demand |
| Dissolved Iron | Dissolved Barium |
| Dissolved Zinc | Dissolved Manganese |
| Volatile Organic Compounds (EPA Method 8260B) | |

b. Ground water analyses shall be performed by a laboratory accredited by the Washington State Department of Ecology. The methods of analysis shall be in accordance with EPA Publication Number SW-846, "Test Methods for Evaluating Solid Waste - Physical/Chemical Methods". For each test method, the lowest achievable detection limit shall be obtained sufficient to verify compliance with the regulatory standards in Section IV.C.1. of this permit.

3. Surface Water Monitoring.

a. Annual Monitoring. Minimum annual monitoring required from surface water station SW-2. The annual event shall occur either between January and March or November and December, and shall be analyzed for:

| | | |
|-------------|----|------------------|
| Temperature | pH | Nitrate-Nitrogen |
|-------------|----|------------------|

Specific Conductance Fecal Coliform

- b. Surface water analysis shall be performed by a laboratory accredited by the Washington State Department of Ecology. The methods of analysis shall be in accordance with 40 CFR Part 136. For each test method, the lowest achievable detection limit shall be obtained sufficient to verify compliance with the standards in Section IV.C.3. of this permit.
4. Landfill Gas Monitoring. Minimum quarterly monitoring of landfill gas is required from gas flares GF-1, GF-2, and GF-3. Individual samples shall be monitored for the following:

| | |
|----------------|-----------------------|
| Methane (%LEL) | Gas Pressure |
| Oxygen (%VOL) | Carbon Dioxide (%VOL) |

E. Reporting Requirements. All reports specified herein must be forwarded to the Health District representative responsible for this permit, and to the Solid Waste Supervisor, Department of Ecology Northwest Regional Office, Solid Waste & Financial Assistance Program, 3190 160th Avenue S.E., Bellevue, WA 98008-5452.

1. Annual Report.
 - a. Annual report due April 1, 2011 (for 2010 calendar year), and April 1, 2012 (for 2011 calendar year).
 - b. The permittee shall prepare and submit one copy and one CD of the annual report to the Health District, and one copy and one CD to Ecology. The annual report will serve as the permit application for the next year's permit. The annual report shall cover facility activities during the previous calendar year and must include, at a minimum, the following information:
 - 1) Name and address of the facility.
 - 2) Calendar year covered by the report.
 - 3) A summary of all ground water, surface water, and landfill gas monitoring will be provided in accordance with the Revised Statistical Approach for Olalla Landfill (June 2008). The summary shall contain the results of applicable statistical analyses, explanations for increasing/decreasing contaminant level trends, confidence intervals and if outliers were discarded, a justification for this action. In addition time series plots for each assessment monitoring well for each parameter that exceeded groundwater quality standards at any time since well installation. All statistically significant increasing/decreasing contaminant level trends and exceedances of ground

water, surface water, or air quality standards as described in Section IV shall be reported and explained to the extent possible, and if deemed necessary, a summary of actions taken to verify results. Additionally, if the trends shows either no improvement or increasing contaminate levels or exceedances are detected at or beyond the point of compliance, a plan of action shall be included which describes what activities will be taken to correct the statistically significant elevated contaminants or exceedances, and the expected time frame required to do so. The summary shall also contain a description of the ground water flow rate and direction in the uppermost aquifer. Any changes in the number or location of monitoring systems shall also be noted and explained.

- 4) A description of the next year's proposed environmental monitoring program, operations, and maintenance activities.
- 5) A summary of all activities related to the inspection, maintenance, and engineering of the facility and its systems. All field data sheets, and inspection and log forms shall be made available to the Health District upon request.
- 6) An explanation of causes and effects of all instances of non-compliance with the provisions of this permit and a summary of how these issues were or are being corrected.

2. Quarterly Environmental Monitoring Reports.

- a. The permittee shall prepare and submit written quarterly environmental monitoring reports to report the information gathered under the "Minimum Environmental Monitoring Requirements" section of this permit. Gas, surface and ground water monitoring data shall be summarized, evaluated, and compared with all applicable standards, and shall also be submitted in hard copy form. All monitoring data shall be reported in detail on a quarterly basis. Results and summary of the fourth quarter sampling event shall be combined with the annual report.
- b. Environmental monitoring reports shall be submitted to the Health District and Ecology within sixty (60) days of the date of sample collection. If the report cannot be submitted within sixty (60) days, the Health District will be notified prior to the end of sixty (60) days as to the status of the report and when it will be submitted. In no event shall monitoring reports be submitted more than seventy (70) days after the date of sample collection. All instances of regulatory standard exceedances shall be faxed to the Health District within ten (10) days of the time of discovery.

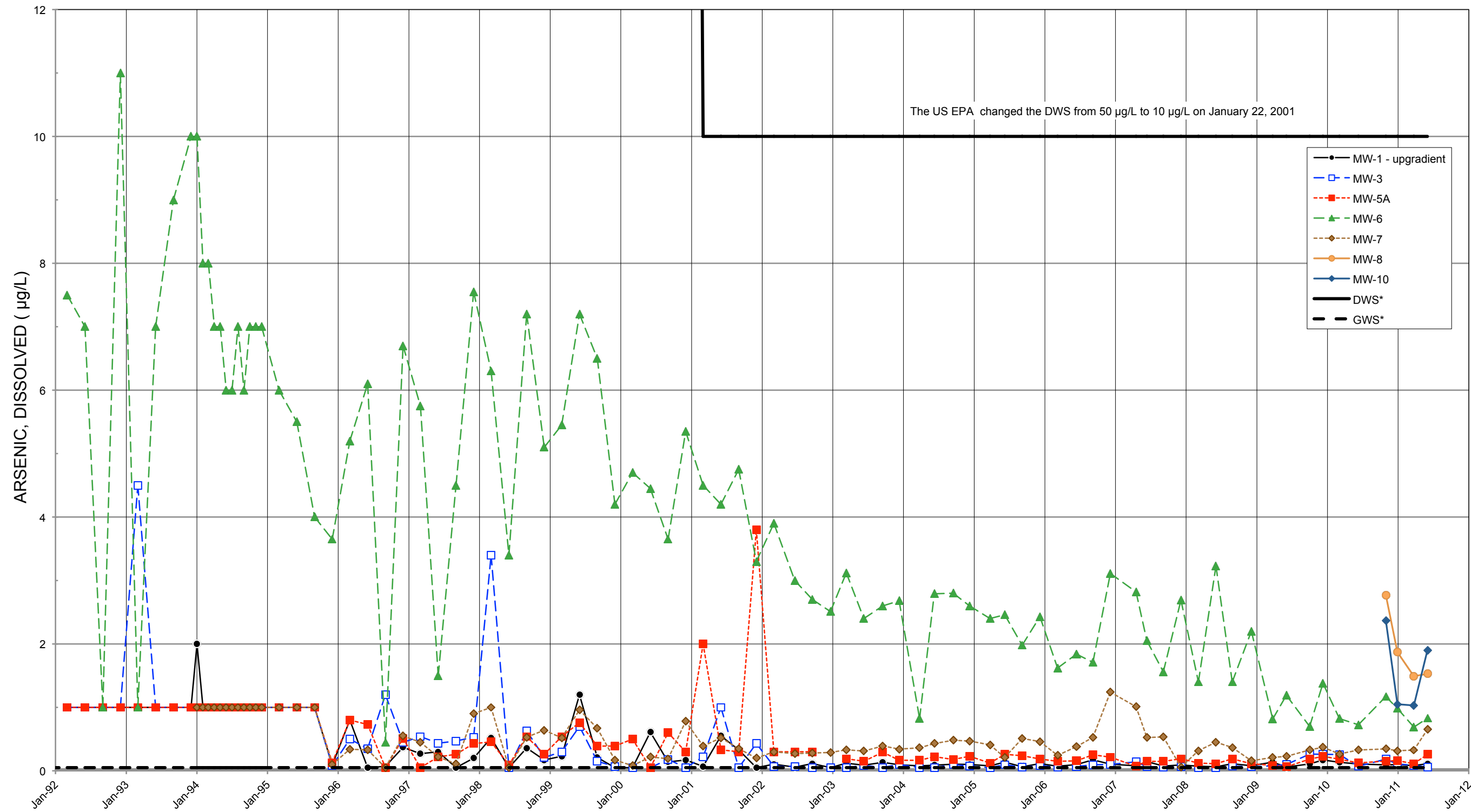
SECTION V. CORRECTIVE ACTION PLAN

| TASKS | Submittal Date |
|---|-----------------------|
| 1. Prepare and submit a work plan for a RI/FS under MTCA which will address water quality violations at the point of compliance under WAC 173-304. | 6/30/2010 |
| 2. MTCA PROCESS SCHEDULE: Prepare and submit a schedule outlining activities to be completed in the RI/FS and anticipated completion dates. Upon approval of schedule by Health District and Department of Ecology, project milestones will be incorporated into this permit. | 6/30/2010 |

Appendix C

Constituents of Concern – Full and Recent (5-Year)
Time Series Graphs

OLALLA LANDFILL Quarterly Monitoring Data

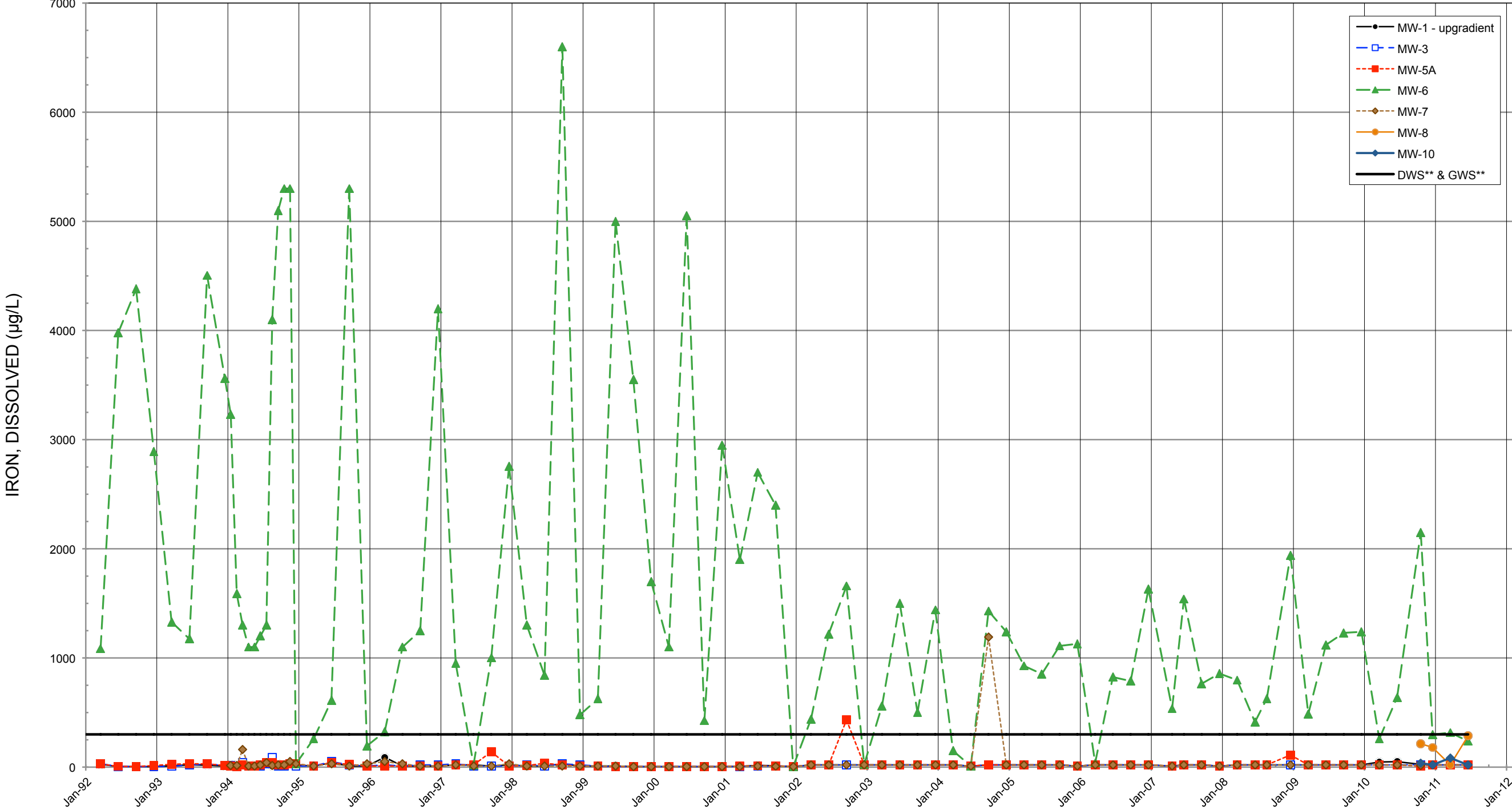


Primary Drinking Water Standard (DWS) = 10 µg/L
 Primary Groundwater Standard (GWS) = 0.05 µg/L

DATE

DISSOLVED ARSENIC
(FULL)

OLALLA LANDFILL Quarterly Monitoring Data

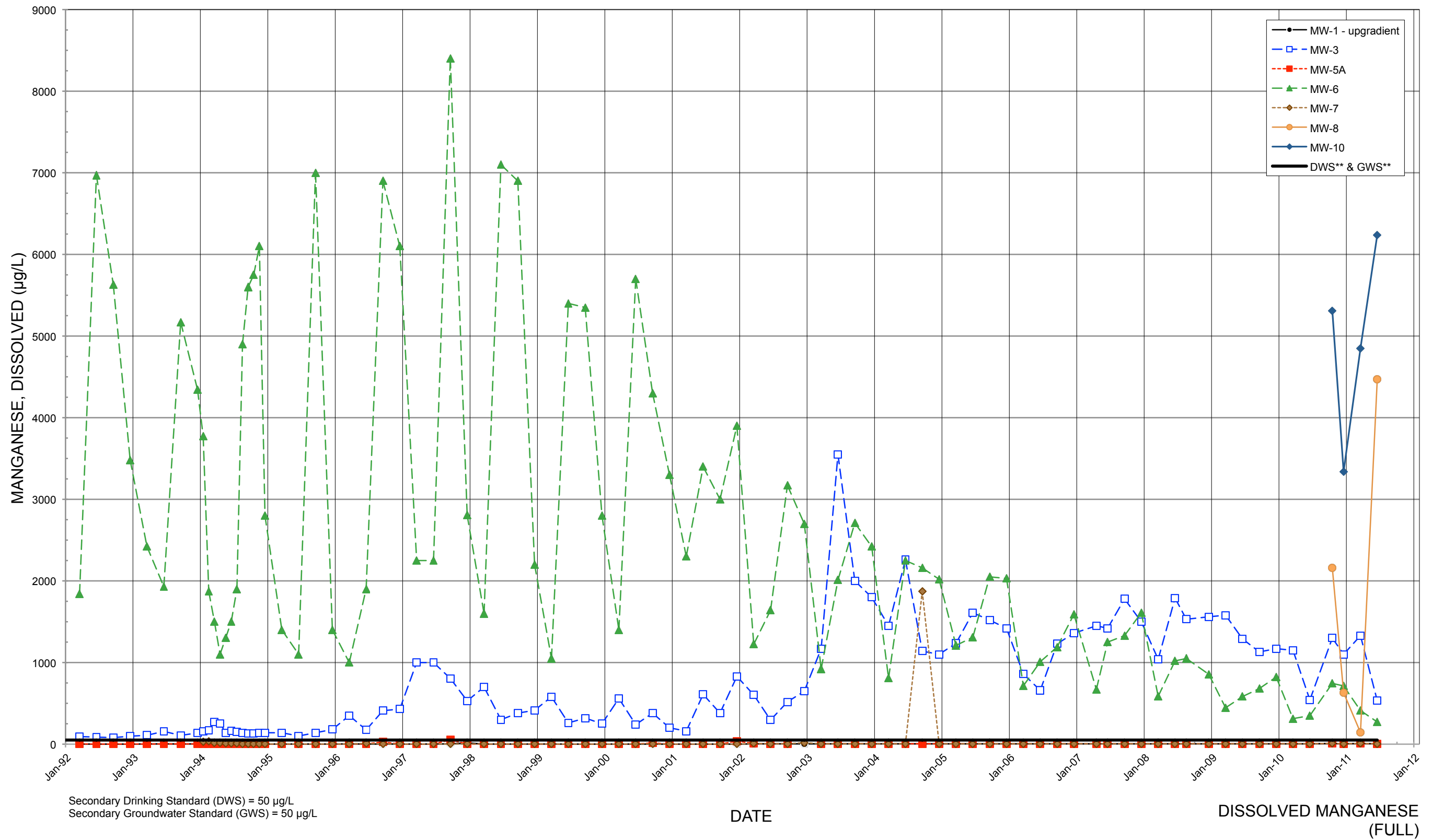


Secondary Drinking Water Standard (DWS) = 300 µg/L
 Secondary Groundwater Standard (GWS) = 300 µg/L

DATE

DISSOLVED IRON
(FULL)

OLALLA LANDFILL Quarterly Monitoring Data

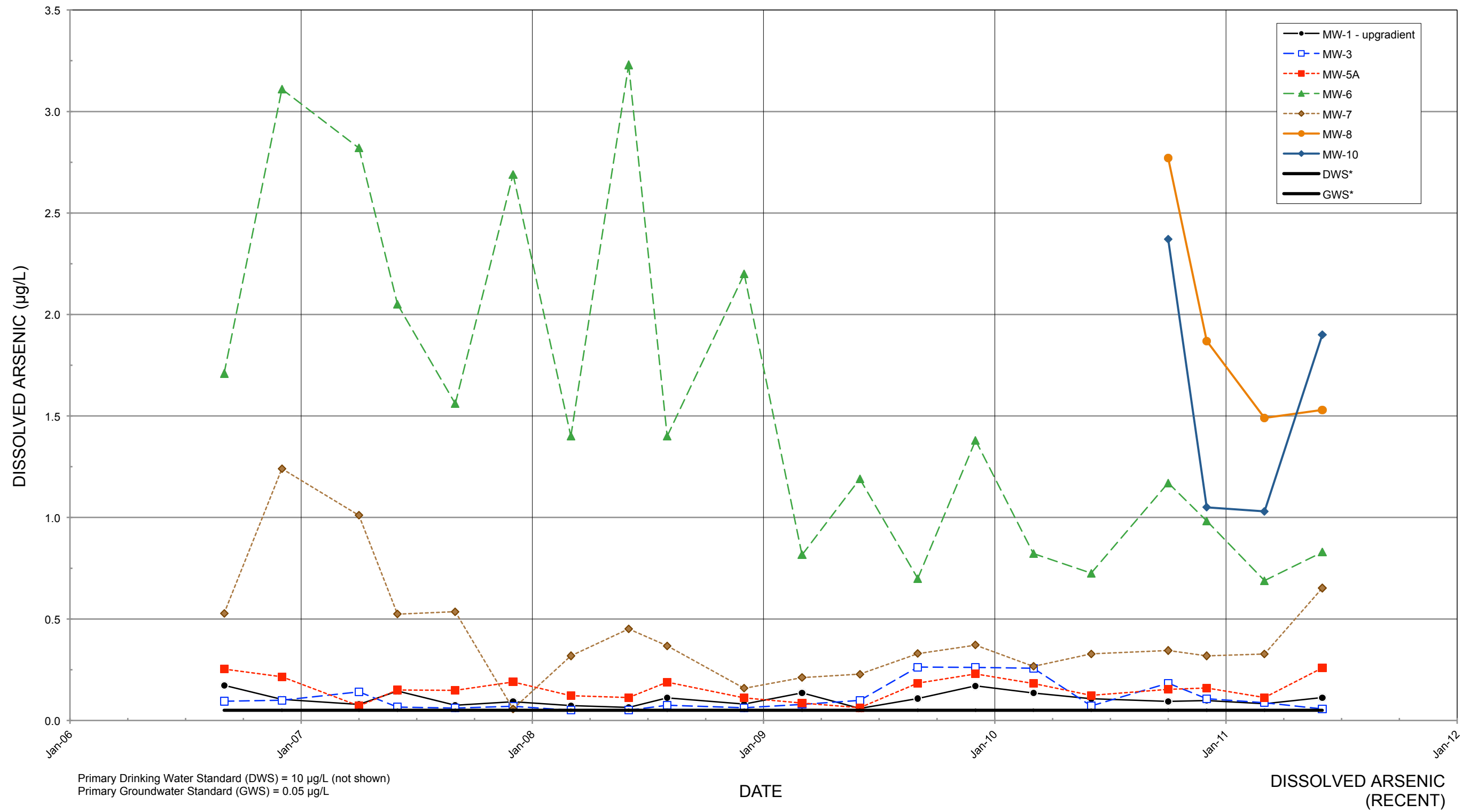


Secondary Drinking Standard (DWS) = 50 µg/L
 Secondary Groundwater Standard (GWS) = 50 µg/L

DISSOLVED MANGANESE
(FULL)

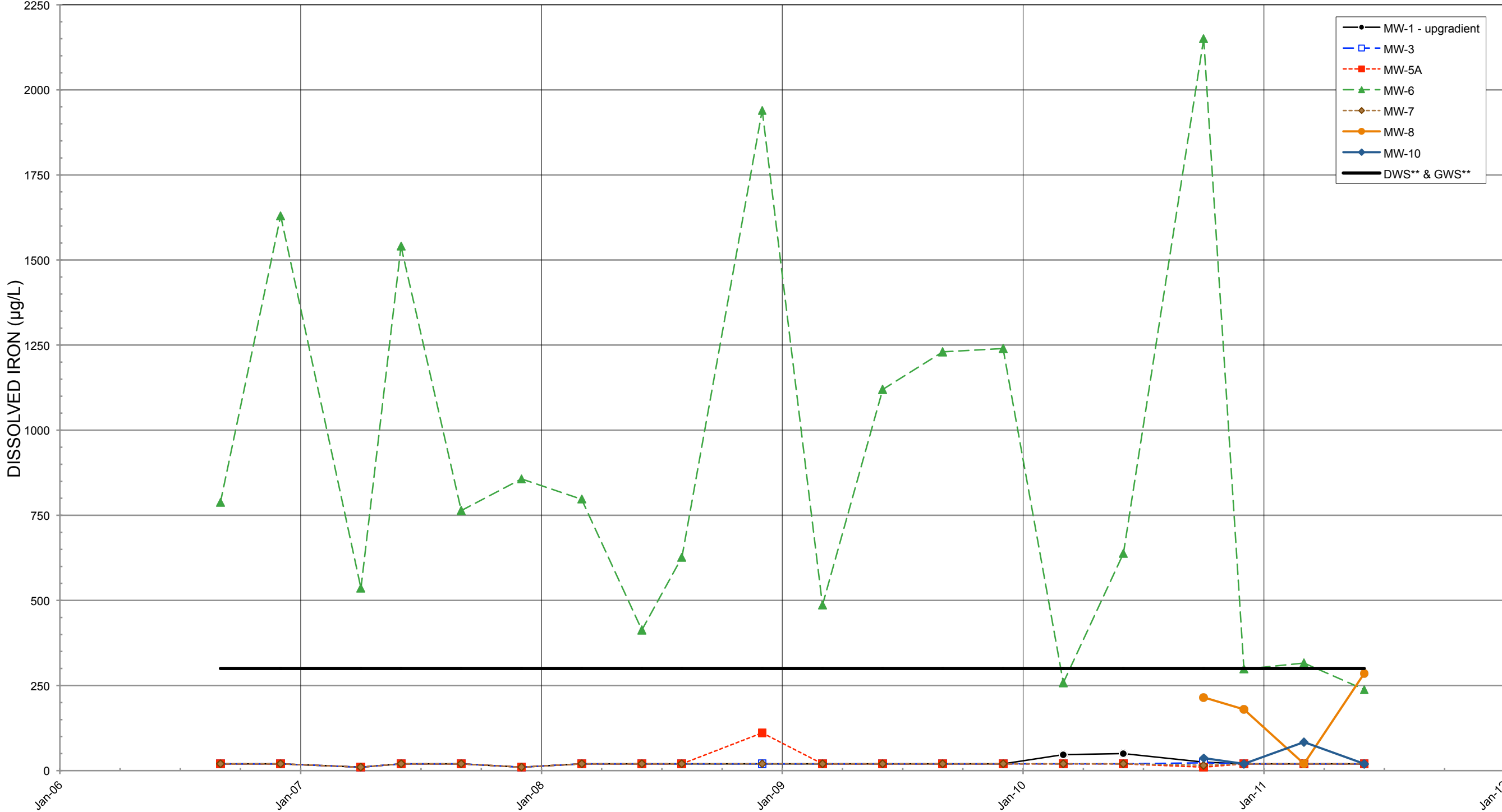
OLALLA LANDFILL

Quarterly Monitoring Data (recent five years)



OLALLA LANDFILL

Quarterly Monitoring Data (recent five years)



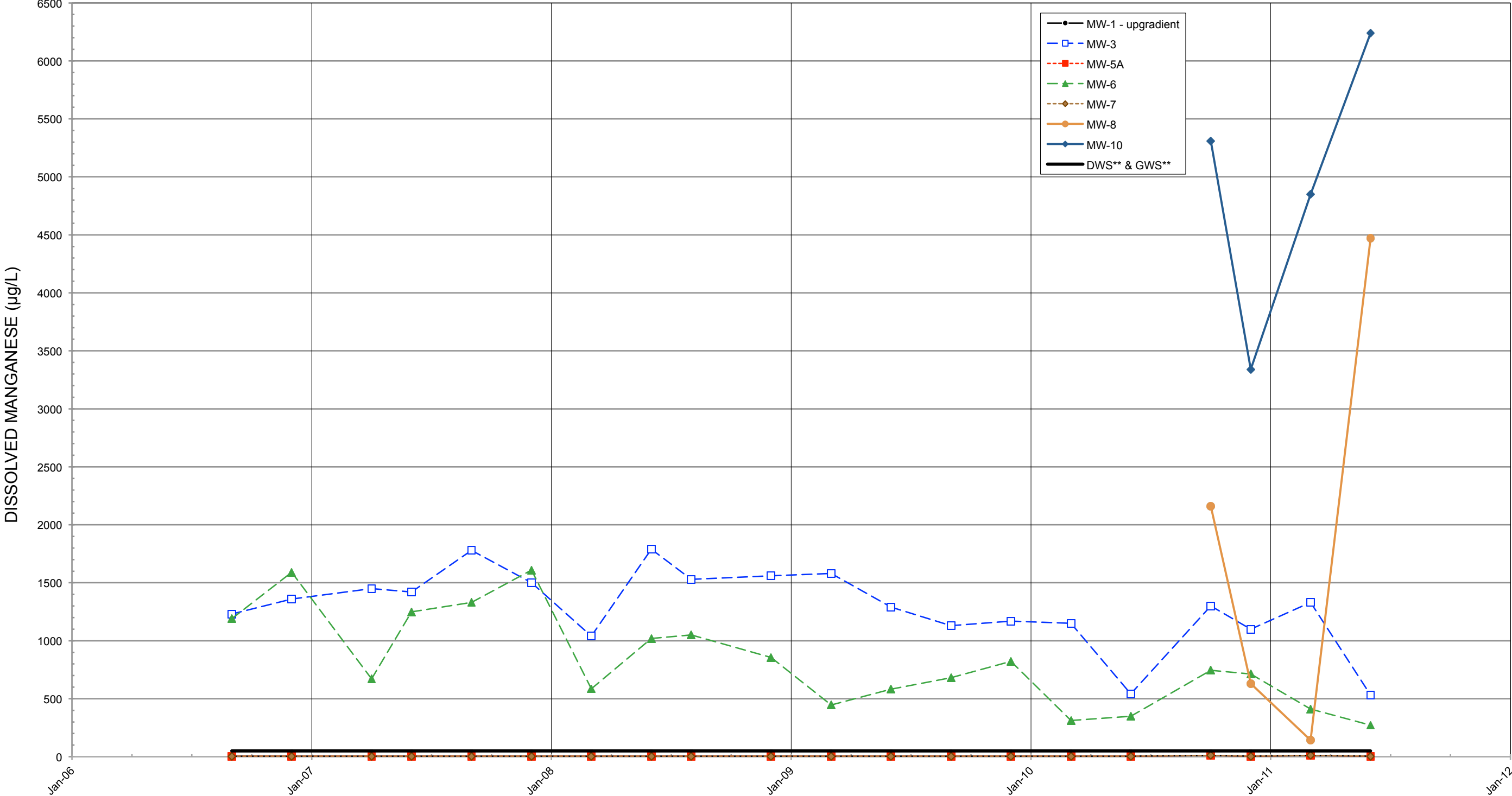
Secondary Drinking Water Standard (DWS) = 300 µg/L
 Secondary Groundwater Standard (GWS) = 300 µg/L

DATE

DISSOLVED IRON (RECENT)

OLALLA LANDFILL

Quarterly Monitoring Data (recent five years)



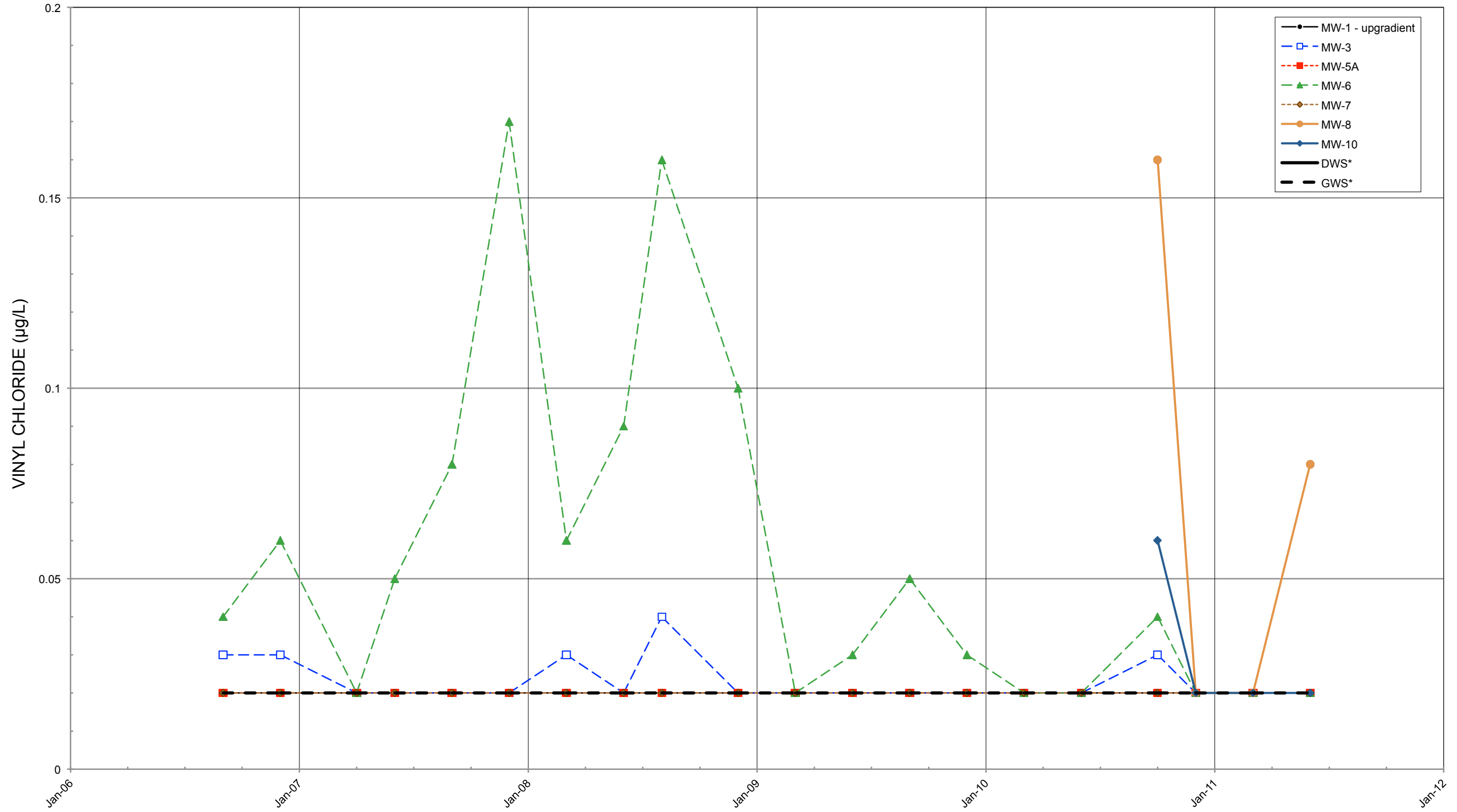
Secondary Drinking Standard (DWS) = 50 µg/L
 Secondary Groundwater Standard (GWS) = 50 µg/L

DATE

DISSOLVED MANGANESE
(RECENT)

OLALLA LANDFILL

Quarterly Monitoring Data (recent five years)



Primary Drinking Water Standard (DWS) = 2 µg/L (not shown)
 Primary Groundwater Standard (GWS) = 0.02 µg/L

DATE

VINYL CHLORIDE
(RECENT)

Appendix D

Geologic Logs and As-Built Diagrams for
MW-8 and MW-10, Split-Spoon Sampler
Photos, Well Development Field Data Sheets,
and Sieve Analysis Curves



MW-8: Sample from 20-21.5 ft. bgs – Poorly-Graded Sand (SP) Gray-Brown, moist to wet, mostly medium sand with trace gravel and trace silt. Photo by EPI.



MW-8: Sample from 20-21.5 bgs – Poorly-Graded Sand (SP) Gray-Brown, moist to wet, mostly medium sand with trace gravel and trace silt. Photo by EPI.



MW-8: Sample from 25-26.5 ft. bgs – Poorly Graded Sand (SP) Gray-Brown, moist to wet, mostly medium sand with trace gravel and trace silt. Photo by EPI.



MW-8: Sample from 37-38.5 ft. bgs – Poorly-Graded Sand (SP) Wet, Grades to well-rounded medium sand with black, gray, and white grains. Photo by EPI.



MW-10: Sample from 45-46.5 ft. bgs – Poorly Graded Sand (SP) Gray-Brown, Wet, mostly medium sand with trace gravel and trace silt. Photo by EPI.



MW-10: Sample from 45-46.5 ft. bgs – Poorly Graded Sand (SP) Gray-Brown, Wet, mostly medium sand with trace gravel and trace silt. Photo by EPI.

Monitoring Well Development Field Measurements Record

Job Name/Location: Olalla Landfill RI/FS
 Job Number: 60101.0 Task 3
 Date: 13-Oct-10

Well #: MW-2
 Page Number: 1 of 1

| Time | Depth to Water (ft.) | Total Depth of Well (ft.) | Volume Purged (gal.) | pH | Cond. (µmhos/cm) | Turbidity (NTU) | Temp. (oC) | Visual Observations |
|-------|----------------------|---------------------------|----------------------|------|------------------|-----------------|------------|--|
| 10:25 | 66.64 | 78.7 | start | | | | | Geliton-like material on bailer, abundant fine sand. |
| 12:00 | 66.65 | | 10 | 7.85 | 85 | 225 | 12.6 | cloudy, some fine sand |
| 12.35 | 66.64 | 78.7 | 15 | 7.77 | 89 | 191 | 12.2 | cloudy, some fine sand, bailing stopped |
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Note: Development started by hand-bailing with stainless steel bailer, then switched to a Waterra inertial pump at 11:00, pump clogged by fine sand, switched back to bailer.



Monitoring Well Development Field Measurements Record

Job Name/Location: Olalla Landfill RI/FS
 Job Number: 60101.0 Task 3
 Date: 13-Oct-10

Well #: MW-4
 Page Numb 1 of 1

| Time | Depth to Water (ft.) | Total Depth of Well (ft.) | Volume Purged (gal.) | pH | Cond. (µmhos/cm) | Turbidity (NTU) | Temp. (oC) | Visual Observations |
|-------|----------------------|---------------------------|----------------------|------|------------------|-----------------|------------|--|
| 8:24 | 63.98 | 73.9 | start | | | | | Geliton-like material on bailer, abundant fine sand. |
| 9:17 | | | 5 | | | | | Cloudy, some silt and sand |
| 9:50 | | | 8 | 7.41 | 62 | 229 | 12.1 | Cloudy, some sand |
| 10:00 | | | 10 | 7.31 | 58 | 35 | 12.2 | Cloudy, some sand |
| 10:10 | 63.98 | | 15 | 7.31 | 57 | 205 | 12.2 | Cloudy, trace sand, stop pumping |
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Note: Development started by hand-bailing with stainless steel bailer until clarity improved, then switched to a Waterra inertial pump.



Monitoring Well Development Field Measurements Record

Job Name/Location: Olalla Landfill RI/FS
 Job Number: 60101.0 Task 3
 Date: 8-Oct-10

Well #: MW-8
 Page Number: 1 of 1

| Time | Depth to Water (ft.) | Total Depth of Well (ft.) | Volume Purged (gal.) | pH | Cond. (µmhos/cm) | Turbidity (NTU) | Temp. (oC) | Visual Observations |
|------|----------------------|---------------------------|----------------------|------|------------------|-----------------|------------|----------------------------------|
| 8:00 | 22.81 | 37.85 | start pump | | | | | |
| 8:08 | | | 10 | | | | | Cloudy with fines |
| 8:17 | | | 15 | | | | | same |
| 8:45 | | | 55 | | | | | same |
| 9:05 | | | 75 | 6.67 | 115 | 39.4 | 10.8 | Cloudy, some fines after surging |
| 9:25 | 22.84 | | 85 | 6.5 | 134 | 19.6 | 10.2 | Clear, stop pumping |
| | | | | | | | | |
| | | | | | | | | |
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| | | | | | | | | |
| | | | | | | | | |

Notes: Development performed using submersible pump



Monitoring Well Development Field Measurements Record

Job Name/Location: Olalla Landfill RI/FS
 Job Number: 60101.0 Task 3
 Date: 8-Oct-10

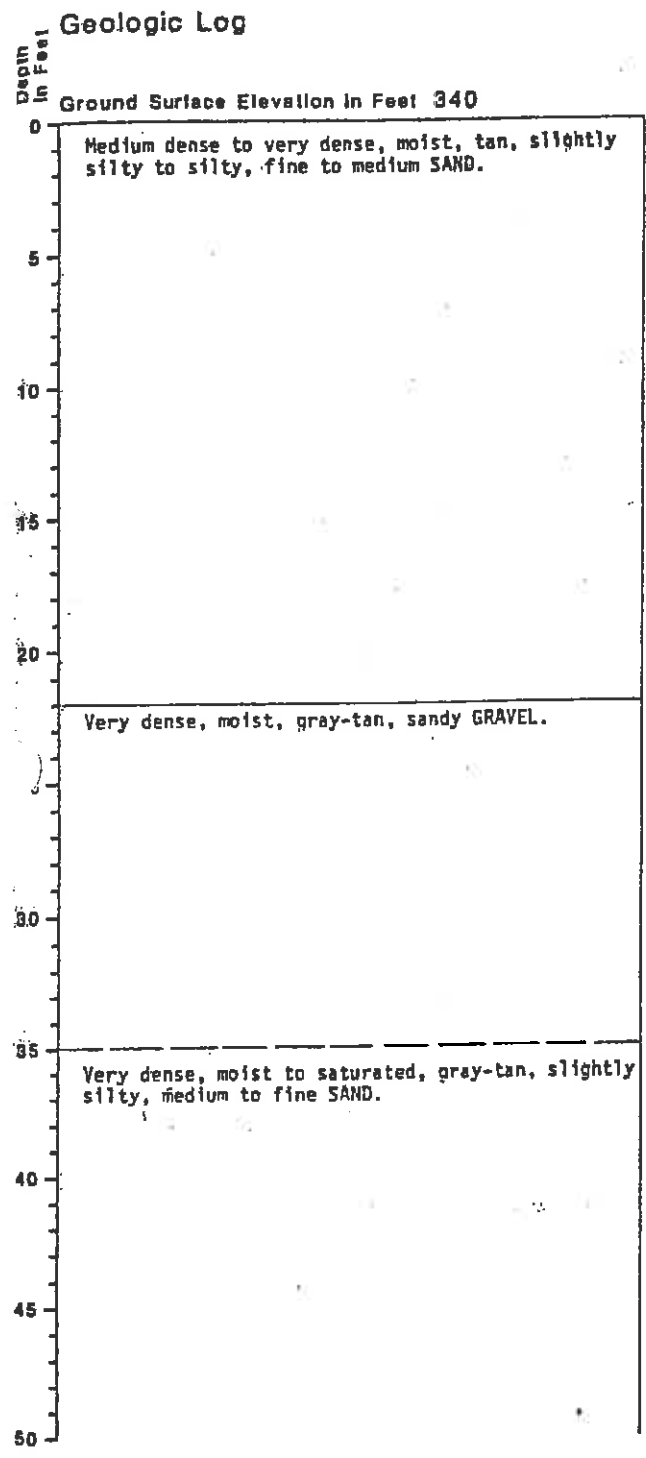
Well #: MW-10

Page Number: _____

| Time | Depth to Water (ft.) | Total Depth of Well (ft.) | Volume Purged (gal.) | pH | Cond. (µmhos/cm) | Turbidity (NTU) | Temp. (oC) | Visual Observations |
|------|----------------------|---------------------------|----------------------|------|------------------|-----------------|------------|----------------------|
| 8:12 | 32.10 | 49.92 | | | | | | start pump |
| 8:22 | | | 15 | | | | | Cloudy |
| 8:38 | | | 22 | | | | | Cloudy after surging |
| 8:55 | | | 35 | 7.82 | 276 | 1.76 | | clear |
| 9:06 | 32.10 | | 50 | 7.62 | 285 | 1.91 | 12.2 | clear, stop pumping |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
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| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |

Notes: Development performed using submersible pump

Boring Log and Construction Data for Well MW-1



Well Design 3
 Top Casing Elevation in Feet 341.41
 Casing Stickup in Feet 1

| Lab | Sample | N | pH | EC |
|------|--------|-------|-----|----|
| S-1 | X | 23 | 6.5 | <5 |
| S-2 | X | 31 | 5.5 | <5 |
| S-3 | X | 50/4" | 5.2 | 5 |
| S-4 | X | 50/6" | 5.3 | 6 |
| S-5 | X | 50/6" | 4.9 | 6 |
| S-6 | X | 50/4" | 5.0 | 19 |
| S-7 | X | 50/4" | 5.2 | 5 |
| S-8 | X | 50/6" | 5.3 | <5 |
| S-9 | X | 84/9" | 5.1 | <5 |
| S-10 | X | 50/5" | 5.1 | <5 |

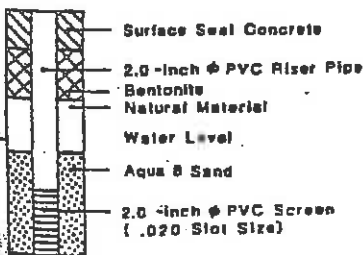
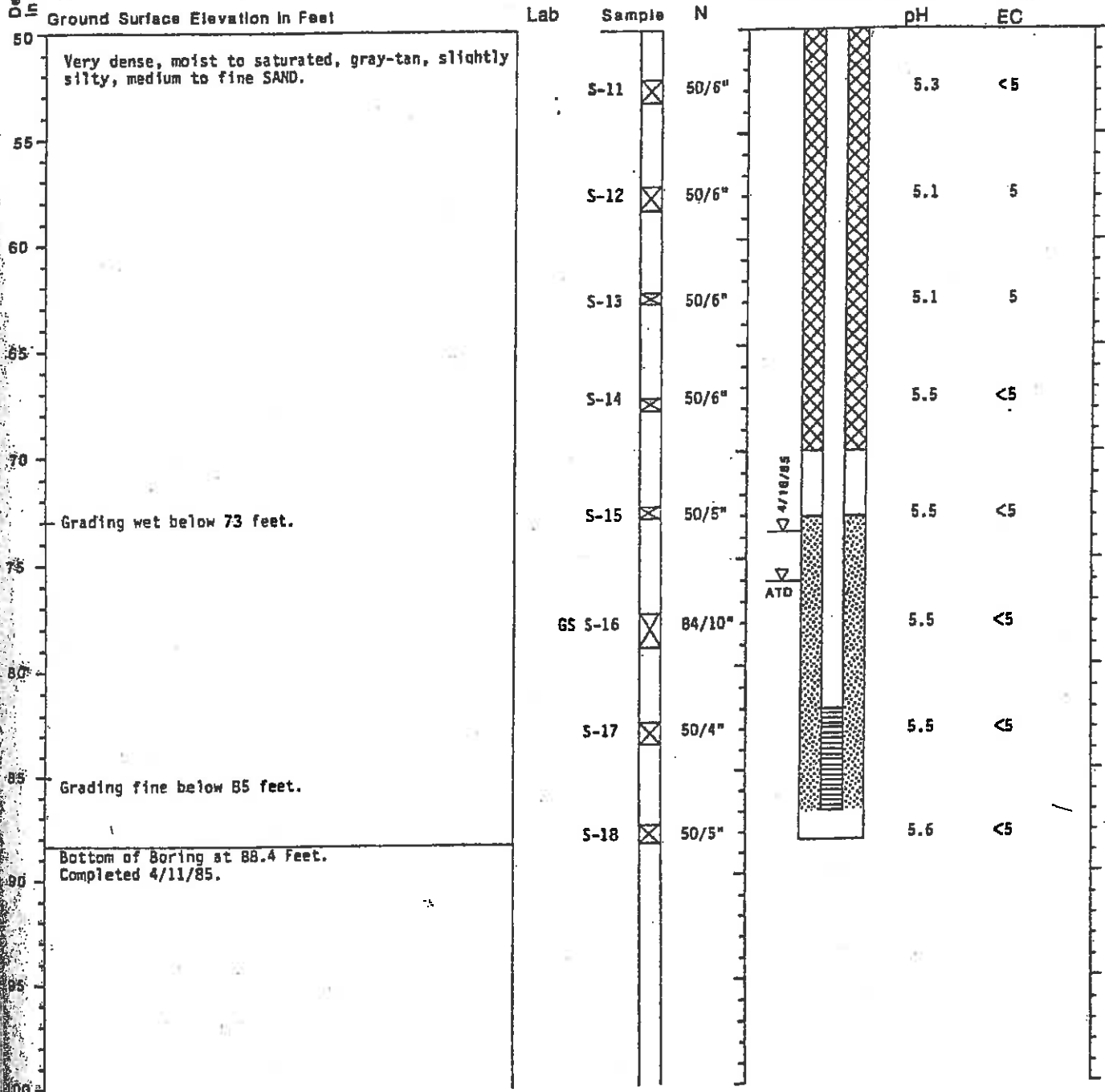
Boring Log and Construction Data for Well MW-1

Geologic Log

Well Design 3

Top Casing Elevation in Feet 341.41

Casing Slickup in Feet 1

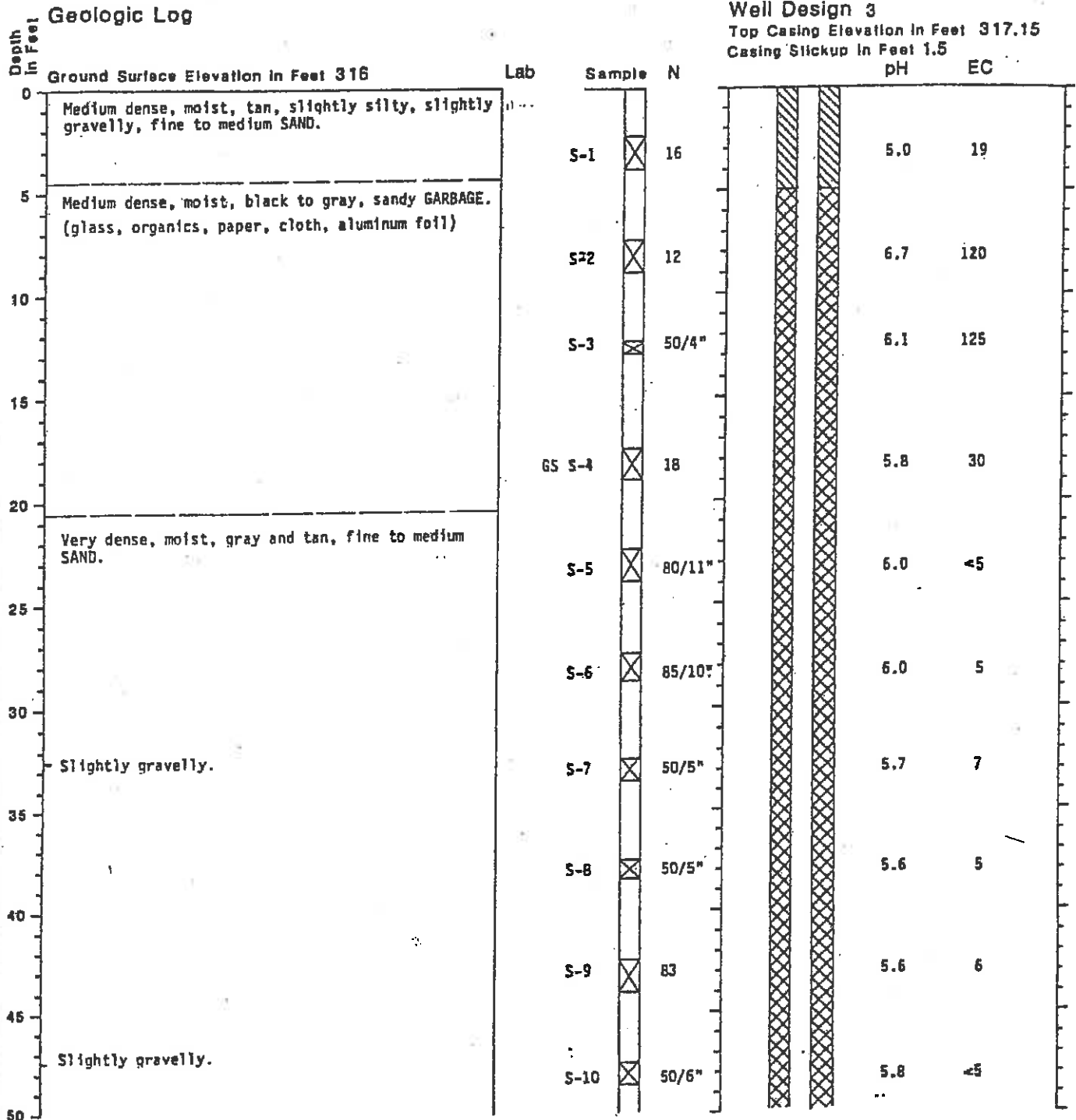


- 2-inch O.D. Split Spoon Sample
- No Sample Recovery
- Standard Penetration Resistance, Blows per foot
- Grain Size Analysis
- Permeability Test

NOTES:

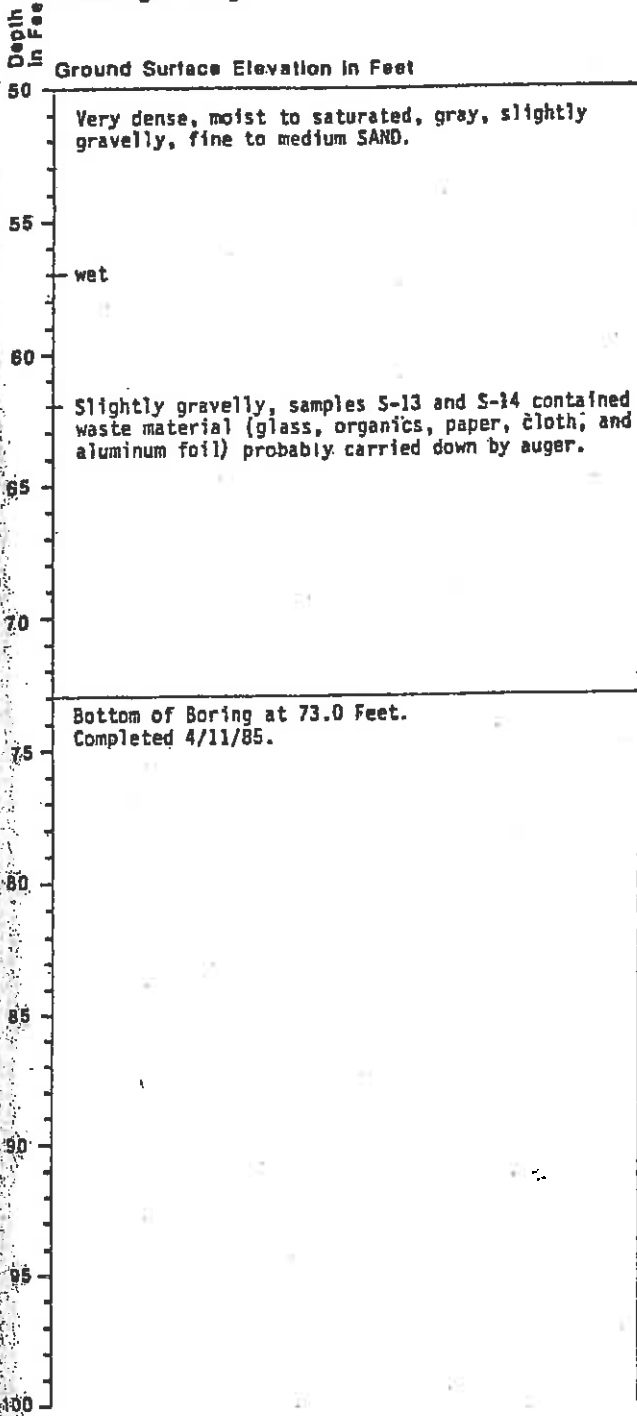
1. Soil descriptions are interpretive and actual changes may be gradual.
2. Water Level is for date indicated and may vary with time of year. ATD: At Time of Drilling
3. Top of PVC casing surveyed by Parametrix to estimate mean sea level datum.

Boring Log and Construction Data for Well MW-2



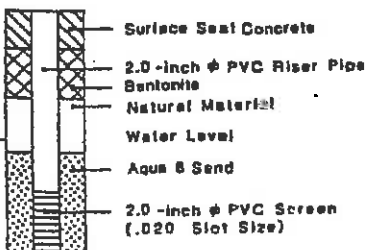
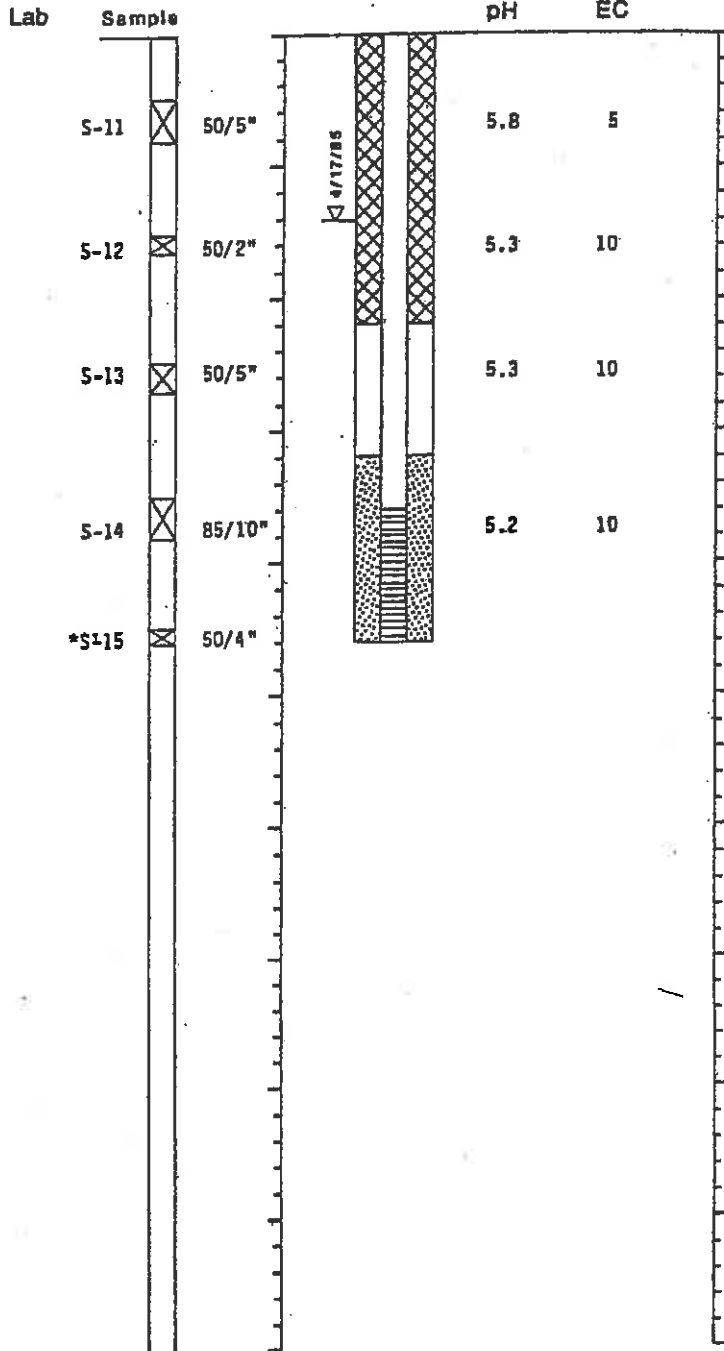
Boring Log and Construction Data for Well MW-2

Geologic Log



Well Design

Top Casing Elevation in Feet
Casing Stickup in Feet

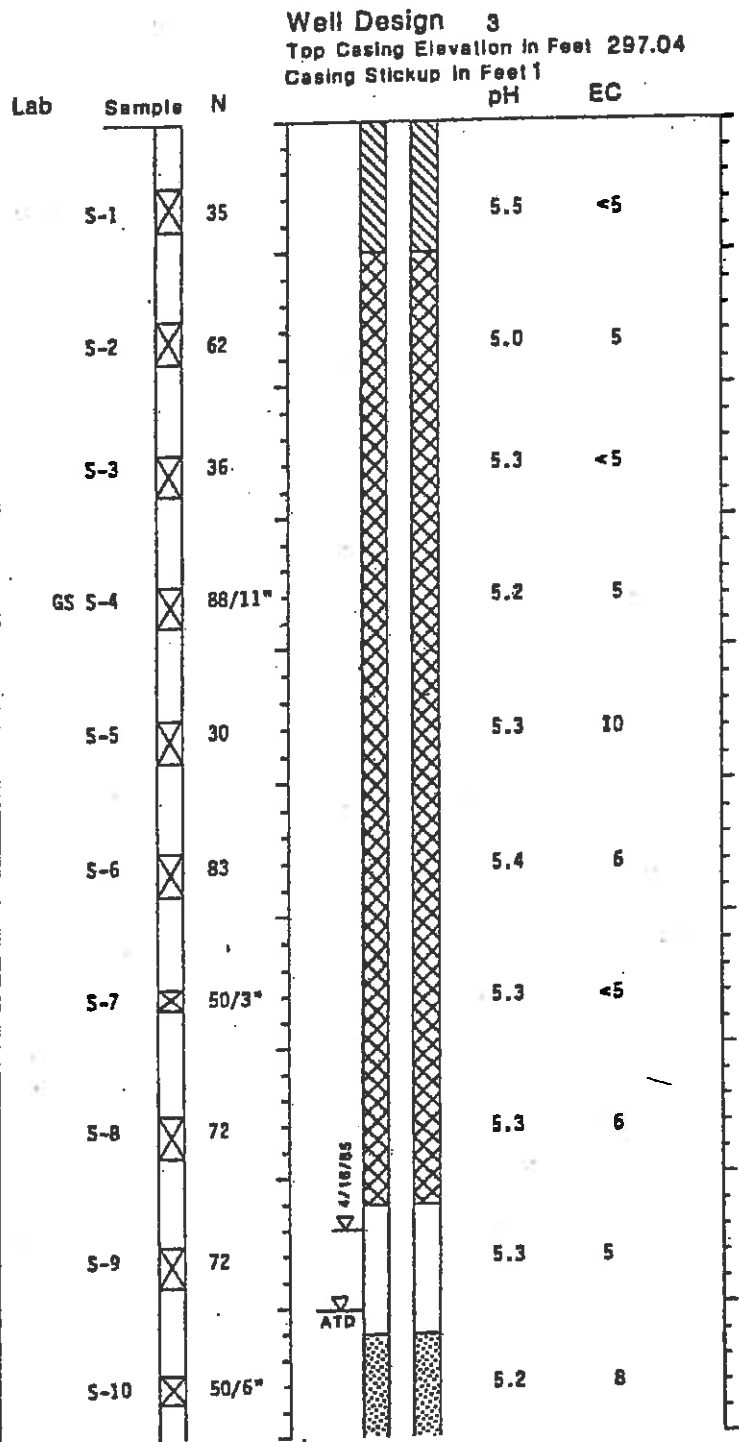
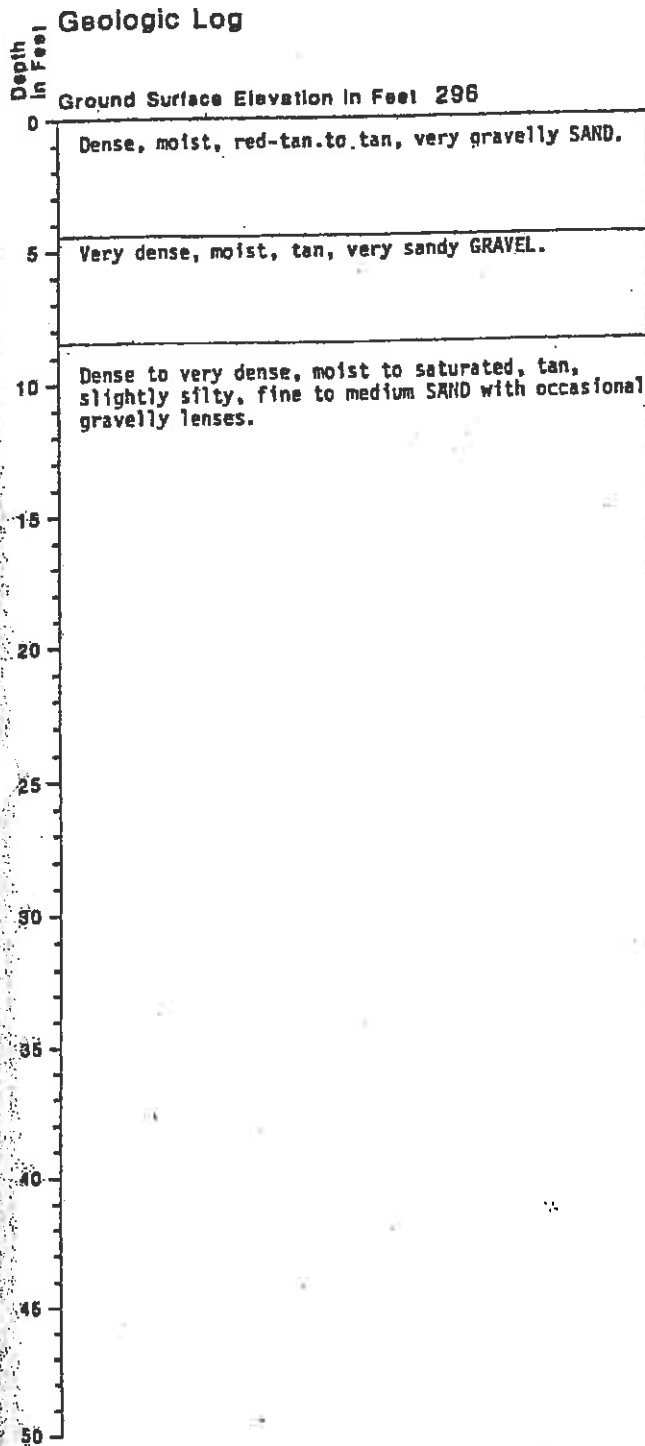


- 2-inch O.D. Split Spoon Sample
- No Sample Recovery
- Standard Penetration Resistance, Blows per foot
- Grain Size Analysis
- Permeability Test

NOTES:

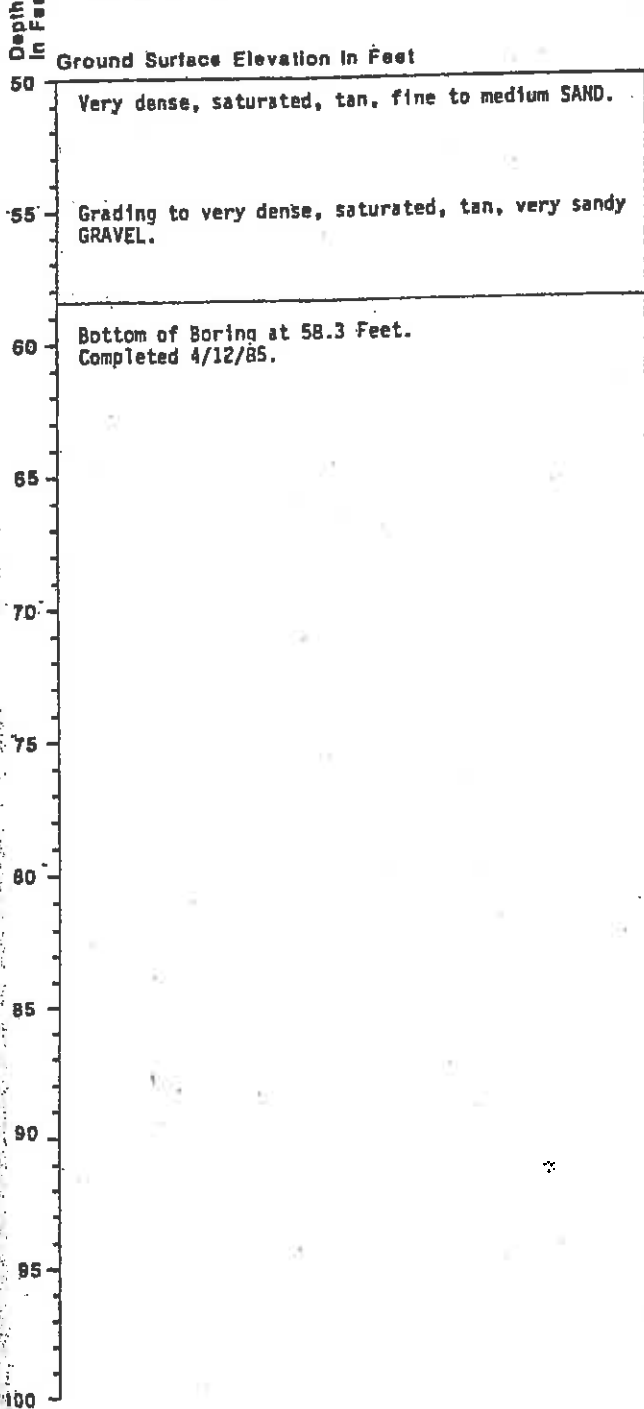
- Soil descriptions are interpretive and actual changes may be gradual.
- Water Level is for date indicated and may vary with time of year. ATD: At Time of Drilling.
- Top of PVC casing surveyed by Parametrix to estimate mean sea level datum.

Boring Log and Construction Data for Well MW-3



Boring Log and Construction Data for Well MW-3

Geologic Log



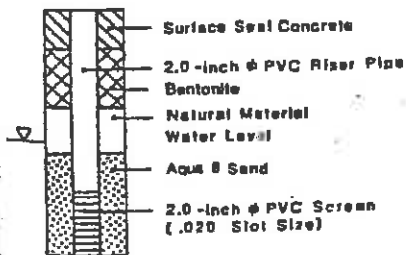
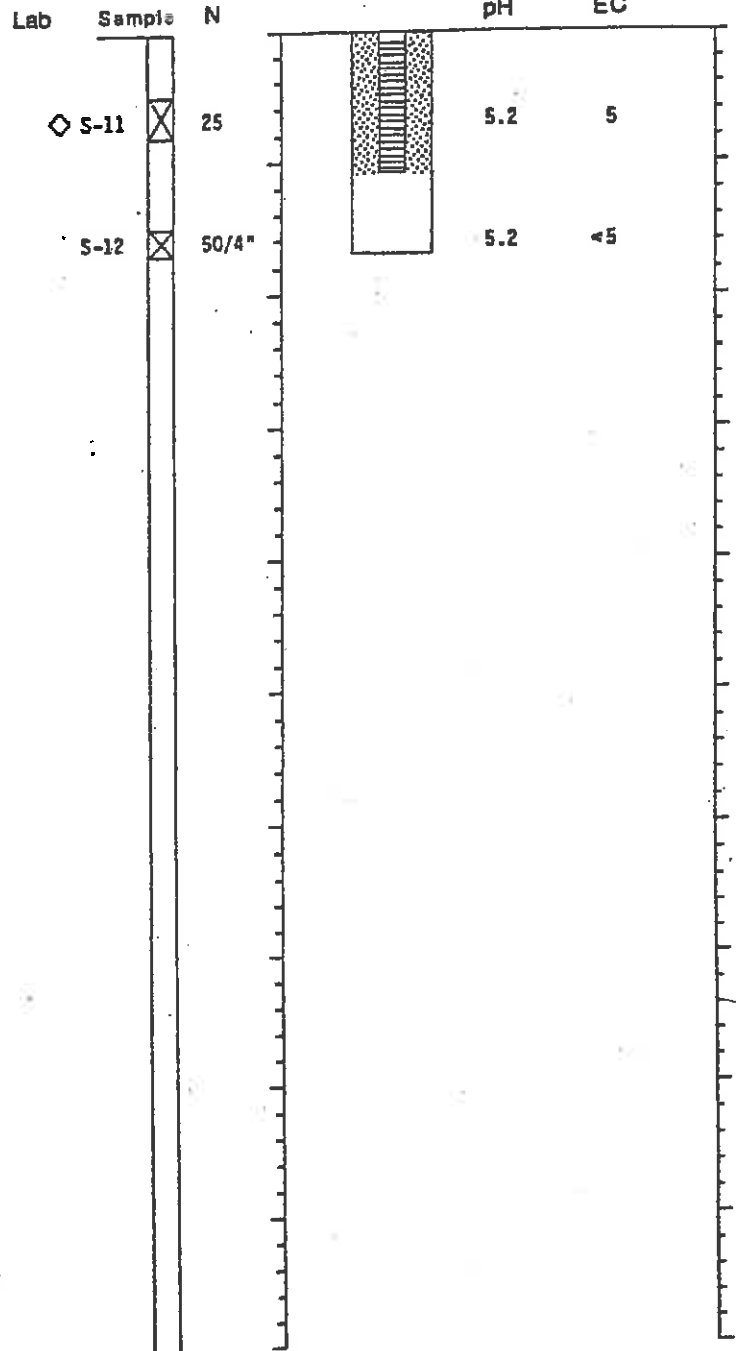
Well Design

Top Casing Elevation In Feet

Casing Stickup In Feet

pH

EC



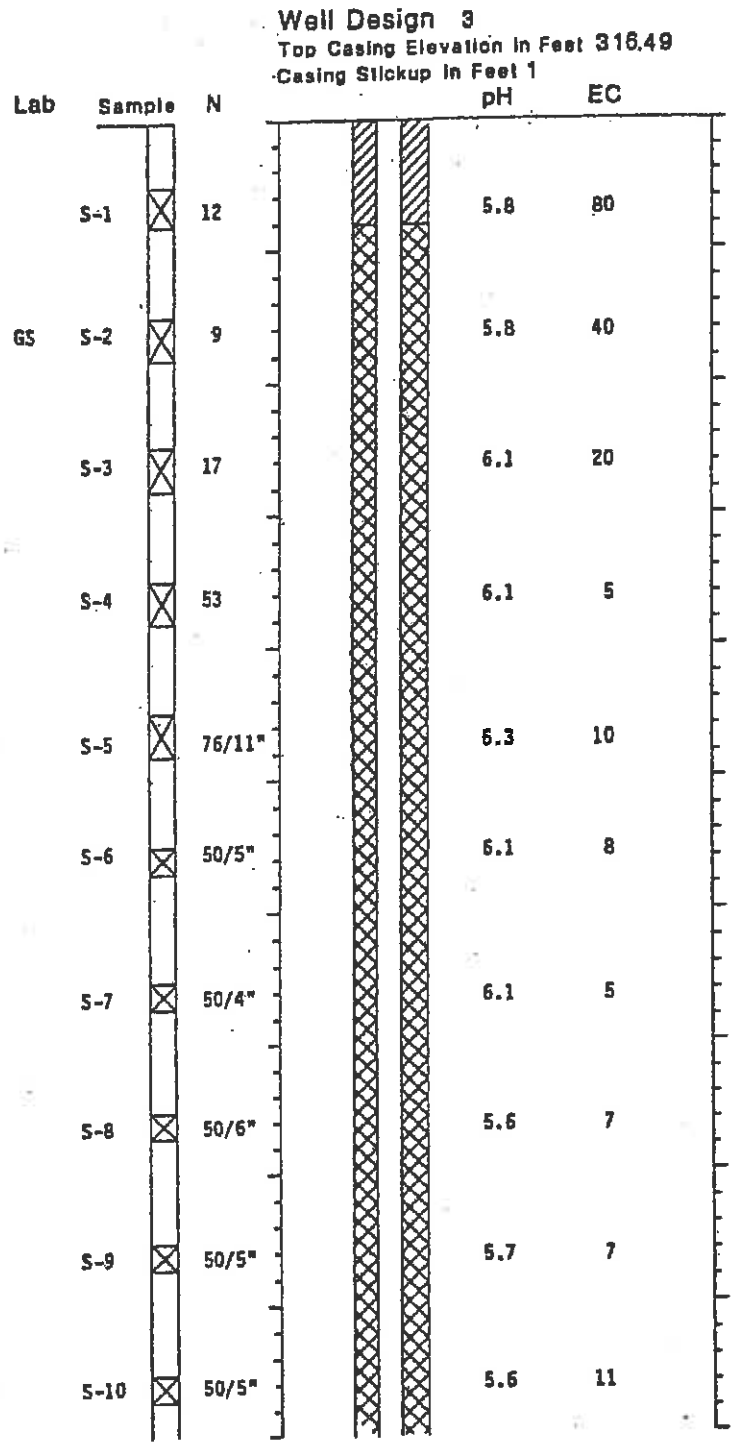
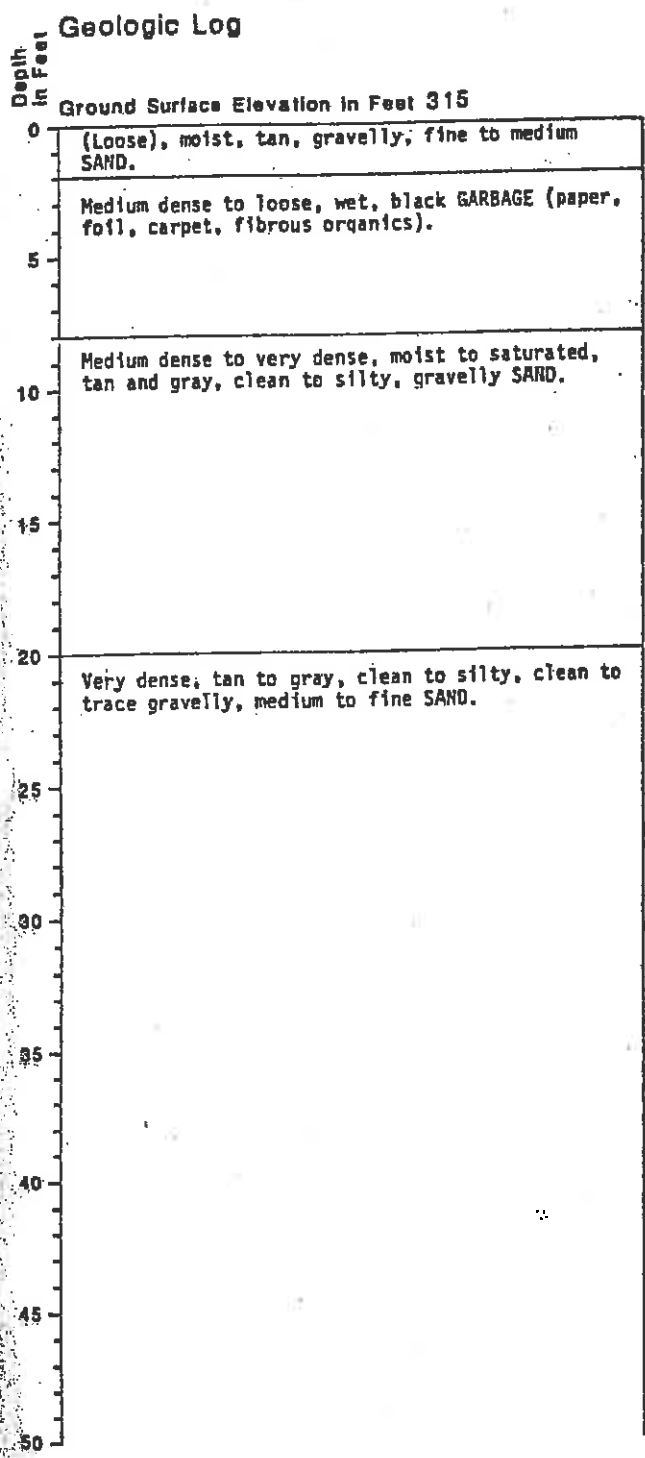
- 2-inch O.D. Split Spoon Sample
- * No Sample Recovery
- N Standard Penetration Resistance, Blows per foot
- GS Grain Size Analysis
- K Permeability Test
- Low blow count due to disturbed material.

NOTES:

- Soil descriptions are interpretive and actual changes may be gradual.
- Water Level is for date indicated and may vary with time of year. ATD:At Time of Drilling
- Top of PVC casing surveyed by Parametrix to estimate mean sea level datum.

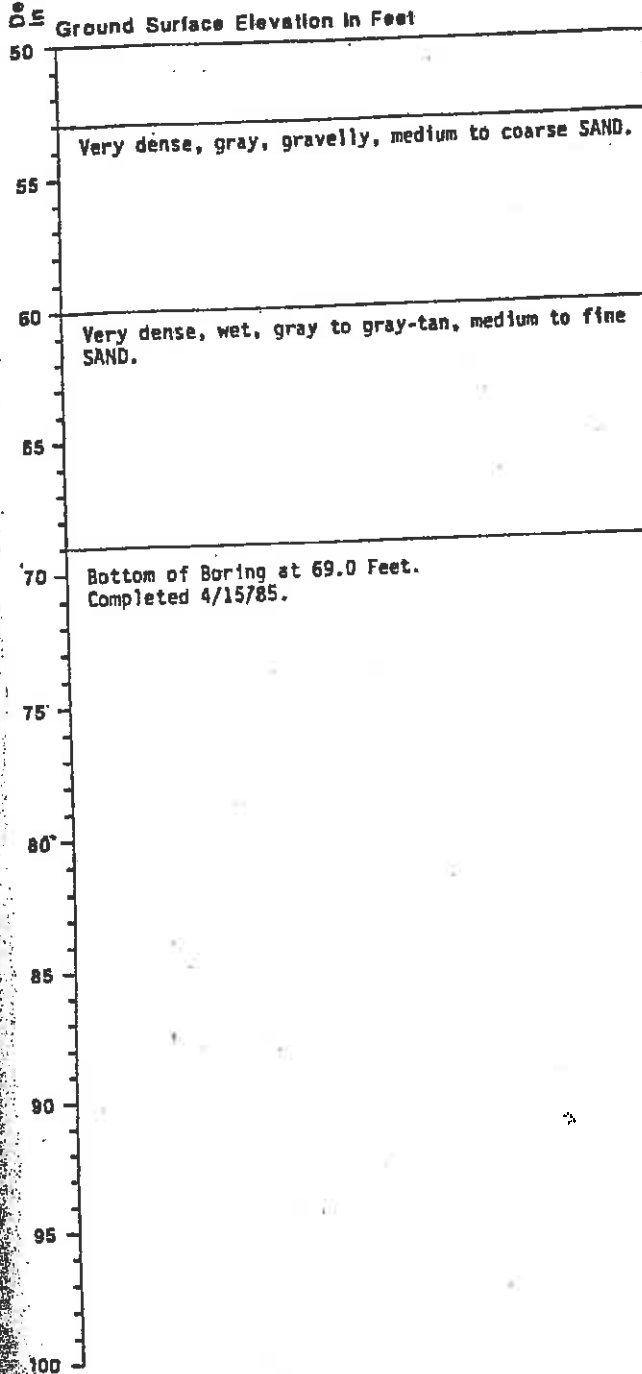
J-1513 April 1985
 HART-CROWSER & associates, inc.
 Sheet 2 of 2 Figure A-4

Boring Log and Construction Data for Well MW-4



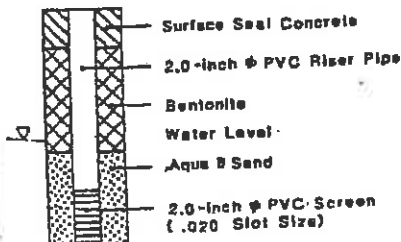
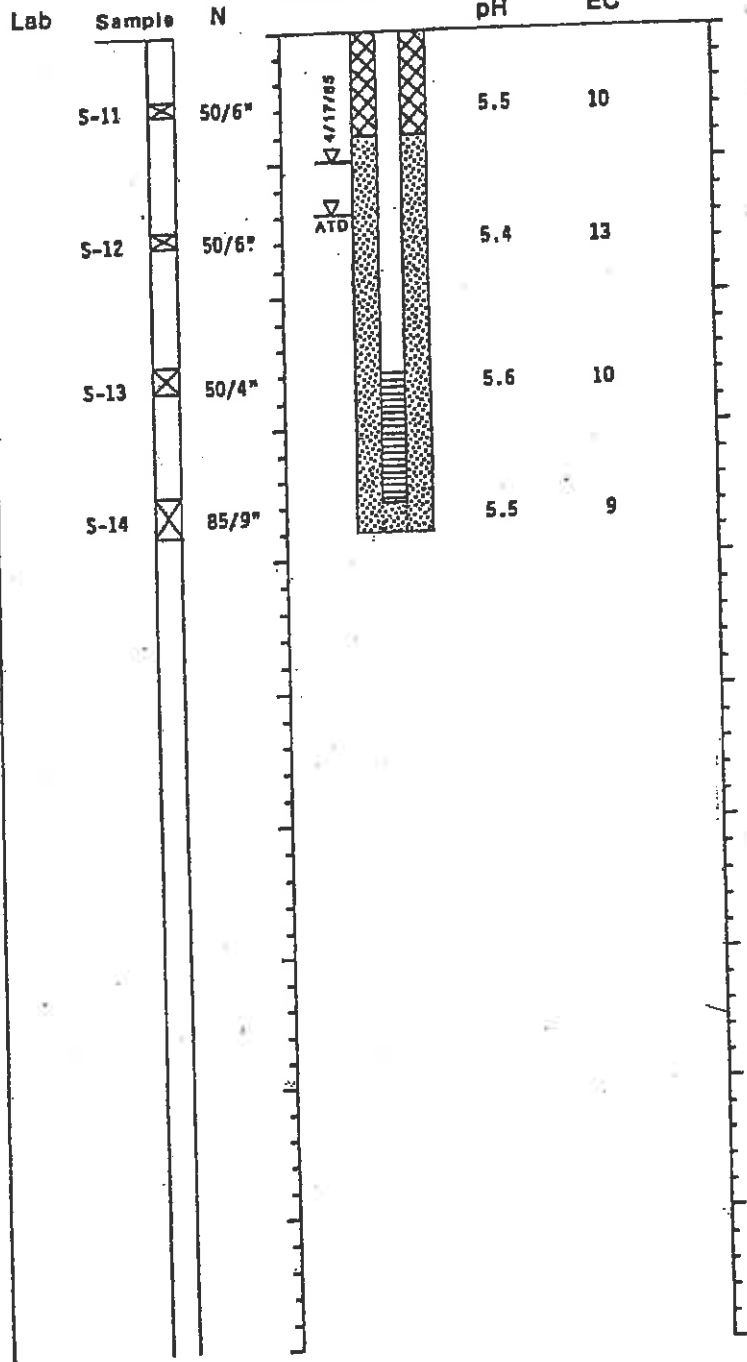
Boring Log and Construction Data for Well MW-4

Geologic Log



Well Design

Top Casing Elevation in Feet
Casing Slickup in Feet



- 2-inch O.D. Split Spoon Sample
- No Sample Recovery
- N Standard Penetration Resistance, Blows per foot
- GS Grain Size Analysis
- X Permeability Test.

NOTES:

1. Soil descriptions are interpretive and actual changes may be gradual.
2. Water Level is for date indicated and may vary with time of year. ATD: At Time of Drilling
3. Top of PVC casing surveyed by Parametrix to estimate mean sea level datum.

J-1513 April 1985
HART-CROWSER & associates, inc.
Sheet 2 of 2 Figure A-5

Well Installation Log

Job No. 31-1578 -25 Client Kitsep County Location Ollala Landfill

| | | | |
|--|------------------------|--|---|
| CASING TYPE <u>2-inch schedule 80 PVC</u> | 10 ft length | DRILLING METHOD <u>Air rotary</u> | WELL NO. <u>MW-E 5A</u> |
| LENGTH CASING <u>0-87ft</u> | Screen <u>87-97 ft</u> | SUMP <u>97-99 1/2</u> | SAMPLING METHOD <u>sediment catch basin</u> |
| JOINT TYPE <u>flush thread</u> | | HAMMER WT. <u>---</u> | DROP <u>---</u> |
| SCREEN TYPE <u>PVC continuous wind</u> | | DATE <u>March 1, 1989</u> | BY <u>James Braddock</u> |
| SLOT SIZE <u>0.020 inch</u> | | DRILLING CONTR. <u>Soil Sampling Service</u> | START <u>Feb 22 1989</u> |
| SEAL TYPE <u>bentonite slurry and pellets</u> | | WATER LEVEL <u>89.7</u> | 89.2 |
| INSTALL METHOD <u>tremie pipe</u> | | TIME <u>0940</u> | 1010 |
| FILTER <u>washed silica sand #10/20</u> | | DATE <u>Feb/28/89</u> | Feb/28/89 |
| INSTALL <u>tremie pipe</u> | | | not static |
| GROUT <u>bentonite slurry (and pellets above sand)</u> | | | |
| | | | FINISH <u>Feb 28/89 drilling</u> |
| | | | <u>Mar 1/89 completed</u> |

| WELL DETAILS | DEPTH | LOG DESCRIPTION | NOTES |
|---|---------|--|--|
| fill casing | 0 | GROUND ELEVATION: | |
| 0-79'2" Above ground to 86'2" bentonite slurry casing | 0-79'2" | 0-2 Organic material | |
| | | 2-6 1/2 GRAVEL with sand & silt | Gravel - Gray, reddish brown, white, brown. Sand - same. Silt - bro |
| | | 6 1/2 - 7 1/2 BOULDER or COBBLES | Granitic material |
| | | 7 1/2 - 16 GRAVEL with sand & silt | Similar to that between 2-6 1/2 : |
| | | | |
| | | 16 - 28 GRAVELLY SAND | Moist, light brown or tan. Water in hole (~2-3 ft) at 16 ft in morning, before drilling. |
| | | | |
| | | 28 - 36 1/2 SILTY GRAVEL | Gravel - dark to medium gray Silt - Grayish brown. Some clay |
| | | | |
| | | 36 1/2 - 48 1/2 CLAY with gravel | Clay - brown. Gravel - blue-gray |
| | | | gradual transition. |
| | | 48 1/2 - 61 CLAY or SILT with some gravel. | Clay or silt - brown. Gravel - blue-gray, less (amount) than above. |
| | | | |
| | | 60 1/2 - 61 - silt layer | |
| | | 61 - 108' SAND | Sand - mostly medium-grained, clear white, lt & dk brown, dark gray - overall brown color. Occasional zones of up to 10-15% gravel - may be bedded |

Well Installation Log

Job No. 31-1578-25

Client Kitsap County

Location Ollala Landfill

| | | | | |
|----------------|--|-----------------|------|----------|
| CASING TYPE | | DRILLING METHOD | | WELL NO. |
| LENGTH | | SAMPLING METHOD | | |
| JOINT TYPE | | HAMMER WT. | DROP | SHEET 2 |
| SCREEN TYPE | | DATE | | OF 2 |
| SLOT SIZE | | BY | | START |
| SEAL TYPE | | DRILLING CONTR. | | FINISH |
| INSTALL METHOD | | WATER LEVEL | | |
| FILTER | | TIME | | |
| INSTALL | | DATE | | |
| GROUT | | | | |

| WELL DETAILS | DEPTH | USCS/ GRAPHIC LOG | LOG DESCRIPTION | NOTES |
|---|-------|-------------------------|----------------------------|---|
| fill casing | | | GROUND ELEVATION: | |
| Above ground to 86'2" | | | | SAND - continued |
| 2" Schedule 80 PVC casing | | | | |
| 79'2" to 82'10" | | | | |
| bentonite pellets | | | | sand becomes a bit coarser, but very gradually |
| 82'10" to 98'7" | | | | (recessional outwash indicate |
| sand filter pack, #10/20 | | | | |
| 86'2" to 96'2" PVC screen 2" OD, 0.020" continuous wound screen | | | | |
| 98'7" to 108' | | | | |
| caved-in sediments | | | | (sand heaves while constructing well, up to 95 foot mark) |
| 96'2" to 98'2" 2" sch. 80 PVC sump | | | | |
| 98'2" to 98'7" - end cap | | | | |
| | | | 108' bottom of well boring | |

SOIL SAMPLING SERVICE, INC.

DATE 11-15-88
 LOG OF BORING NO. 2 WELL ID # 6 START CARD # 011125
 C. NT. Scuccola Const CLIENT JOB # _____ S.S.S. JOB # 10-2581
 LOCATION 114 NE 114 SEC. 1 TOWNSHIP 28 N RANGE 1 E
 CITY Ohalla STATE WASH COUNTY KITSAF
 DRILLING METHOD AUGER DATUM _____ WATER ELEVATION _____
 RIG # 11 TRUCK # 16 OTHER STEAM CLEANER DATE INSTALLED 11-15-88

| TIME & CHARGE DISTRIBUTION | TOTAL TIME | DRILLER <u>TERRY SHEPPY</u> | HOURS <u>9.5</u> | TOTAL FOOTAGE <u>35'</u> |
|----------------------------|------------|---|------------------|------------------------------------|
| DRILLING | <u>2.5</u> | HELPER <u>JOE RICHARD</u> | | |
| CREWED STANDBY | | HELPER | | |
| NON-CREWED STANDBY | | | | |
| MOVING | | | | |
| INSTALLATION | <u>4.5</u> | DRILLER'S SIGNATURE <u>Terry Sheppy</u> <small>The figures herein agree with payments.</small> | | SCALE 1" = <u>5</u> FEET |
| WATER | <u>1.5</u> | BY <u>Terry Scuccola Const</u> Inspector or Client's Representative | | PAGE _____ OF _____ |
| CLEAN + PATCH | | | | |
| REDRILL | | | | |

| | | PVC BLANK <u>3 X 2 inch X 10 FT</u> | ANNULUS MATERIAL | AS BUILT | FORMATION DESCRIPTION | | | |
|---------------|----------|--------------------------------------|------------------|-----------------------|---------------------------|---|----------|------|
| | | PVC SCREEN <u>1 X 2 inch X 10 FT</u> | | | | | | |
| | | MONUMENT TYPE <u>1 stand up</u> | | | | | | |
| | | <u>2 bags 8'</u> | vented SLIC CAP | | | | | |
| | | <u>6 Portland Cement</u> | | CEMENT BENTONITE SEAL | | 8.5' 10-10-12 | | |
| | | <u>1/2 BKT 3/8 PELLETS</u> | | | | 9.5' LT BRN SAND 45 SAND | | |
| | | <u>1/2 kg WYO BENTONITE</u> | | | | 9.0' 9.15-12 | SAME | |
| | | <u>1 kg ENVIRO SEL.</u> | | | | 12.5' | 12-32-24 | SAME |
| | | | | | | 14.5' | | |
| | | | | | 12.5' | 38-58 | | |
| TRAVEL | | | | | 18.5' SAND WITH some sand | | | |
| DEMOB | | | | | 22.5' 18-58 | LT BRN SAND | | |
| SAMPLE METHOD | ATTEMPTS | <u>1 X 2 inch TE CAP</u> | | | 26.5' 20-58 | WATER | | |
| DRIVE SAMPLES | <u>8</u> | <u>1 X 2 inch SLIP CAP, VENTED</u> | | | 28.5' 20-58 | LT BRN TO LT SAND | | |
| GSTERBURG | | <u>1 MONUMENT PROTECTORS</u> | Bentonite Seal | | 29.5' 24-50% | bottom of hole | | |
| | | | SAND PACK | | | water at 23.0 12 LT BRN + LT SAND | | |

COMMENTS

10 FT hard drilling

7:30-9:30, installing #5

9:30-11:00, water

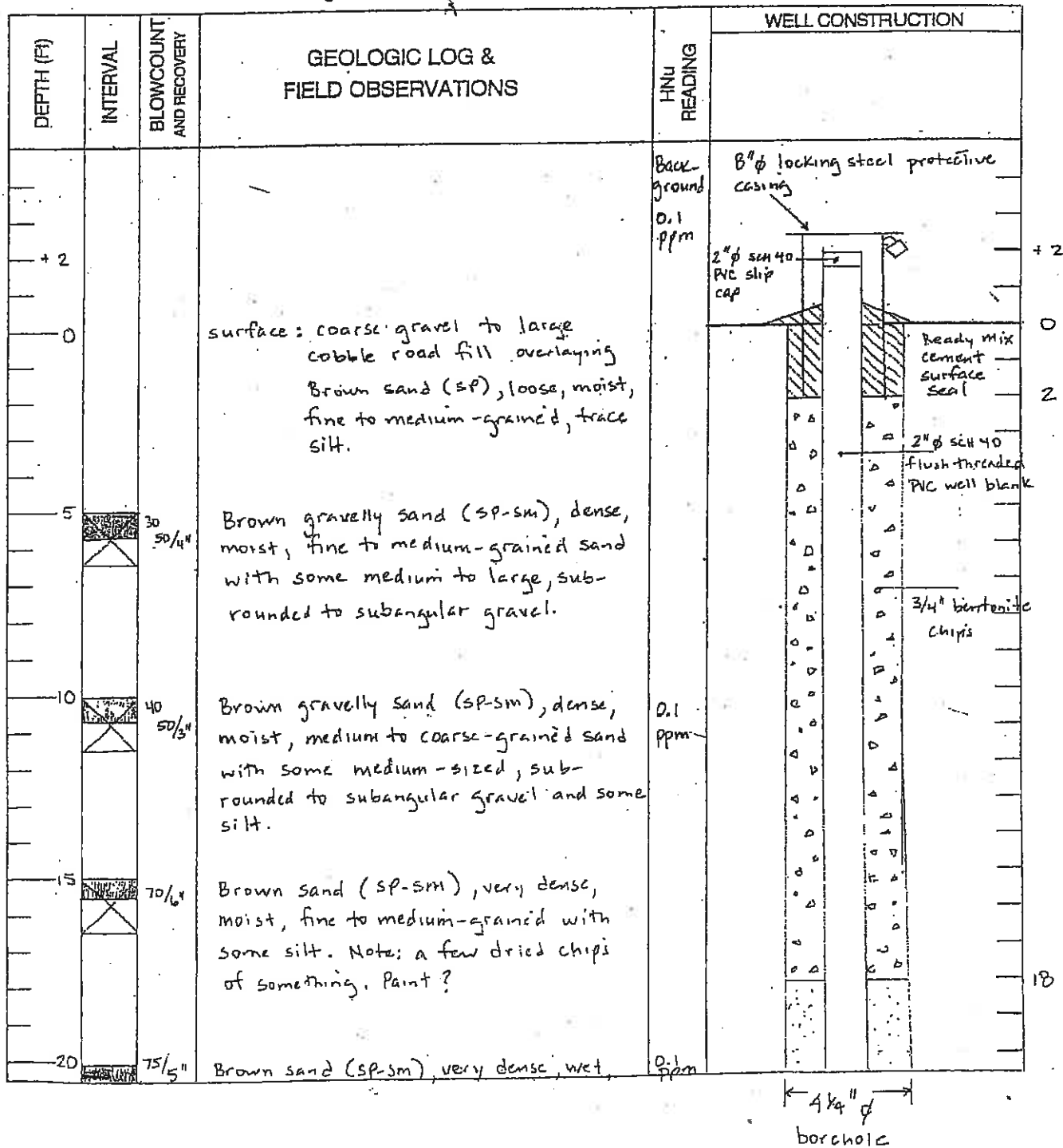
11:00-1:30, drilling #6, from 0' to 35'

1:30-4:00, installing #6

| MONITORING WELL GEOLOGIC & CONSTRUCTION LOG | | |
|---|---------------------|--------------|
| PROJECT NUMBER SEA 35781.82 | WELL NUMBER MW-7 | SHEET 1 OF 2 |

PROJECT Dialla Landfill
 ELEVATION (TOP OF WELL CASING) _____
 WATER LEVEL ELEVATION 26.09 ft. from top of PVC casing
 DRILLING CONTRACTOR Tacoma Pump and Drill
 DRILLING METHOD Hollow stem auger

LOCATION Dialla - Burley Road
 SURFACE ELEVATION _____
 START DATE 12-21-93
 FINISH DATE 12-21-93
 HYDROGEOLOGIST K. Gehweiler



MONITORING WELL GEOLOGIC & CONSTRUCTION LOG

| | | |
|----------------------------------|----------------------|--------------|
| PROJECT NUMBER SEA 3578 I. B2 | WELL NUMBER MW -7 | SHEET 2 OF 2 |
|----------------------------------|----------------------|--------------|

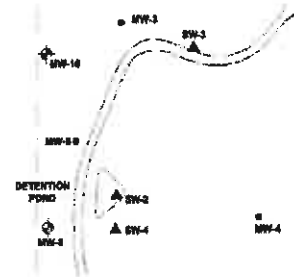
| | |
|---|--------------------------------------|
| PROJECT <u>Olalla Landfill</u> | LOCATION <u>Olalla - Burley Road</u> |
| ELEVATION (TOP OF WELL CASING) | SURFACE ELEVATION |
| WATER LEVEL ELEVATION <u>26.09 ft. from top of PVC casing</u> | START DATE <u>12-21-93</u> |
| DRILLING CONTRACTOR <u>Tacoma Pump and Drill</u> | FINISH DATE <u>12-21-93</u> |
| DRILLING METHOD <u>Hollow stem auger</u> | HYDROGEOLOGIST <u>K. Getweiler</u> |

| DEPTH (FT) | INTERVAL | BLOWCOUNT AND RECOVERY | GEOLOGIC LOG & FIELD OBSERVATIONS | PUMP READING | WELL CONSTRUCTION | |
|------------|----------|------------------------|--|--------------|---|----|
| | | | | | | |
| 21 | X | | fine to medium-grained with some silt. No gravel. Sand becomes saturated at approx. 22 ft. bgs. | | 10-20 sand pack | 21 |
| 25 | X | 75/5 | Brown sand (SP-SM), very dense, saturated, fine to medium-grained with some silt. No gravel. | | 2" φ SCH 40 flush threaded PVC well screen (0.020" slot) | 31 |
| 30 | X | 75/4.5 | Brown sand (SP-SM), very dense, saturated, fine to medium-grained with some silt. No gravel. | 0.2 ppm | 2" φ SCH 40 flush threaded PVC sump | 33 |
| 35 | | | stopped drilling at 33.1 ft. bgs. Water level measured at 30.3 ft bgs at 1100, at 26.09 ft bgs at 1345. | | | |

Boring/Well Designation: MW-8

Client: Kitsap County Public Works
Logged By: E. Caddey, L.G.
Date of Drilling: 10/7/10
Location: Olalla Landfill
Drilling Contractor: Cascade Drilling
Drill Rig: CME-75
Method: Hollow-Stem Auger
Borehole: 8" diameter

Site Representation:

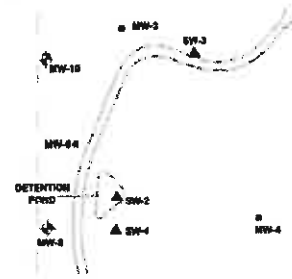


| Depth | SUBSURFACE PROFILE | | | SAMPLE | | PID (ppm) | Well Data | Comments |
|----------|--------------------|-----------|--|----------|----------|-----------|-----------|-------------------------------------|
| | Log | USCS Code | Description | Interval | Recovery | | | |
| 0 | | | Ground Surface | | | | | |
| 0 | | | Gravel | | | | | Above-ground monument with bollards |
| 5 | SW | | Well-Graded Sand with Gravel Tan; moist; mostly fine to coarse sand with little gravel | 5-6-6 | | 5-6-6 | 0.0 | |
| 9-12-15 | | | | 9-12-15 | | 9-12-15 | 0.0 | |
| 50-5" | | | | 50-5" | | 50-5" | 0.0 | |
| 16-50-4" | SP-SM | | Poorly-Graded Sand with Silt and Gravel Brown; slightly moist; mostly medium sand with little silt and little gravel | 16-50-4" | | 16-50-4" | 0.0 | |
| 50-4" | | | | 50-4" | | 50-4" | 0.0 | |
| 50-6" | SP | | Poorly-Graded Sand with Gravel Gray-Brown; moist to wet; mostly medium sand with little gravel at 21' | 50-6" | | 50-6" | 0.0 | 2" Sch 40 PVC blank |
| 10-17-23 | | | | 10-17-23 | | 10-17-23 | 0.0 | Hydrated bentonite chips |
| 6-8-25 | SP | | Poorly-Graded Sand Gray-brown; moist to wet; medium sand with trace gravel and trace silt | 6-8-25 | | 6-8-25 | 0.0 | Water at 22' bgs ATD |
| 12-50-6" | | | Becomes gray at 27' | 12-50-6" | | 12-50-6" | 0.0 | Stainless steel centralizer |
| 8-21-23 | | | Becomes brown at 29' | 8-21-23 | | 8-21-23 | 0.0 | #2/12 silica sand pack |
| 12-16-20 | | | Grades to well-rounded medium sand, black, gray and white grains | 12-16-20 | | 12-16-20 | 0.0 | 0.010" slot sch 40 PVC |
| 2-4-4 | | | | 2-4-4 | | 2-4-4 | 0.0 | Stainless steel centralizer |
| 40 | | | End of Borehole | | | | | |

Boring/Well Designation: MW-10

Client: Kitsap County Public Works
Logged By: E. Caddey, L.G.
Date of Drilling: 10/7/10
Location: Olalla Landfill
Drilling Contractor: Cascade Drilling
Drill Rig: CME-75
Method: Hollow-Stem Auger
Borehole: 8" diameter

Site Representation:



| Depth | SUBSURFACE PROFILE | | | SAMPLE | | PID (ppm) | Well Data | Comments |
|-------|--------------------|-----------|--|----------|----------|-----------|-----------|-------------------------------------|
| | Log | USCS Code | Description | Interval | Recovery | | | |
| 0 | | | Ground Surface | | | | | |
| 0-5 | | | Organic Soil | | | | | Above-ground monument with bollards |
| 5 | SW | | Well-Graded Sand with Gravel Tan; moist; mostly fine to coarse sand with little gravel | 8-18-27 | | 1.2 | | |
| 5 | | | | 12-14-16 | | 4.4 | | |
| 5 | | | | 9-16-19 | | 5.5 | | 2" Sch 40 PVC blank |
| 10 | GP | | Poorly-Graded Gravel Gray; moist; mostly fine gravel with little medium sand | 9-16-18 | | 0.0 | | |
| 10 | | | | 50-6" | | 0.0 | | |
| 15 | SW | | Well-Graded Sand with Gravel Tan; moist; mostly fine to coarse sand with little gravel | 50-6" | | 0.0 | | Hydrated bentonite chips |
| 20 | SP | | Poorly-Graded Sand with Gravel Tan; moist; mostly medium sand with little gravel | 50-6" | | 0.0 | | |
| 20 | SP | | Poorly-Graded Sand Gray-Brown; moist to wet; mostly medium sand with trace gravel and trace silt | 50-5" | | 0.0 | | |
| 25 | | | | 50-6" | | 0.0 | | |
| 30 | | | | 50-6" | | 0.0 | | |
| 35 | | | Becomes saturated at 32.4' | 8-19-20 | | 0.0 | | Water at 32' bgs |
| 40 | | | | 22-50-5" | | 0.0 | | Stainless steel centralizer |
| 40 | | | | | | | | #2/12 silica sand pack |
| 45 | | | | 50-6" | | 0.0 | | 0.010" slot Sch 40 PVC |
| 45 | | | | 12-20-22 | | 0.0 | | Stainless steel centralizer |

Project No. 60101.0

OW-1

22/01-1A

File Original and First Copy with Department of Ecology Second Copy - Owner's Copy Third Copy - Driller's Copy

WATER WELL REPORT STATE OF WASHINGTON

Application No. Permit No.

(1) OWNER: South County Transfer Station Olalla-Burley Landfill (2) LOCATION OF WELL: County Kitsap NE 1/4 NE 1/4 Sec 1 T. 22N., R. 1 W.M.

(3) PROPOSED USE: Domestic [X] Industrial [] Municipal [] Irrigation [] Test Well [] Other []

(4) TYPE OF WORK: Owner's number of well (if more than one) New well [X] Method: Dug [] Bored [] Deepened [] Cable [] Driven [] Reconditioned [] Rotary [X] Jetted []

(5) DIMENSIONS: Diameter of well 6 inches. Drilled 180 ft. Depth of completed well 159 ft.

(6) CONSTRUCTION DETAILS: Casing installed: 6" diam. from 0 ft. to 143 ft. Perforations: Yes [] No [X] Screens: Yes [X] No [] Manufacturer's Name Johnson

Gravel packed: Yes [] No [X] Surface seal: Yes [X] No [] To what depth? 18 ft. Material used in seal Bentomite

(7) PUMP: Manufacturer's Name Berkeley Type: Submersible HP 3

(8) WATER LEVELS: Static level 86 ft. below top of well Date 10-15-84

(9) WELL TESTS: Was a pump test made? Yes [] No [] Yield: gal./min. with ft. drawdown after hrs.

Table with 6 columns: Time, Water Level, Time, Water Level, Time, Water Level. Recovery data (time taken as zero when pump turned off) (water level measured from well top to water level)

Date of test Bailer test 30 gal./min. with 21 ft. drawdown after 1 hrs. Artesian flow g.p.m. Date Temperature of water Was a chemical analysis made? Yes [] No []

(10) WELL LOG: 22-01-1A Formation: Describe by color, character, size of material and structure, and show thickness of aquifers and the kind and nature of the material in each stratum penetrated, with at least one entry for each change of formation.

Table with 3 columns: MATERIAL, FROM, TO. Fill 0-3, Top soil and logs 3-6, Brown clay some gravels sand 6-20, Brown sand water 20-32, Brown clay 32-36, Blue clay 36-42, Brown gravel 42-43, Brown sandy clay 43-59, Brown sand gravel 59-61, Brown sand layered 61-87, Brown sand med. with water 87-115, Brown sand water 115-159, Blue sand water fine layers clay 159-180

Well total 159 ft. 6 1/8 in. To top of packer 142 ft. 7/8 in. Riser pipe 1 ft. 6 in. #.012 slot screen 5 ft. 3 in. #.012 slot screen 5 ft. 2 3/4 in. #.010 slot screen 5 ft. 2 1/2 in. Tailpipe 3 in. Screen overall 17 ft. 5 1/4 in. Six in. casing 143 ft. 8 1/4 in. Static head 86 ft. 20 G.P.M. 99 ft. 6 in. 30 G.P.M. 107 ft. 6 in.

DEPARTMENT OF ECOLOGY NORTHWEST REGION Work started 10-11-84 Completed 10-15-84

WELL DRILLER'S STATEMENT: This well was drilled under my jurisdiction and this report is true to the best of my knowledge and belief.

NAME Nicholson Drilling Company (Person, firm, or corporation) (Type or print) Address P.O. Box 123 Port Orford, WA 98366 [Signed] (Well Driller)

License No. 0520 Date 10-23-83

(USE ADDITIONAL SHEETS IF NECESSARY)

The Department of Ecology does NOT warranty the Data and/or the Information on this Report.



WATER WELL REPORT

Original & 1st copy - Ecology, 2nd copy - owner, 3rd copy - driller

297695

OW-5

22-1E-1B

Construction/Decommission ("x" in circle)

- Construction
 Decommission ORIGINAL INSTALLATION Notice of Intent Number _____

CURRENT

Notice of Intent No. W216878
 Unique Ecology Well ID Tag No. APR627
 Water Right Permit No. _____
 Property Owner Name Gene Ryker
 Well Street Address 13041 Olympic Dr. SE
 City Olalla County Kitsap
 Location NW 1/4-1/4 NE 1/4 Sec 1 Twn 22 NR 1E EWM or WWM circle one
 Lat/Long (s, t, r) Lat Deg _____ Lat Min/Sec _____
 Still REQUIRED) Long Deg _____ Long Min/Sec _____
 Tax Parcel No. 01220110322008

PROPOSED USE: Domestic Industrial Municipal
 DeWater Irrigation Test Well Other _____

TYPE OF WORK: Owner's number of well (if more than one) _____
 New well Reconditioned Method: Dug Bored Driven
 Deepened Cable Rotary Jetted

DIMENSIONS: Diameter of well -- 6 inches, drilled 279 ft.
 Depth of completed well 279 ft.

CONSTRUCTION DETAILS
 Casing Welded 6 " Diam. from +1 ft. to 279 ft.
 Installed: Liner installed " Diam. from _____ ft. to _____ ft.
 Threaded " Diam. from _____ ft. to _____ ft.

Perforations: Yes No
 Type of perforator used _____
 SIZE of perfs _____ in. by _____ in. and no. of perfs _____ from _____ ft. to _____ ft.

Screens: Yes No K-Pac Location 274
 Manufacturer's Name Johnson
 Type Stainless steel Model No. _____
 Diam. 5 Slot size 18 from 274 ft. to 279 ft.
 Diam. _____ Slot size _____ from _____ ft. to _____ ft.

Gravel/Filter packed: Yes No Size of gravel/sand _____ ft.
 Materials placed from _____ ft. to _____ ft.

Surface Seal: Yes No To what depth? 18 ft.
 Material used in seal Bentonite
 Did any strata contain unusable water? Yes No
 Type of water? _____ Depth of strata _____
 Method of sealing strata off _____

PUMP: Manufacturer's Name Flint Walling
 Type: Submersible H.P. 1 1/2

WATER LEVELS: Land-surface elevation above mean sea level _____ ft.
 Static level 100 ft. below top of well Date 5/28/08
 Artesian pressure _____ lbs. per square inch Date _____
 Artesian water is controlled by _____ (cap, valve, etc.)

WELL TESTS: Drawdown is amount water level is lowered below static level
 Was a pump test made? Yes No If yes, by whom? _____
 Yield: 22 gal./min. with 6 ft. drawdown after 1 hrs.
 Yield: _____ gal./min. with _____ ft. drawdown after _____ hrs.
 Yield: _____ gal./min. with _____ ft. drawdown after _____ hrs.
 Recovery data (time taken as zero when pump turned off) (water level measured from well top to water level)

| Time | Water Level | Time | Water Level | Time | Water Level |
|------|-------------|------|-------------|------|-------------|
| 0 | 106' | 3 m | 101' | 10 m | 100 |
| 1 m | 102' | 4 m | 101' | | |
| 2 m | 101' | 5 m | 101' | | |

 Date of test 5/28/08
 Bailer test 12 gal./min. with 3 ft. drawdown after 2 hrs.
 Airtest _____ gal./min. with stem set at _____ ft. for _____ hrs.
 Artesian flow _____ g.p.m. Date _____
 Temperature of water _____ Was a chemical analysis made? Yes No

CONSTRUCTION OR DECOMMISSION PROCEDURE

Formation: Describe by color, character, size of material and structure, and the kind and nature of the material in each stratum penetrated, with at least one entry for each change of information. (USE ADDITIONAL SHEETS IF NECESSARY.)

| MATERIAL | FROM | TO |
|------------------------|------|-----|
| Brown top | 0 | 2 |
| Brown sand/clay | 2 | 57 |
| Grey clay/sand | 57 | 82 |
| Grey silt | 82 | 110 |
| Grey clay/fine sand | 110 | 226 |
| Brown clay | 226 | 227 |
| Grey clay/fine sand | 227 | 273 |
| Grey sand w/b | 273 | 275 |
| Grey sand & gravel w/b | 275 | 279 |
| Grey till | 279 | |

RECEIVED

JUN 11 2008

DEPT. OF ECOLOGY

Start Date 5/6/08 Completed Date 5/28/08

WELL CONSTRUCTION CERTIFICATION: I constructed and/or accept responsibility for construction of this well, and its compliance with all Washington well construction standards. Materials used and the information reported above are true to my best knowledge and belief.

Driller Engineer Trainee Name (Print) Matt Olsen
 Driller/Engineer/Trainee Signature Matt Olsen
 Driller or trainee License No. 2337

Drilling Company Olsen Drilling
 Address PO Box 1554
 City, State, Zip: Port Orchard, WA 98366
 Contractor's Registration No. OLSEND101LJ Date 6/10/08

If TRAINEE:
 Driller's Licensed No. _____
 Driller's Signature _____

Ecology is an Equal Opportunity Employer.

WATER WELL REPORT

STATE OF WASHINGTON

Start Card No. 29474
Water Right Permit No. 22.E.1F

(1) OWNER: Name Shoemaker Address _____
(2) LOCATION OF WELL: County Kitsap SE & NW & Sec 1 T. 22N., R. 1E W.M.
(2a) STREET ADDRESS OF WELL (or nearest address) 13320 Olympic Port Orchard

(3) PROPOSED USE: Domestic Industrial Municipal
 Irrigation Test Well Other
 DeWater

(4) TYPE OF WORK: Owner's number of well (if more than one) _____
Abandoned New well Deepened Reconditioned Method: Dug Cable Rotary Bored Driven Jettied

(5) DIMENSIONS: Diameter of well 6 inches.
Drilled _____ feet. Depth of completed well 61 ft.

(6) CONSTRUCTION DETAILS:
Casing installed: 6 ft. Diam. from 0 ft. to 56 ft.
Welded _____ ft. Diam. from _____ ft. to _____ ft.
Liner installed _____ ft. Diam. from _____ ft. to _____ ft.
Threaded _____ ft. Diam. from _____ ft. to _____ ft.
Perforations: Yes No
Type of perforator used _____
SIZE of perforations _____ in. by _____ in.
_____ perforations from _____ ft. to _____ ft.
_____ perforations from _____ ft. to _____ ft.
_____ perforations from _____ ft. to _____ ft.

Screens: Yes No
Manufacturer's Name Howard Smith
Type 35 Model No. _____
Diam. 5 Slot size 16 from 56 ft. to 61 ft.
Diam. _____ Slot size _____ from _____ ft. to _____ ft.

Gravel packed: Yes No Size of gravel _____
Gravel placed from _____ ft. to _____ ft.
Surface seal: Yes No To what depth? 18 ft.
Material used in seal Bentonite
Did any strata contain unusable water? Yes No
Type of water? _____ Depth of strata _____
Method of sealing strata off _____

(7) PUMP: Manufacturer's Name _____
Type: _____ H.P. _____

(8) WATER LEVELS: Land-surface elevation above mean sea level _____ ft.
Static level 32 ft. below top of well Date 5/4
Artesian pressure _____ lbs. per square inch Date _____
Artesian water is controlled by _____ (Cap, valve, etc.)

(9) WELL TESTS: Drawdown is amount water level is lowered below static level
Was a pump test made? Yes No If yes, by whom? _____
Yield: _____ gal./min. with _____ ft. drawdown after _____ hrs.

| Recovery data (time taken as zero when pump turned off) (water level measured from well top to water level) | | | | | |
|---|-------------|------|-------------|------|-------------|
| Time | Water Level | Time | Water Level | Time | Water Level |
| | | | | | |
| | | | | | |

Date of test _____
Ball test 12 gal./min. with 6' ft. drawdown after 2 hrs.
Airtest _____ gal./min. with stem set at _____ ft. for _____ hrs.
Artesian flow _____ g.p.m. Date _____
Temperature of water _____ Was a chemical analysis made? Yes No

(10) WELL LOG or ABANDONMENT PROCEDURE DESCRIPTION
Formation: Describe by color, character, size of material and structure, and show thickness of aquifers and the kind and nature of the material in each stratum penetrated, with at least one entry for each change of information.

| MATERIAL | FROM | TO |
|-------------------|------|----|
| Brown Top | 0 | 3 |
| Br Sand | 3 | 8 |
| Br Till | 8 | 18 |
| Br Sand Clay | 18 | 40 |
| Br Sand Clay Seep | 40 | 51 |
| Br Sand w/B | 51 | 61 |

RECEIVED
 MAY 08 1992
 DEPT. OF ECOLOGY

Work started 4/30, 19. Completed 5/4, 1992

WELL CONSTRUCTOR CERTIFICATION:
I constructed and/or accept responsibility for construction of this and its compliance with all Washington well construction standards. Materials used and the information reported above are true to knowledge and belief.

NAME Olsen Drilling
(PERSON, FIRM, OR CORPORATION)
Address 1598 SE Oregon Por
(Signed) Rupe Howie
(WELL DRILLER)
Contractor's Registration No. 0752ND101T Date _____

(USE ADDITIONAL SF)



AQUATIC RESEARCH INCORPORATED

LABORATORY & CONSULTING SERVICES

3927 AURORA AVENUE NORTH, SEATTLE, WA 98103

PHONE: (206) 632-2715 FAX: (206) 632-2417

ASTM D422 PARTICLE SIZE

ANALYSIS DATE: 11/2/2010

LAB SAMPLE ID: BKS00142A1

ANALYST: S. NIELSON

CLIENT SAMPLE ID: OL-MW-08-5

TOTAL WET WEIGHT (g) 117.28
 TOTAL DRY WEIGHT (g) 93.64

PERCENT SOLIDS @ 105C 79.84%

SIEVE FRACTION - SAND GRAVEL

BEAKER # 1

SIEVE FRACTION

| | |
|------------------|----------|
| BEAKER TARE: | 112.6400 |
| TOTAL DRY WT (g) | 203.7400 |
| SAMPLE WT (g) | 91.1000 |

| SIEVE SIZE | (mm) | (g) | (%) | % FINER |
|------------|--------|---------|--------|---------|
| 0.375 | 9.50 | 0.0000 | 0.00% | 100.00% |
| 0.25 | 6.30 | 0.0000 | 0.00% | 100.00% |
| 4 | 4.75 | 0.0000 | 0.00% | 100.00% |
| 10 | 2.00 | 0.2700 | 0.29% | 99.71% |
| 20 | 0.850 | 0.4300 | 0.46% | 99.25% |
| 40 | 0.425 | 2.9600 | 3.16% | 96.09% |
| 60 | 0.250 | 37.7630 | 40.33% | 55.76% |
| 140 | 0.106 | 44.7900 | 47.83% | 7.93% |
| 200 | 0.075 | 2.9600 | 3.16% | 4.77% |
| <200 | <0.075 | 2.1900 | | |

COMMENTS

TOTAL WT 91.3630
 RECOVERY (%) 100.29%

HYDROMETER FRACTION - SILT CLAY

| SAMPLE TIME (20C) | BLANK | Δ wt (g) | (%) | % FINER |
|-------------------|------------|----------|--------|---------------|
| | 1.0 | | | |
| TIME MIN | HYDROMETER | TEMP C | K | DIAM (mm) |
| 2 | 6.0 | 21.0 | 0.0137 | 0.0378 |
| 5 | 5.0 | | 0.0137 | 0.0240 |
| 15 | 4.0 | | 0.0137 | 0.0139 |
| 30 | 3.5 | | 0.0137 | 0.0099 |
| 60 | 3.0 | | 0.0137 | 0.0070 |
| 250 | 2.0 | 21.0 | 0.0137 | 0.0035 |
| 1440 | 2.0 | 21.0 | 0.0137 | 0.0014 |

PHYSICAL PROPERTIES

DISTRIBUTION % GRAVEL 0.29% % SAND 94.94% % CLAY/SILT 4.77% MEAN (mm) 0.1796

% SOLIDS 79.84% % WATER 20.16%



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PHONE: (206) 632-2715 FAX: (206) 632-2417

CLIENT SAMPLE ID: OL-MW-08-5





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LABORATORY & CONSULTING SERVICES
 3927 AURORA AVENUE NORTH, SEATTLE, WA 98103
 PHONE: (206) 632-2715 FAX: (206) 632-2417

ASTM D422 PARTICLE SIZE

ANALYSIS DATE: 11/2/2010

LAB SAMPLE ID: BKS00142A2

ANALYST: S. NIELSON

CLIENT SAMPLE ID: OL-MW-10-5

TOTAL WET WEIGHT (g) 119.78
 TOTAL DRY WEIGHT (g) 97.68

PERCENT SOLIDS @ 105C 81.55%

SIEVE FRACTION - SAND GRAVEL

BEAKER # 2

SIEVE FRACTION

| | |
|------------------|----------|
| BEAKER TARE: | 117.4700 |
| TOTAL DRY WT (g) | 212.5600 |
| SAMPLE WT (g) | 95.0900 |

| SIEVE SIZE | (mm) | (g) | (%) | % FINER |
|------------|--------|---------|--------|---------|
| 0.375 | 9.50 | 0.0000 | 0.00% | 100.00% |
| 0.25 | 6.30 | 0.0000 | 0.00% | 100.00% |
| 4 | 4.75 | 0.0000 | 0.00% | 100.00% |
| 10 | 2.00 | 0.4800 | 0.49% | 99.51% |
| 20 | 0.850 | 1.1300 | 1.16% | 98.35% |
| 40 | 0.425 | 11.5600 | 11.83% | 86.52% |
| 60 | 0.250 | 40.1000 | 41.05% | 45.47% |
| 140 | 0.106 | 37.8300 | 38.73% | 6.74% |
| 200 | 0.075 | 2.3500 | 2.41% | 4.33% |
| <200 | <0.075 | 1.6200 | | |

COMMENTS

TOTAL WT 95.0700
 RECOVERY (%) 99.98%

HYDROMETER FRACTION - SILT CLAY

| SAMPLE TIME (20C) | | BLANK | |
|-------------------|------------|--------|--------|
| TIME MIN | HYDROMETER | TEMP C | K |
| | | | 1.0 |
| 2 | 6.0 | 21.0 | 0.0137 |
| 5 | 5.0 | | 0.0137 |
| 15 | 5.0 | | 0.0137 |
| 30 | 4.0 | | 0.0137 |
| 60 | 3.5 | | 0.0137 |
| 250 | 3.0 | 21.0 | 0.0137 |
| 1440 | 3.0 | 21.0 | 0.0137 |

| Δ wt (g) | (%) | % FINER |
|----------|-------|---------|
| 5.0000 | 5.12% | 5.12% |
| 4.0000 | 1.02% | 4.09% |
| 4.0000 | 0.00% | 4.09% |
| 3.0000 | 1.02% | 3.07% |
| 2.5000 | 0.51% | 2.56% |
| 2.0000 | 0.51% | 2.05% |
| 2.0000 | 0.00% | 2.05% |

PHYSICAL PROPERTIES

DISTRIBUTION % GRAVEL 0.49% % SAND 95.18% % CLAY/SILT 4.33% MEAN (mm) 0.2178

% SOLIDS **81.55%** % WATER 18.45%



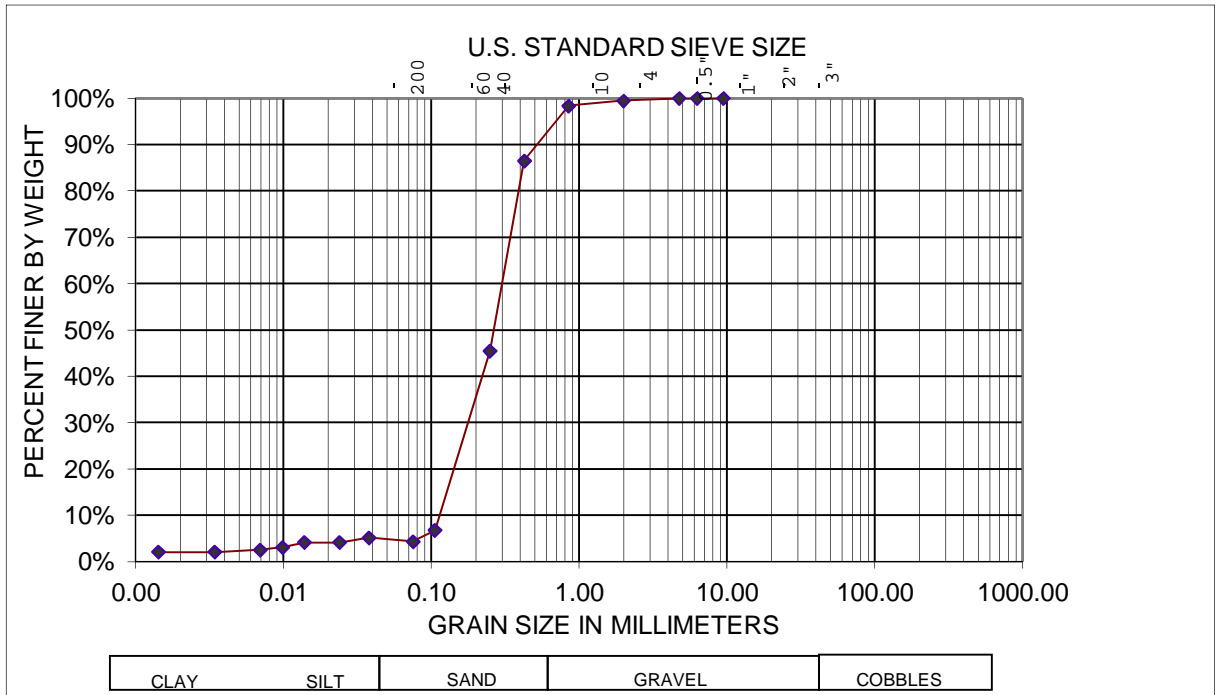
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PHONE: (206) 632-2715 FAX: (206) 632-2417

CLIENT SAMPLE ID: OL-MW-10-5





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3927 AURORA AVENUE NORTH, SEATTLE, WA 98103

PHONE: (206) 632-2715 FAX: (206) 632-2417

ASTM D422 PARTICLE SIZE

ANALYSIS DATE: 11/2/2010

LAB SAMPLE ID: BKS00142A2-dup

ANALYST: S. NIELSON

CLIENT SAMPLE ID: OL-MW-10-5-dup

TOTAL WET WEIGHT (g) 117.47
 TOTAL DRY WEIGHT (g) 96.51

PERCENT SOLIDS @ 105C 82.16%

SIEVE FRACTION - SAND GRAVEL

BEAKER # 2

SIEVE FRACTION

BEAKER TARE: #####
 TOTAL DRY WT (#####)
 SAMPLE WT (g) 92.9700

| SIEVE SIZE | (mm) | (g) | (%) | % FINER | COMMENTS |
|------------|--------|---------|--------|---------|----------|
| 0.375 | 9.50 | 0.0000 | 0.00% | 100.00% | |
| 0.25 | 6.30 | 0.0000 | 0.00% | 100.00% | |
| 4 | 4.75 | 0.3400 | 0.35% | 99.65% | |
| 10 | 2.00 | 0.7700 | 0.80% | 98.85% | |
| 20 | 0.850 | 1.2200 | 1.26% | 97.59% | |
| 40 | 0.425 | 15.4300 | 15.99% | 81.60% | |
| 60 | 0.250 | 41.3500 | 42.84% | 38.75% | |
| 140 | 0.106 | 29.8300 | 30.91% | 7.85% | |
| 200 | 0.075 | 2.0700 | 2.14% | 5.70% | |
| <200 | <0.075 | 1.9800 | | | |

TOTAL WT 92.9900
 RECOVERY (%) 100.02%

HYDROMETER FRACTION - SILT CLAY

| SAMPLE TIME (20C) | | BLANK | | DIAM (mm) | Δ wt (g) | (%) | % FINER |
|-------------------|------------|--------|--------|-----------|----------|-------|---------|
| TIME MIN | HYDROMETER | TEMP C | K | | | | |
| | | | | 1.0 | | | |
| 2 | 6.0 | 21.0 | 0.0137 | 0.0378 | 5.0000 | 5.18% | 5.18% |
| 5 | 5.0 | | 0.0137 | 0.0240 | 4.0000 | 1.04% | 4.14% |
| 15 | 5.0 | | 0.0137 | 0.0139 | 4.0000 | 0.00% | 4.14% |
| 30 | 4.0 | | 0.0137 | 0.0099 | 3.0000 | 1.04% | 3.11% |
| 60 | 3.5 | | 0.0137 | 0.0070 | 2.5000 | 0.52% | 2.59% |
| 250 | 3.0 | 21.0 | 0.0137 | 0.0034 | 2.0000 | 0.52% | 2.07% |
| 1440 | 3.0 | 21.0 | 0.0137 | 0.0014 | 2.0000 | 0.00% | 2.07% |

PHYSICAL PROPERTIES

DISTRIBUTION % GRAVEL 1.15% % SAND 93.15% % CLAY/SILT 5.70% MEAN (mm) 0.2552

% SOLIDS 82.16% % WATER 17.84%

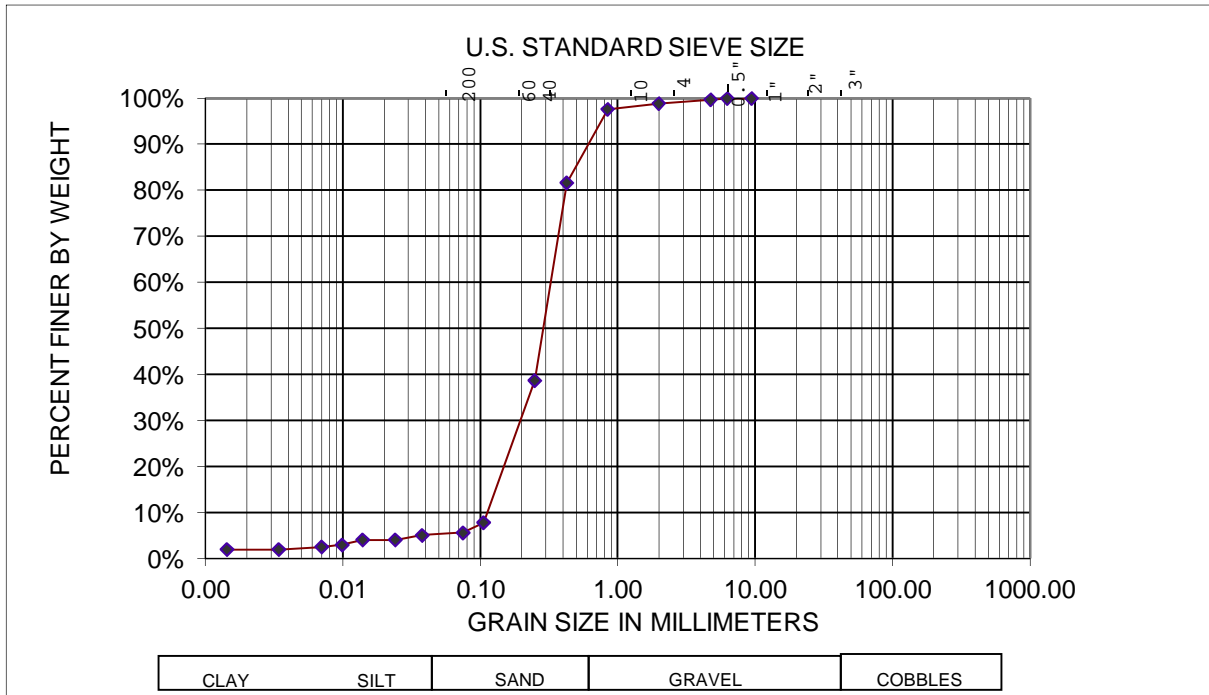


AQUATIC RESEARCH INCORPORATED
LABORATORY & CONSULTING SERVICES

3927 AURORA AVENUE NORTH, SEATTLE, WA 98103

PHONE: (206) 632-2715 FAX: (206) 632-2417

CLIENT SAMPLE ID: **OL-MW-10-5-dup**



60101.0

Clear, 65-70°F Today.

E. Caddy w/ EPT.

Olalla Landfill, SWC of Budy Olalla Rd + Bendix Rd

07:00 - E. Caddy arrives at site, Glacier operator onsite, no equipment yet.

07:10 - Mike Washel w/ Parametrix onsite, open gates on Bendix Rd side.

07:30 - John w/ UDS onsite to check ~~address~~ for buried utilities. Checked drill sites for MU-10 and MU-8, no utilities found. Mike w. had UDS check per meter of area to be trenched, no utilities found.

07:50 - Mike + Glacier operator go down to MU-10 location and start cutting trees w/ chainsaw.

08:15 - Kelli w/ Kitsap Co. Public Works onsite.

08:25 - John w/ UDS offsite.

08:50 - Truck + pup w/ gravel onsite, no machinery yet.

09:00 - Randal Rand + gravel T&T offsite, - flatbed onsite w/ excavator

09:15 - Jan w/ Health Dept onsite, Noel w/ Ecology onsite.

09:30 - Mary Holder EPT onsite

10/7/10

- Mike & Keli made decision w/ Noel's approval to start drilling MW-8 first while MW-10 road & pad being constructed
- 10:00 - Tail gate Health & Safety meeting
- 10:27 - Start drilling MW-8, Mike took Mary up to sample MW-2 & MW-4.
- 12:00 - Almost done w/ MW-8, hydrated chips well specs: - Screen 27-37' olo slot, 2/12 silica sand 26-37' hydrated chips 26-5' gravel to surface.
- Glacier finished w/ drill rod to MW-10, needs Cascade to look at and approve.
 - Noel w/ Ecology offsite at ~11:20, he stated he will be back ~1:00
 - No PFD ready above 0.0
 - Cascade down again in high pressure wash steam clean trailer.
- 12:45 - Noel back on site
- 13:00 - Mike & Mary took truck up to start digging trenches.
- Mike & Mary finished sampling both MW-2 & MW-4, Mike will take samples to lab.
 - crew installing ballards around MW-8
- 13:45 - Move rig onto MW-10 pad.
- 14:00 - Start drilling

- 14:48 - DTW at 32.4 in MW-10, Noel offsite. will drill to 47' & set well.
- Noel back, locked in
- 15:00 - Noel returned keys offsite
- 15:10 - TD MW-10 at 47'
- 15:45 - Mike offsite w/ samples, talking to lab.
- 16:00 - Keli offsite, will be back in the morning for the keys.
- Glacier Tech still backfilling excavations, Mary H. supervising.
- 17:10 - Cascade offsite, Mary Holder offsite Bon w/ Glacier filling in pits - has 1 pit to go. EC to stay and lock gate.
- 17:45 - Bon w/ Glacier offsite, lot 1 excavator unit, not sure who is calling in the pickup.
- 17:47 - EC lock gate, offsite.

7. 

10/8/10

E. Caddy w/EPI

07:30 - arrive To develop ground water wells, Cascade crew arrive, will set pump in MW-8 & MW-10 each.
 DPW - 22.82 from TOC MW-8
 MW-10 - 32.12 from TOC

08:05 - Tailgate H&S meeting
 - down pumps for use in MW-8, new pump in MW-10

collect DPW measurements prior the develop, down water level meter prior to each use.
 - see Monitor Well Development sheets for field measurements.

- In general, set pump in middle of screen & moved down to remove sed in bottom.

Lot clear, then surge up & down. Lot clear, move pump towards top, surge, lot clear. move pump in middle, surge & lot clear.

MW-8 removed \approx 80 gallons until it cleared & no sediment fines.

MW-10 - removed about 50 gallons, was very clear.

- EC measured total depth of each well, MW-8 was at \approx 35' bgs, not 37 ft.

David w/ Cascade stated he knew it was

10/8/10 7

at 37' bgs.

\approx 10:00 - Done with MW-8 & 10

Drum Count - MW-8 - 2 develop water
 - Keli w/ KC arrive 2 soils
 1 down.
 MW-10 - 2 soils
 1 develop water

- Crew moved to develop MW-2 & MW-4, started w/ MW-4. Total depth of well 78', DPW 66' bgs. David said the electrical wires on pump don't go that deep - we measured them - 62' long, won't make it to water

- I called D. Kunkel, we will have to develop wells next week ourselves. Tell Cascade to leave 2 drums.

10:20 - crew cleaning up, loading equip.
 EC check Area/Truck by MW-3 for sample locations.

11:00 - Cascade arrive

- EC arrive, return PWD to rental business in Lakewood.

} T. [Signature]

8 60101.0

10/13/10

Sunny, 60°F today.

E. Caddy w/ EPT

08:00 - E. Caddy on site to develop MW-2 and MW-4. Meet Keli, she opened gate & provided keys to wells.

- set up on MW-4 first. Start w/ stainless steel bailer, remove sand from bottom of well, with 55 gallon, switch to using the Waterloo in-situ pump. Remove an additional 5 gallons, clean, check NTA and parameter w/ Hanna multi-meter. (See field log table.)

- signed w/ the ss bailer, says lock for the Waterloo was too tight fitting in well - I did not want to stick it.

Develop water placed into 55-gallon drum along road.

10:00 - set up on MW-2, ^{used} ~~needed~~ step ladder to reach top of steel up monument that is 5' above the ground. Again, started w/ ss bailer (do not fit) to remove sand in bottom. Removed ~ 8 gallons, put in Waterloo, it would not work - filled up w/ sand. Removed and continued w/ bailer.

10:00 - up to 10 gallons removed, still sandy, some of the brown "gelatin" stuff on side of bailer. Check parameters,

60101.0

10/13/10⁹

Then continue develop w/ bailer. 12:35 - 15 gallons removed, still cloudy w/ some sand. Arms exhausted. May need to come back w/ a submersible pump to clean out well completely.

- Monroed state up monument - hole between well casing and monument - it is open to 5' below ground surface. ~~Fit~~ Both wells have the grey PVC on top.
- Develop water from both wells placed into one 55 gallon drum, labeled, and placed along gravel perimeter road.
- 13:00 - EC off site, "dummy" lock gate as requested by Keli so rental truck can get in to pick up excavator.

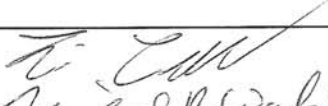
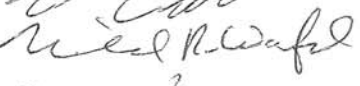
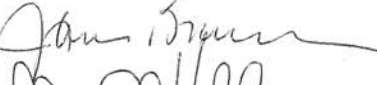








Environmental Partners, Inc.

Tail Gate Health and Safety Meeting

Olalla Landfill Drilling

Site Personnel

| Printed Name and Company | Approvals Signature | Date |
|--------------------------|---|------------|
| Eric Caddy EPA |  | 10/7/10 |
| Mike Ward Parvato |  | 10/7/2010 |
| Jan Brewer |  | 10/7/2010 |
| Mary Holder |  | 10/7/10 |
| KELI McKay-MEANS |  | 10/7/10 |
| Noel Philip |  | 10/07/2010 |
| RACE LAUER COOP |  | 10/07/10 |
| DAVID GOSSE |  | 10/7/10 |
| Daryl Dietrich |  | 10/7/10 |

Monitoring Well Development Field Measurements Record

Job Name/Location: Olalla Landfill RI/FS
 Job Number: 60101.0 Task 3
 Date: 10/8/10

Well #: MW-8

Page Number: 1

| Time | Depth to Water (ft.) | Total Depth of Well (ft.) | Volume Purged (gal.) | pH | Cond. | Turbidity (NTU) | Temp. (oC) | Visual Observations |
|-------|----------------------|---------------------------|----------------------|------|----------------------------------|-----------------|------------|---------------------------------|
| 08:00 | 22.81 | 27.99 | — | — | 115 EC 10/19/10 | — | — | — |
| 08:08 | | 27.85 from | 10 | | | | | |
| 08:17 | | TUC | 15 | | | | | |
| 08:45 | | | 55 | | | | | cloudy, some fines |
| 09:05 | | | 75 | 6.67 | 115 | 29.4 | 10.8 | cloudy, some fines after sample |
| 09:25 | 22.84 | | 85 | 6.50 | 134 | 19.6 | 10.2 | clear |
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1 well volume = 2.41 gallons
 10 well volume = 24.10 gallons



Monitoring Well Development Field Measurements Record

Job Name/Location: Olalla Landfill RI/FS
 Job Number: 60101.0 Task 3
 Date: 10/8/10

Well #: MW-10

Page Number: 1

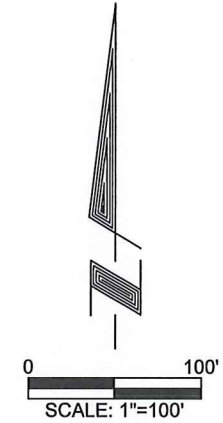
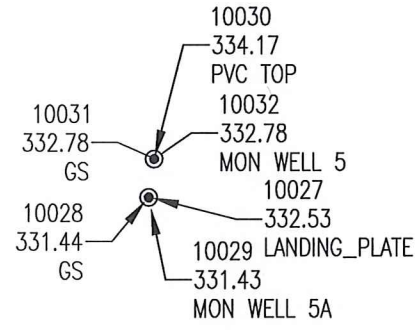
| Time | Depth to Water (ft.) | Total Depth of Well (ft.) | Volume Purged (gal.) | pH | Cond. | Turbidity (NTU) | Temp. (oC) | Visual Observations |
|-------|----------------------|---------------------------|----------------------|------|-----------------------|-----------------|------------|----------------------|
| 08:12 | 32.10 | 49.92 from | — | — | 276 276 | — | — | |
| 08:22 | | TOC | 15 | | | | | cloudy |
| 08:38 | | | 20 | | | | | cloudy after surging |
| 08:55 | | | 35 | 7.80 | 276 | 1.76 | — | clear |
| 09:06 | 32.10 | | 50 | 7.62 | 285 | 1.91 | 12.2 | clear |
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1 well volume = 2.53 gallons
 10 well volumes = 25.3 gallons.

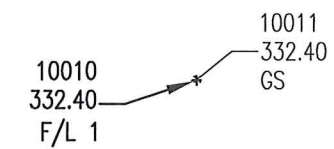
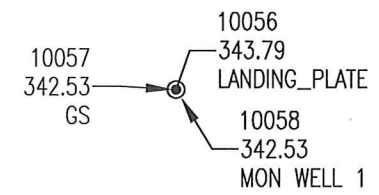
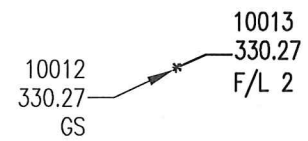
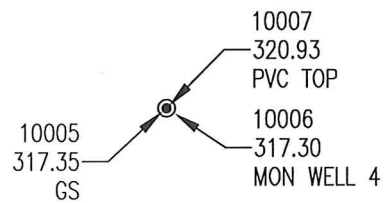
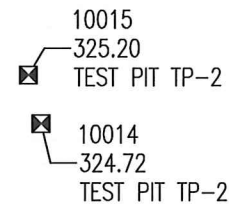
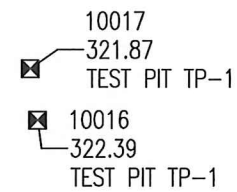
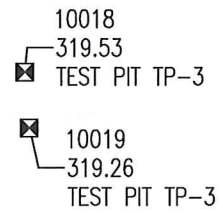
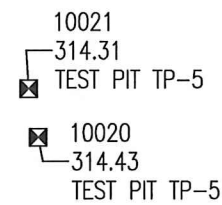
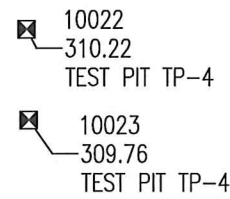
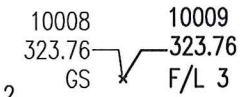
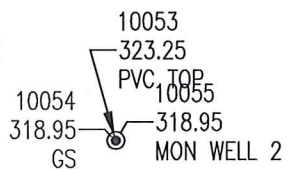
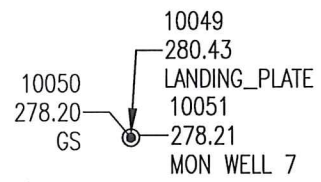
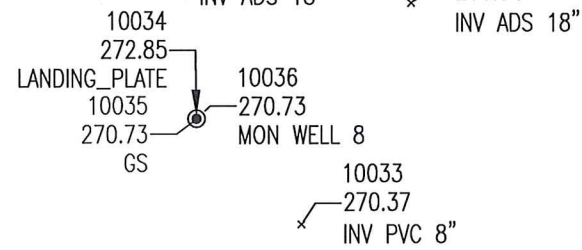
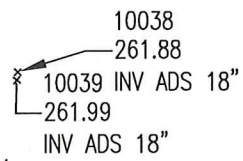
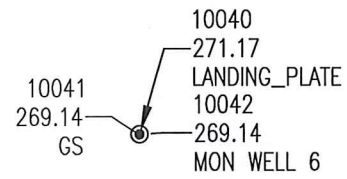
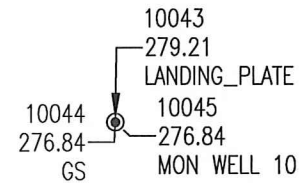
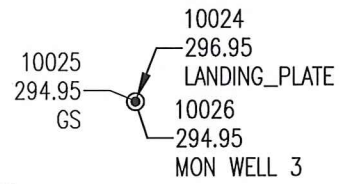
Appendix E

Survey Data

10003
342.53
TOP



DATUM:
HORIZONTAL: PROJECT
(PER 1997 SURVEY BY PMX)
VERTICAL: PROJECT
CONTROL POINT #1578-25
ELEV. 333.43



OLALLA LANDFILL SURVEY DATA.xls

| ID # | PMX # | NORTHING | EASTING | ELEVATION | DESCRIPTION | ELEV. | GROUND | ELEV. | DESC. |
|--------|-------|------------|------------|-----------|---|--------|------------|--------|---|
| F-1 | 10010 | 161775.469 | 560218.662 | 332.40 | FLARE 1 | 332.40 | GS FLARE 1 | | |
| F-2 | 10013 | 161922.669 | 560207.322 | 330.27 | FLARE 2 | 330.27 | GS FLARE 2 | | |
| F-3 | 10009 | 161707.727 | 559732.444 | 323.76 | FLARE 3 | 323.76 | GS FLARE 3 | | |
| MW-01 | 10058 | 161858.133 | 560525.840 | 342.53 | MONITORING WELL 1 | 342.53 | GS MW 1 | 343.79 | LANDING PLATE MW 1 |
| MW-02 | 10055 | 161704.534 | 559572.839 | 318.95 | MONITORING WELL 2 | 318.95 | GS MW 2 | 323.25 | PVC TOP MW 2 |
| MW-03 | 10026 | 162333.903 | 559463.060 | 294.95 | MONITORING WELL 3 | 294.95 | GS MW 3 | 296.95 | LANDING PLATE MW 3 |
| MW-04 | 10006 | 161911.192 | 559787.735 | 317.30 | MONITORING WELL 4 | 317.35 | GS MW 4 | 320.93 | PVC TOP MW 4 |
| MW-05 | 10032 | 162510.115 | 559878.901 | 332.78 | MONITORING WELL 5 | 332.78 | GS MW 5 | 334.17 | PVC TOP MW 5 |
| MW-05A | 10029 | 162487.878 | 559875.742 | 331.43 | MONITORING WELL 5A | 331.44 | GS MW 5A | 332.53 | LANDING PLATE MW 5A |
| MW-06 | 10042 | 162077.699 | 559358.970 | 269.14 | MONITORING WELL 6 | 269.14 | GS MW 6 | 271.17 | LANDING PLATE MW 6 |
| MW-07 | 10051 | 161723.016 | 559398.979 | 278.21 | MONITORING WELL 7 | 278.20 | GS MW 7 | 280.43 | LANDING PLATE MW 7 |
| MW-08 | 10036 | 161897.813 | 559350.147 | 270.73 | MONITORING WELL 8 | 270.73 | GS MW 8 | 272.85 | LANDING PLATE MW 8 |
| MW-10 | 10045 | 162218.490 | 559340.899 | 276.84 | MONITORING WELL 10 | 276.84 | GS MW 10 | 279.21 | LANDING PLATE MW 10 |
| STM-C1 | 10037 | 161964.884 | 559475.019 | 266.54 | INVERT 18" ADS - POND INLET | | | | |
| STM-C2 | 10038 | 162010.704 | 559336.808 | 261.88 | INVERT 18" ADS - OUTFALL FROM POND | | | | |
| STM-C3 | 10039 | 162006.004 | 559336.340 | 261.99 | INVERT 18" ADS - OUTFALL FROM POND | | | | |
| STM-C4 | 10033 | 161836.404 | 559411.172 | 270.37 | INVERT 8" PVC CULVERT - 90' SE. OF MW-7 | | | | |
| TP-1 | 10017 | 162289.111 | 559988.875 | 321.87 | TEST PIT #1 - NORTH PIN FLAG | | | | |
| TP-1 | 10016 | 162259.235 | 559993.205 | 322.39 | TEST PIT #1 - SOUTH PIN FLAG | | | | |
| TP-2 | 10015 | 162135.942 | 560120.834 | 325.20 | TEST PIT #2 - NORTH PIN FLAG | | | | |
| TP-2 | 10014 | 162107.804 | 560128.212 | 324.72 | TEST PIT #2 - SOUTH PIN FLAG | | | | |
| TP-3 | 10018 | 162123.185 | 559944.476 | 319.53 | TEST PIT #3 - NORTH PIN FLAG | | | | |
| TP-3 | 10019 | 162090.903 | 559947.172 | 319.26 | TEST PIT #3 - SOUTH PIN FLAG | | | | |
| TP-4 | 10022 | 162118.139 | 559675.923 | 310.22 | TEST PIT #4 - NORTH PIN FLAG | | | | |
| TP-4 | 10023 | 162065.499 | 559677.316 | 309.76 | TEST PIT #4 - SOUTH PIN FLAG | | | | |
| TP-5 | 10021 | 162201.740 | 559810.304 | 314.31 | TEST PIT #5 - NORTH PIN FLAG | | | | |
| TP-5 | 10020 | 162173.560 | 559815.981 | 314.43 | TEST PIT #5 - SOUTH PIN FLAG | | | | |
| WELL | 10003 | 162761.407 | 559934.278 | 342.53 | TOP CENTER OF WELL CASING - 260' N. OF MW-5 | | | 341.34 | SCRIBE 'X' CONC. FOUNDATION OF WELL |
| | | | | | | | | 342.31 | SCRIBED 'X' NE. COR. OF POWER J-BOX AT WELL |

Appendix F
Field Data Sheets

Parametrix, Inc.

Well #: MW-1
 Sample #: OK-MW-

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|--|
| Project Number | <u>215-1578-119</u> | Date | <u>6/1/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde / M. Baxter / J. Bennett</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde / Same</u> |

Casing Diameter: 2" 4" 6" Other

Depth to Water (feet) 25.90 Purge Vol. Meas. Method Grad Cyl/Stop Watch
 Depth of Well (feet) N/A Date Purged 6/1/2011
 Reference Point (surveyors notch/etc) TOC Purge Time (from/to) 1525 - 1602
 Date/Time Sampled 6/1/2011 1555 Flow = 460 ml/min

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 4

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|--------------|----------------|-------------|--------------------|
| <u>1525</u> | <u>26.71*</u> | <u>6.85</u> | <u>16.3</u> | <u>9.31</u> | <u>10.02</u> | <u>8.4</u> | <u>201</u> | |
| <u>1532</u> | <u>26.71</u> | <u>6.10</u> | <u>16.9</u> | <u>9.77</u> | <u>10.55</u> | <u>15.7</u> | <u>197</u> | |
| <u>1536</u> | <u>26.75</u> | <u>5.96</u> | <u>16.9</u> | <u>9.94</u> | <u>12.20</u> | <u>4.6</u> | <u>190</u> | |
| <u>1540</u> | <u>26.75</u> | <u>6.16</u> | <u>16.8</u> | <u>10.01</u> | <u>12.80</u> | <u>3.0</u> | <u>178</u> | |
| <u>1544</u> | <u>26.57</u> | <u>6.33</u> | <u>16.8</u> | <u>9.93</u> | <u>12.92</u> | <u>5.9</u> | <u>170</u> | |
| <u>1548</u> | <u>26.69</u> | <u>6.37</u> | <u>16.8</u> | <u>9.83</u> | <u>12.72</u> | <u>6.6</u> | <u>169</u> | |
| <u>1552</u> | <u>26.53</u> | <u>6.39</u> | <u>16.7</u> | <u>9.67</u> | <u>12.58</u> | <u>8.4</u> | <u>162</u> | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>6/2/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity hooked, clear of debris
 Remarks kink in tubing about 1.5' from top of casing.
 Signature _____ Page 1 of 1

* convenience depth

Parametrix, Inc.

Well #: MW-2
Sample #: Q-MW-2

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|--------------------------------------|
| Project Number | <u>215-1578-119</u> | Date | <u>6/1/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/ M Bostar/ J Bennett</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde</u> |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>63.20'</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | <u>N/A</u> | Date Purged | <u>6/1/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>0930 - 1015</u> |
| Date/Time Sampled | <u>6/1/2011 1010</u> | Flow = | <u>400 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 5

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|--------------|----------------|-------------|--------------------|
| <u>0940</u> | <u>63.35</u> | <u>5.69</u> | <u>14.2</u> | <u>8.51</u> | <u>11.41</u> | <u>1.4</u> | <u>213</u> | |
| <u>0944</u> | <u>63.33</u> | <u>6.09</u> | <u>13.7</u> | <u>8.53</u> | <u>13.96</u> | <u>1.2</u> | <u>182</u> | |
| <u>0948</u> | <u>63.34</u> | <u>6.35</u> | <u>13.6</u> | <u>8.61</u> | <u>14.13</u> | <u>3.3</u> | <u>169</u> | |
| <u>0952</u> | <u>63.35</u> | <u>6.52</u> | <u>13.5</u> | <u>8.61</u> | <u>14.05</u> | <u>5.6</u> | <u>166</u> | |
| <u>0956</u> | <u>—</u> | <u>6.55</u> | <u>13.4</u> | <u>8.50</u> | <u>14.04</u> | <u>8.7</u> | <u>165</u> | |
| <u>1000</u> | <u>—</u> | <u>6.67</u> | <u>13.4</u> | <u>8.56</u> | <u>14.01</u> | <u>13.9</u> | <u>163</u> | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>6/6/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity packed, clear of debris

Remarks Soft pump at #1 - water level indicator stopped working correctly after 0952 reading.

Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-3
 Sample #: 01-MW-3

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|--|
| Project Number | <u>215-1578-119</u> | Date | <u>6/2/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde / M. Baxter / C. Brito</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde / Same</u> |

Casing Diameter: 2" X 4" _____ 6" _____ Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>43.19</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | <u>N/A</u> | Date Purged | <u>6/2/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1133 - 1217</u> |
| Date/Time Sampled | <u>6/2/2011 1200</u> | Flow = | <u>320 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 3.5

| TIME (2400 hr) | WATER LEVEL (feet) \downarrow | pH (units) ± 0.1 | COND (mS/cm) $\pm 3\%$ | DO (mg/L) $\pm 10\%$ | TEMP $^{\circ}\text{C}$ | TURB. $\pm 10\%$ | ORP (mV) | CUM. VOL. (gal) |
|-------------------|---------------------------------------|----------------------------|------------------------------|----------------------------|----------------------------|---------------------|-------------|--------------------|
| <u>1136</u> | <u>43.84</u> | <u>6.78</u> | <u>24.3</u> | <u>0.22</u> | <u>10.51</u> | <u>0.0</u> | <u>189</u> | |
| <u>1140</u> | <u>43.80</u> | <u>6.28</u> | <u>23.2</u> | <u>0.00</u> | <u>11.20</u> | <u>0.3</u> | <u>195</u> | |
| <u>1144</u> | <u>43.80</u> | <u>6.22</u> | <u>23.3</u> | <u>0.00</u> | <u>12.15</u> | <u>0.0</u> | <u>191</u> | |
| <u>1148</u> | <u>43.80</u> | <u>6.22</u> | <u>23.4</u> | <u>0.00</u> | <u>12.69</u> | <u>0.0</u> | <u>187</u> | |
| <u>1152</u> | <u>-</u> | <u>6.21</u> | <u>23.4</u> | <u>0.00</u> | <u>12.89</u> | <u>0.0</u> | <u>186</u> | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>6/2/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>01-MW-01</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

| | | | |
|----------------|---|------|----------------------|
| Well Integrity | <u>locked, clear of debris.</u> | | |
| Remarks | <u>collected field duplicate, 01-MW-01. (set time @ 1245)</u> | | |
| Signature | _____ | Page | <u>1</u> of <u>1</u> |

* Convenience depth

Parametrix, Inc.

Well #: MW-4
 Sample #: OL-MW-4

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|---|
| Project Number | <u>215-1578-119</u> | Date | <u>6/1/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/ M. Baxter / J. Bennett</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/ Sumic</u> |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>60.25</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | <u>N/A</u> | Date Purged | <u>6/1/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1110-1157</u> |
| Date/Time Sampled | <u>6/1/2011 1145</u> | Flow = | <u>320 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 4

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|--------------|----------------|-------------|--------------------|
| <u>1115</u> | <u>60.51</u> | <u>6.87</u> | <u>15.1</u> | <u>8.99</u> | <u>11.48</u> | <u>2.9</u> | <u>159</u> | |
| <u>1119</u> | <u>60.50</u> | <u>6.86</u> | <u>15.1</u> | <u>8.86</u> | <u>13.24</u> | <u>3.8</u> | <u>165</u> | |
| <u>1123</u> | <u>60.31</u> | <u>6.97</u> | <u>19.0</u> | <u>8.84</u> | <u>13.98</u> | <u>3.3</u> | <u>165</u> | |
| <u>1127</u> | <u>60.32</u> | <u>7.01</u> | <u>15.2</u> | <u>8.77</u> | <u>14.91</u> | <u>9.6</u> | <u>144</u> | |
| <u>1131</u> | <u>60.52</u> | <u>7.09</u> | <u>19.9</u> | <u>9.09</u> | <u>15.60</u> | <u>6.1</u> | <u>136</u> | |
| <u>1135</u> | <u>-</u> | <u>7.07</u> | <u>15.0</u> | <u>9.04</u> | <u>14.60</u> | <u>4.9</u> | <u>142</u> | |
| <u>1139</u> | <u>-</u> | <u>7.00</u> | <u>15.0</u> | <u>8.92</u> | <u>14.16</u> | <u>10.5</u> | <u>147</u> | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>6/1/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Caps locked, lots of little black ants!
 Remarks Sampling not done
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: WA
 Sample #: 06-MW-WA

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|--|
| Project Number | <u>215-1578-119</u> | Date | <u>6/2/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Lindet - M Baxter / C Burtrigo</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Lindet same</u> |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>74.10</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | <u>N/A</u> | Date Purged | <u>6/2/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>0900 - 1005</u> |
| Date/Time Sampled | <u>6/2/2011 0940</u> | Flow = | <u>360 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 5

| TIME (2400 hr) | WATER LEVEL (feet) * | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|----------------|----------------------|------------------|-------------------|-----------------|--------------|-------------|------------|-----------------|
| 0905 | <u>75.87</u> | <u>6.67</u> | <u>15.7</u> | <u>9.22</u> | <u>11.58</u> | <u>4.1</u> | <u>204</u> | |
| 0909 | <u>75.87</u> | <u>6.81</u> | <u>15.2</u> | <u>9.12</u> | <u>11.65</u> | <u>1.2</u> | <u>181</u> | |
| 0913 | <u>76.10</u> | <u>6.00</u> | <u>14.9</u> | <u>9.01</u> | <u>12.57</u> | <u>0.0</u> | <u>159</u> | |
| 0917 | <u>76.76</u> | <u>6.32</u> | <u>14.7</u> | <u>8.84</u> | <u>13.46</u> | <u>0.0</u> | <u>148</u> | |
| 0921 | <u>75.90</u> | <u>6.49</u> | <u>14.6</u> | <u>9.12</u> | <u>14.00</u> | <u>0.0</u> | <u>140</u> | |
| 0925 | <u>75.90</u> | <u>6.51</u> | <u>14.6</u> | <u>9.00</u> | <u>14.32</u> | <u>0.0</u> | <u>140</u> | |
| 0929 | <u>75.85</u> | <u>6.58</u> | <u>14.7</u> | <u>9.23</u> | <u>14.59</u> | <u>1.5</u> | <u>140</u> | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>6/3/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity checked, clear of debris. lots of small bugs!
 Remarks None
 Signature [Signature] Page 1 of 1

* Convenience depth

Parametrix, Inc.

Well #: MW-6
 Sample #: Oh-MW-6

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|---|
| Project Number | <u>215-1578-119</u> | Date | <u>6/2/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/ M. Burtner/ C. Burtner</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/ SWMC</u> |

Casing Diameter: 2" X 4" 6" Other

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>19.16</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | <u>N/A</u> | Date Purged | <u>6/2/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1310 - 1355</u> |
| Date/Time Sampled | <u>6/2/2011 1355</u> | Flow = | <u>350 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) Actual Purge Volume (gallons) 5

| TIME (2400 hr) | WATER LEVEL (feet) * | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ±10% | TEMP °C | TURB. ±10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|----------------------------|------------------------|-------------------------|----------------------|------------|---------------|-------------|--------------------|
| 1315 | 19.70 | 6.21 | 12.0 | 2.35 | 11.72 | 103.0 | 51 | |
| 1319 | 19.76 | 5.98 | 12.5 | 0.88 | 12.70 | 26.1 | 48 | |
| 1323 | | 6.01 | 12.9 | 0.43 | 12.94 | 26.4 | 46 | |
| 1327 | 19.60 | 6.15 | 13.4 | 0.25 | 13.00 | 32.7 | 38 | |
| 1331 | 19.60 | 6.37 | 13.7 | 0.10 | 13.00 | 34.3 | 24 | |
| 1335 | 19.64 | 6.53 | 13.8 | 0.03 | 13.04 | 6.0 | 16 | |
| 1339 | 19.58 | 6.61 | 13.9 | 0.00 | 13.10 | 29.0 | 10 | |
| 1343 | 19.58 | 6.61 | 14.0 | 0.00 | 13.60 | 41.7 | 7 | |
| 1347 | 19.58 | 6.67 | 14.1 | 0.00 | 12.96 | 55.8 | 6 | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>6/3/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

| | |
|----------------|--|
| Well Integrity | <u>locked, clear of debris.</u> |
| Remarks | <u>sand visible in outlet tube causing turbidity to jump around.</u> |
| Signature | <u> </u> Page <u>1</u> of <u>1</u> |

* Convenience depth.

Parametrix, Inc.

Well #: MW-7
 Sample #: OL-MW-7

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|--|
| Project Number | <u>215-1578-119</u> | Date | <u>6/1/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/ M. Bunker/ J. Bennett</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/ M. Bunker/ J. Bennett</u> |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>23.65</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | <u>N/A</u> | Date Purged | <u>6/1/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1342 - 1422</u> |
| Date/Time Sampled | <u>6/1/2011 1420</u> | Flow = | <u>320 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 3.0

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|--------------|----------------|-------------|--------------------|
| <u>1345</u> | <u>23.74</u> | <u>7.06</u> | <u>17.0</u> | <u>4.58</u> | <u>10.49</u> | <u>9.1</u> | <u>147</u> | |
| <u>1349</u> | <u>23.72</u> | <u>6.41</u> | <u>16.9</u> | <u>4.13</u> | <u>11.44</u> | <u>4.6</u> | <u>153</u> | |
| <u>1353</u> | <u>-</u> | <u>6.45</u> | <u>16.0</u> | <u>4.10</u> | <u>11.85</u> | <u>4.4</u> | <u>148</u> | |
| <u>1357</u> | <u>23.71</u> | <u>6.61</u> | <u>17.0</u> | <u>4.01</u> | <u>12.12</u> | <u>5.2</u> | <u>144</u> | |
| <u>1401</u> | <u>23.69</u> | <u>6.78</u> | <u>16.9</u> | <u>3.97</u> | <u>12.37</u> | <u>9.0</u> | <u>134</u> | |
| <u>1405</u> | <u>23.70</u> | <u>6.85</u> | <u>16.9</u> | <u>3.77</u> | <u>12.64</u> | <u>13.1</u> | <u>132</u> | |
| <u>1409</u> | <u>23.70</u> | <u>6.88</u> | <u>16.8</u> | <u>3.94</u> | <u>12.45</u> | <u>19.0</u> | <u>118</u> | |
| <u>1413</u> | <u>23.70</u> | <u>6.90</u> | <u>16.8</u> | <u>3.89</u> | <u>12.70</u> | <u>22.3</u> | <u>119</u> | |
| <u>1417</u> | <u>-</u> | <u>6.92</u> | <u>16.8</u> | <u>3.86</u> | <u>12.81</u> | <u>26.3</u> | <u>122</u> | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>6/1/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

| | | | |
|----------------|--------------------------------|------|----------------------|
| Well Integrity | <u>locked, clear of debris</u> | | |
| Remarks | <u>None</u> | | |
| Signature | <u>[Signature]</u> | Page | <u>1</u> of <u>1</u> |

Parametrix, Inc.

Well #: MW-8
 Sample #: OL-MW-8

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|--|
| Project Number | <u>215-1578-119</u> | Date | <u>6/2/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde / M Baxter / C Brittrago</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde / Same</u> |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>19.65</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | <u>N/A</u> | Date Purged | <u>6/2/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1413 - 1510</u> |
| Date/Time Sampled | <u>6/2/2011 1500</u> | Flow = | <u>400 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 5

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ±10% | TEMP °C | TURB. ±10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|----------------------|------------|---------------|-------------|--------------------|
| 1415 | 20.60 | 6.37 | 60.5 | 1.15 | 10.06 | 96.9 | 20 | |
| 1422 | 20.58 | 6.13 | 59.1 | 0.00 | 10.82 | 65.3 | 28 | |
| 1426 | 20.59 | 6.19 | 55.2 | 0.00 | 11.11 | 57.1 | 31 | |
| 1430 | 20.58 | 6.31 | 50.4 | 0.00 | 11.24 | 60.3 | 31 | |
| 1434 | 20.59 | 6.49 | 46.4 | 0.00 | 11.33 | 48.9 | 26 | |
| 1438 | 20.59 | 6.61 | 44.6 | 0.05 | 11.34 | 45.5 | 24 | |
| 1442 | - | 6.66 | 42.2 | 0.13 | 11.42 | 42.8 | 24 | |
| 1446 | 20.59 | 6.68 | 41.7 | 0.19 | 11.39 | 41.1 | 23 | |
| 1450 | 20.59 | 6.68 | 40.6 | 0.24 | 11.41 | 41.6 | 25 | |
| 1454 | 20.59 | 6.68 | 39.1 | 0.29 | 11.42 | 42.2 | 26 | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>6/3/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity locked, clear of debris.
 Remarks None
 Signature [Signature] Page 1 of 1

* Convenience depth.

Parametrix, Inc.

Groundwater Sampling Field Data Sheet

MW-10
 Well #: ~~MW-3~~ MW-3
 Sample #: ~~OK-MW-3~~ OK-MW-3
 OK-MW-10

| | | | |
|-----------------|-----------------------------|------------|-----------------------------------|
| Project Number | 215-1578-119 | Date | 6/2/2011 |
| Project Name | Olalla LF Qtrly GW Sampling | Location | Olalla LF |
| Project Address | Bandix Rd | Sampled By | L. Linde/ M. Baxter / C. Buttrick |
| Client Name | KC SWD | Purged By | L. Linde/ same |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|---------------|-------------------------|---------------------|
| Depth to Water (feet) | 20.18 | Purge Vol. Meas. Method | Grad Cyl/Stop Watch |
| Depth of Well (feet) | N/A | Date Purged | 6/2/2011 |
| Reference Point (surveyors notch/etc) | TOC | Purge Time (from/to) | 1029 - 1110 |
| Date/Time Sampled | 6/2/2011 1105 | Flow | = 400 ml/min |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 15

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|------------|----------------|-------------|--------------------|
| 1032 | 29.02* | 6.30 | 76.6 | 0.00 | 10.40 | 19.9 | 205 | |
| 1036 | 29.04 | 6.35 | 76.7 | 0.00 | 11.40 | 0.0 | 206 | |
| 1040 | 29.04 | 6.04 | 76.4 | 0.00 | 11.84 | 0.0 | 194 | |
| 1044 | 29.06 | 6.23 | 76.9 | 0.00 | 11.94 | 0.0 | 179 | |
| 1048 | 29.08 | 6.39 | 76.0 | 0.00 | 11.97 | 0.0 | 168 | |
| 1052 | 29.08 | 6.47 | 76.0 | 0.00 | 11.93 | 0.0 | 162 | |
| 1056 | — | 6.49 | 76.9 | 0.00 | 11.94 | 0.0 | 159 | |
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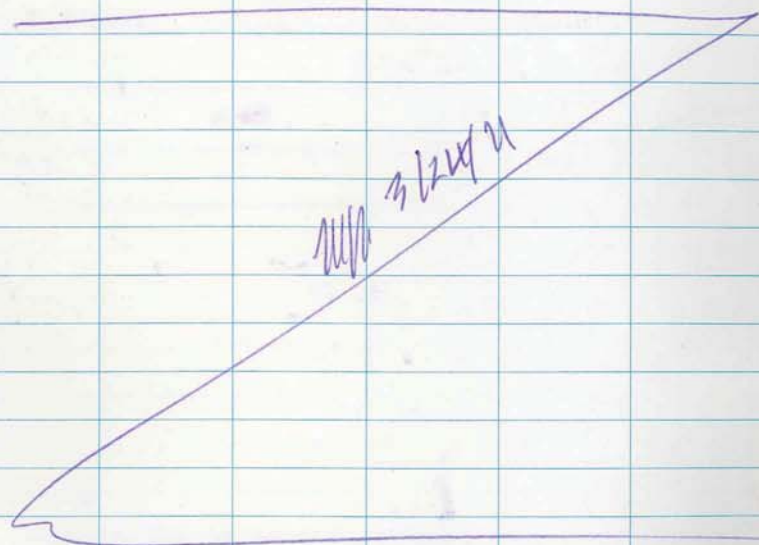
| | | | |
|---------------------------|----------------------|-----------------------------------|----------------------|
| Purge Equipment | Grundfos Ready-flo 2 | Sampling Equipment | Grundfos Ready-flo 2 |
| Laboratory | Aquatic Research | Date Sent to Lab | 6/3/2011 |
| Chain-of-Custody (yes/no) | Yes | Field QC Sample Number | N/A |
| Shipment Method | Hand Delivered | Split with (name(s)/organization) | N/A |

Well Integrity checked, r. clear of debris.
 Remarks None
 Signature [Signature] Page 1 of 1

* Convenience depth.

Thursday, March 24, 2011

- 1335 Collect MW-10, plus M6/M6D
- 1420 End purge
- 1435 Depart MW-10
- 1440 Arrive at MW-3
- 1445 Start purge
- 1430 Collect MW-3 sample
- 1540 End purge
- 1545 Depart MW-3
- 1550 Transfer samples and equipment to Jesse's truck, complete COC
- 1600 Depart project site, lock gate behind us.



Wednesday - June 1, 2011

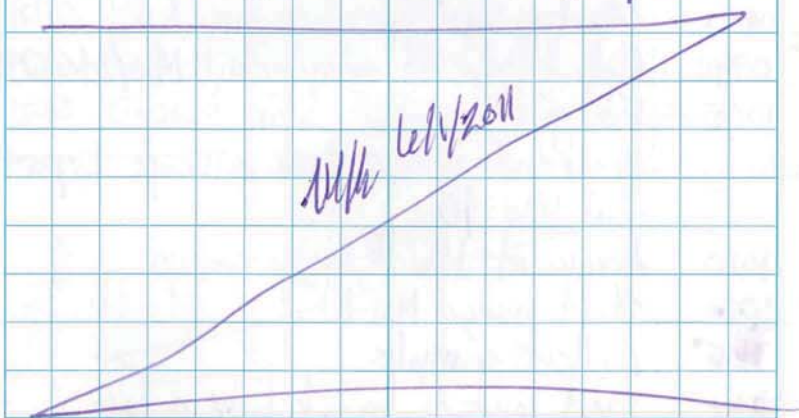
Samplers - Jesse Bennett + Mike Baxter
 Forecast - Hi 58° to 48° Rain showers
 Current - 52° Overcast + sprinkles

- 0800 Meet with J Bennett and L Hinde at Bandix Rd gate. Receive field journal and tubing for MW-2 + MW-4 from Kara. Review operation of gas meter.
 - 0815 L Hinde departs. Unlock gate and Jesse and Mike enter. Close gate behind us (not locked).
 - 0825 Plan for day and health + safety meeting.
 - 0840 Depart for MW-2.
 - 0845 Arrive at staging area on road to get to MW-2 and MW-4. Pre-label sample containers for wells and calibrate fluorometer.
 - 0905 Begin walking equipment out and setting up at MW-2
 - 0920 Measure MW-2 DTW (63.20')
 - 0930 Begin purging MW-2
- (over)

Wednesday, June 1, 2011

- 0915 Water level indicator stopped working correctly. Would work when test button was pressed, but would not buzz to indicate water level.
- 1010 Collect MW-2 sample.
- 1015 End purge. Breakdown setup and begin decon. Also start moving equipment to MW-4.
- 1055 Finish decon. Depart MW-2 for MW-4.
- 1045 Arrive at MW-4 and begin setup.
- 1050 Measure MW-4 DTW (60.25)
- 1105 Visit from Charlie (?) from KCSW
- 1110 Begin purging MW-4.
- 1145 Collect MW-4
- 1157 End purge. Breakdown setup. Move pumping equipment off caps to where we can pull vehicles around to pick up samples.
- 1125 Decon pump. Put away MW-2 and MW-4 samples.
- 1235 Break for lunch.
- 1315 Arrive at MW-7 start set up.
- 1320 Measure MW-7 DTW (25.65)

- 1343 Start purging MW-7
- 1420 Collect MW-7 sample
- 1427 End purge. Breakdown setup.
- 1435 Depart MW-7 for gas monitoring.
- 1445 Measure Flare 3
- 1455 Measure Flare 1
- 1505 Measure Flare 2, depart for MW-1
- 1510 Arrive at MW-1, begin setup.
- 1515 Measure DTW (75.90)
- 1525 Begin purge
- 1555 Collect MW-1 sample
- 1602 End purge, breakdown setup.
- 1615 Collect rinse blank (Oh-EK)
- 1635 Secure gate on Bandia Rd, depart site.



Thursday - June 2, 2011

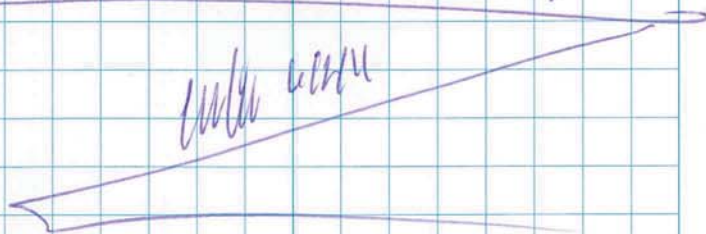
- 0803 Meet up with Craig Bairrigo
 0810 Unlock gate on Bandix Rd to access site, lock gate behind us.
 0815 Transfer gear, etc, to Highlander. Depart for MW-6A.
 0820 Arrive at MW-6A, Conduct H₂S meeting, label sample containers.

Samplers: Mike Buxton and Craig Bairrigo
 Forecast: Hi 67° to 49° Rain Showers
 Current: 64° Mostly Cloudy, Slight Breeze

- 0840 Calibrate Horiba WQ Meter
 0850 setup for purge.
 0900 start purge
 0940 Collect MW-6A sample and MS/MSD
 1005 End purge.
 1010 Breakdown setup, lock well cap. Depart for MW-10.
 1020 Arrive at MW-10, begin setup.
 1029 start purge MW-10
 1055 Collect sample
 1110 End purge, breakdown setup

Thursday - June 2, 2011

- 1120 Depart MW-10 for MW-3
 1124 Arrive at MW-3 and setup for purge
 1133 start purge.
 1200 Collect MW-3 and field duplicate (MW-9)
 1217 End purge, breakdown setup
 1225 Break for lunch.
 1300 Arrive at MW-6 and setup for purge
 1310 Begin purge
 1355 Collect MW-6 sample
 1358 End purge, breakdown setup.
 1405 Depart for MW-8
 1408 Arrive at MW-8, set up for purge
 1413 start Purge
 1500 Collect MW-8 sample
 1510 End purge, breakdown setup
 1520 Depart MW-8 for Bandix Rd gate.
 1530 Depart site, lock gate behind us.





Kitsap County Public Works
Solid Waste Division
 614 Division Street (MS-27)
 Port Orchard, WA 98366
 Phone: 360-337-5777
 FAX: 360-337-4867

CHAIN OF CUSTODY RECORD

FOR LAB USE ONLY

| | |
|-----------------|---|
| LOGIN COMMENTS: | Sample(s) checked/logged by: _____ |
| _____ | Sample temp on receipt at lab: _____ COMPLIANCE 4 +/- 4° C |
| _____ | Are custody seals intact? <input type="checkbox"/> YES <input type="checkbox"/> NO |
| _____ | Sample receipt (day/time): _____ / _____ |
| _____ | Ice? <input type="checkbox"/> NONE <input type="checkbox"/> BLUE <input type="checkbox"/> WET |

TO BE COMPLETED BY THE SAMPLER:

| | | |
|---|---------------------------|--|
| TAT REQUESTED: STANDARD: <input checked="" type="checkbox"/> OTHER: _____ | PROJECT: <u>Olalla WF</u> | SAMPLER PRINTED NAME AND SIGNATURE <u>Mike Baxter / Jesse Bennett</u> |
|---|---------------------------|--|

| SAMPLE DATE | SAMPLE TIME | SITE NAME OR LOCATION | SAMPLE IDENTIFIER | MATRIX * | GRAB | COMP | ANALYSES REQUIRED—MARK AN X IN ALL TESTS REQUIRED FOR EACH SAMPLE LINE | | | | | | | | | | | | | | | |
|-------------|-------------|-----------------------|-------------------|----------|------|------|--|-----------|-----|---|----------------------|---------|-------------|--|---------|--|--|--|--|--|--|-----------------|
| | | | | | | | VOC Full list + SEM-VC | ALK-T,B,C | NH3 | Cl, NO ₂ , NO ₃ , SO ₄ | CO ₂ /TAC | pH, TDS | T. Coliform | Pres. Ag. Bact. Cont. E. Coli, W. Coli | Na, Zn. | | | | | | | |
| 6/1 | 1010 | | OL-MW-2 | GW | X | | | | | | | | | | | | | | | | | Pres. Metalsure |
| | 1145 | | OL-MW-4 | | | | | | | | | | | | | | | | | | | Field Filtered |
| | 1420 | | OL-MW-7 | | | | | | | | | | | | | | | | | | | |
| | 1605 | | OL-MW-1 | | | | | | | | | | | | | | | | | | | |
| | 1615 | | OL-ER | DIW | | | | | | | | | | | | | | | | | | Equip. Rinse |
| 6/24 | 0000 | | OL-TB | DIW | | | | | | | | | | | | | | | | | | Vocstrip Blank |
| | | | | | | | | | | | | | | | | | | | | | | |
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* MATRIX TYPES: SW = Surface Water GW = Ground Water SO = Soil

| SIGNATURE | PRINTED NAME | COMPANY AND TITLE | DATE | TIME |
|-------------------------------------|------------------|-------------------------|----------|------|
| RELINQUISHED BY: <u>[Signature]</u> | Mike Baxter | Parametrix / Env Tech I | 6/1/2011 | 1900 |
| RECEIVED BY: <u>[Signature]</u> | Michael R. Wafel | Parametrix / Project | 6/1/2011 | 1900 |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |



Kitsap County Public Works
Solid Waste Division
 614 Division Street (MS-27)
 Port Orchard, WA 98366
 Phone: 360-337-5777
 FAX: 360-337-4867

CHAIN OF CUSTODY RECORD

FOR LAB USE ONLY

| | |
|-----------------|---|
| LOGIN COMMENTS: | Sample(s) checked/logged by: _____ |
| | Sample temp on receipt at lab: _____ COMPLIANCE 4 +/- 4° C |
| | Are custody seals intact? <input type="checkbox"/> YES <input type="checkbox"/> NO |
| | Sample receipt (day/time): _____ / _____ |
| | Ice? <input type="checkbox"/> NONE <input type="checkbox"/> BLUE <input type="checkbox"/> WET |

TO BE COMPLETED BY THE SAMPLER:

| | | |
|---|---------------------------|--|
| TAT REQUESTED: STANDARD: <input checked="" type="checkbox"/> OTHER: _____ | PROJECT: <u>Oralla LE</u> | SAMPLER PRINTED NAME AND SIGNATURE <u>Mike Baxter / Craig Brittrago</u> |
|---|---------------------------|--|

| SAMPLE DATE | SAMPLE TIME | SITE NAME OR LOCATION | SAMPLE IDENTIFIER | MATRIX * | GRAB | COMP | ANALYSES REQUIRED—MARK AN X IN ALL TESTS REQUIRED FOR EACH SAMPLE LINE | | | | | | | | | | | |
|-------------|-------------|-----------------------|-------------------|----------|------|------|--|----------|-----------------|---|----------------------|--------|-------------|--------------------------------------|---|---|--|--|
| | | | | | | | VE Full list + SIM-VC | Alk-TB,C | NH ₃ | Cl, NO ₂ , NO ₃ , SO ₄ | CO ₂ /Toc | pH/TDS | T. Coliform | Dist. As, Ba, Ca, Fe, Mn, Ni, Na, Zn | | | | |
| 4/2 | 0940 | | Oh-MW-5A | GW | X | | X | X | X | X | X | X | X | X | X | X | MS/MEO set | |
| | 1105 | | Oh-MW-10 | | | | | | | | | | | | | | Dist. Metals samples have been field filtered. | |
| | 1200 | | Oh-MW-3 | | | | | | | | | | | | | | | |
| | 1245 | | Oh-MW-9 | | | | | | | | | | | | | | | |
| | 1355 | | Oh-MW-6 | | | | | | | | | | | | | | | |
| | 1500 | | Oh-MW-8 | | | | | | | | | | | | | | | |

* MATRIX TYPES: SW = Surface Water GW = Ground Water SO = Soil

| SIGNATURE | PRINTED NAME | COMPANY AND TITLE | DATE | TIME |
|--------------------|-------------------|-------------------------|----------|------|
| <i>[Signature]</i> | Mike Baxter | Parametrix / Env Tech I | 4/2/2011 | 1830 |
| <i>[Signature]</i> | Michael R. Warfel | Parametrix / Proj Mgr | 6/2/2011 | 1830 |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |

Parametrix, Inc.

Well #: MW-1
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-1

| | | | |
|-----------------|--|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>3/23/11</u> |
| Project Name | <u>Olalla LF Qtrly GW/RIFSSampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Lind / M. Baxter</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>"</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|---------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>78.12</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | <u>-</u> | Date Purged | <u>3/23/11</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>0941-1010</u> |
| Date/Time Sampled | <u>3/23/11 1010</u> | | <u>450 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|------------|-------------------------------|------------------------|--------------------|
| 0941 | 78.38 | 5.44 | 0.231 | 10.36 | 9.94 | < 1 | 186 | |
| 0945 | 78.37 | 5.64 | 0.232 | 10.38 | 10.47 | < 1 | 179 | |
| 0949 | 78.37 | 5.70 | 0.233 | 10.55 | 12.20 | < 1 | 155 | |
| 0953 | 78.38 | 5.78 | 0.234 | 10.87 | 13.80 | < 1 | 140 | |
| 0957 | 78.38 | 5.86 | 0.234 | 10.95 | 13.31 | < 1 | 137 | |
| 1001 | 78.38 | 5.92 | 0.234 | 10.87 | 13.55 | < 1 | 137 | |
| 1005 | 78.38 | 5.95 | 0.234 | 10.99 | 13.66 | < 1 | 137 | 4.5 gal |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>3/23/11</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good

Remarks _____

Signature Lind Page 1 of 1

Parametrix, Inc.

Well #: MW-2
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-2

| | | | |
|-----------------|--|------------|-----------------------------|
| Project Number | 215-1578-121 | Date | <u>3/23/11</u> |
| Project Name | <u>Olalla LF Qtrly GW/RIFSSampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde / M. Baxter</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>11</u> |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|---------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>63.67</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | <u>-</u> | Date Purged | <u>3/23/11</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1336 - 1352</u> |
| Date/Time Sampled | <u>3/23/11 1400</u> | | <u>420 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|--------------|-------------------------------|------------------------|--------------------|
| <u>1336</u> | <u>63.77</u> | <u>6.83</u> | <u>0.185</u> | <u>9.46</u> | <u>13.62</u> | <u>< 1</u> | <u>229</u> | |
| <u>1342</u> | <u>63.79</u> | <u>6.73</u> | <u>0.193</u> | <u>9.80</u> | <u>13.61</u> | <u>< 1</u> | <u>98</u> | |
| <u>1344</u> | <u>63.81</u> | <u>6.71</u> | <u>0.195</u> | <u>10.29</u> | <u>13.79</u> | <u>< 1</u> | <u>72</u> | |
| <u>1346</u> | <u>63.80</u> | <u>6.72</u> | <u>0.196</u> | <u>10.46</u> | <u>13.58</u> | <u>< 1</u> | <u>72</u> | |
| <u>1352</u> | <u>63.80</u> | <u>6.73</u> | <u>0.197</u> | <u>10.50</u> | <u>13.40</u> | <u>< 1</u> | <u>72</u> | <u>4 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>3/24/11</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-3
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-3

| | | | |
|-----------------|---------------------------------|------------|-----------------------------|
| Project Number | 215-1578-121 | Date | <u>3/24/2011</u> |
| Project Name | Olalla LF Qtrly GW/RIFSSampling | Location | Olalla LF |
| Project Address | Bandix Rd | Sampled By | <u>J Bennett / M Parker</u> |
| Client Name | KC SWD | Purged By | <u>same</u> |

Casing Diameter: 2" X 4" 6" Other _____

| | | | |
|---------------------------------------|-----------------------|-------------------------|----------------------|
| Depth to Water (feet) | <u>41.79</u> | Purge Vol. Meas. Method | Meas. Cup/Stop Watch |
| Depth of Well (feet) | _____ | Date Purged | <u>3/24/2011</u> |
| Reference Point (surveyors notch/etc) | TOC | Purge Time (from/to) | <u>1448 - 1540</u> |
| Date/Time Sampled | <u>3/24/2011 1530</u> | Flow = | <u>420 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 4.15

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|--------------|-------------------------------|------------------------|--------------------|
| <u>1451</u> | <u>42.31*</u> | <u>6.34</u> | <u>0.692</u> | <u>1.63</u> | <u>11.39</u> | <u>70.8</u> | <u>154</u> | |
| <u>1455</u> | <u>42.35</u> | <u>6.04</u> | <u>0.689</u> | <u>0.75</u> | <u>11.27</u> | <u>68.5</u> | <u>159</u> | |
| <u>1459</u> | <u>42.38</u> | <u>5.96</u> | <u>0.690</u> | <u>0.59</u> | <u>12.29</u> | <u>66.9</u> | <u>160</u> | |
| <u>1503</u> | <u>42.37</u> | <u>5.93</u> | <u>0.706</u> | <u>0.33</u> | <u>12.41</u> | <u>61.8</u> | <u>161</u> | |
| <u>1507</u> | <u>42.37</u> | <u>5.94</u> | <u>0.708</u> | <u>0.12</u> | <u>12.46</u> | <u>55.2</u> | <u>161</u> | |
| <u>1511</u> | <u>42.37</u> | <u>5.92</u> | <u>0.708</u> | <u>0.00</u> | <u>12.49</u> | <u>54.9</u> | <u>161</u> | |
| <u>1515</u> | — | <u>5.93</u> | <u>0.709</u> | <u>0.00</u> | <u>12.49</u> | <u>61.1</u> | <u>161</u> | |
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|---------------------------|----------------------|-----------------------------------|----------------------|
| Purge Equipment | Grundfos Ready-flo 2 | Sampling Equipment | Grundfos Ready-flo 2 |
| Laboratory | Aquatic Research | Date Sent to Lab | <u>3/24/2011</u> |
| Chain-of-Custody (yes/no) | Yes | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | Hand Delivered | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

* Convenience Depth

Parametrix, Inc.

Well #: mw-4
 Sample #: _____
OL-mw-4

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|---------------------------------|------------|-----------------------------|
| Project Number | 215-1578-121 | Date | <u>3/23/11</u> |
| Project Name | Olalla LF Qtrly GW/RIFSSampling | Location | Olalla LF |
| Project Address | Bandix Rd | Sampled By | <u>L. Lunde / M. Baxter</u> |
| Client Name | KC SWD | Purged By | <u>"</u> |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|---------------------|-------------------------|----------------------|
| Depth to Water (feet) | <u>61.00</u> | Purge Vol. Meas. Method | Meas. Cup/Stop Watch |
| Depth of Well (feet) | <u>-</u> | Date Purged | <u>3/23/11</u> |
| Reference Point (surveyors notch/etc) | TOC | Purge Time (from/to) | <u>1443 - 1510</u> |
| Date/Time Sampled | <u>3/23/11 1510</u> | | <u>300 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|------------|-------------------------------|-----------------------|--------------------|
| 1443 | 61.11 | 6.92 | 0.217 | 9.75 | 15.77 | < 1 | 233 | |
| 1447 | 61.08 | 6.85 | 0.231 | 9.27 | 13.45 | < 1 | 172 | |
| 1451 | 61.10 | 6.80 | 0.240 | 9.30 | 14.65 | < 1 | 117 | |
| 1455 | 61.09 | 6.78 | 0.243 | 9.43 | 14.51 | < 1 | 89 | |
| 1459 | 61.09 | 6.80 | 0.245 | 9.41 | 14.46 | < 1 | 82 | |
| 1503 | 61.09 | 6.81 | 0.244 | 9.45 | 14.46 | < 1 | 79 | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>3/24/11</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good

Remarks _____

Signature L. Lunde Page 1 of 1

Parametrix, Inc.

Well #: MW-5A
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-5A

| | | | |
|-----------------|--|------------|-----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>3/23/11</u> |
| Project Name | <u>Olalla LF Qtrly GW/RIFSSampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde / M. Baxter</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>"</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|---------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>75.14</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>3/23/11</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1047 - 1110</u> |
| Date/Time Sampled | <u>3/23/11 1110</u> | | <u>300 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|--------------|-------------------------------|------------------------|--------------------|
| <u>1047</u> | <u>76.78</u> | <u>6.36</u> | <u>0.295</u> | <u>10.78</u> | <u>11.48</u> | <u>< 1</u> | <u>138</u> | |
| <u>1051</u> | <u>76.80</u> | <u>6.39</u> | <u>0.305</u> | <u>11.00</u> | <u>11.56</u> | <u>< 1</u> | <u>117</u> | |
| <u>1055</u> | <u>76.77</u> | <u>6.43</u> | <u>0.288</u> | <u>11.25</u> | <u>12.50</u> | <u>< 1</u> | <u>98</u> | |
| <u>1100</u> | <u>76.78</u> | <u>6.47</u> | <u>0.280</u> | <u>11.44</u> | <u>13.37</u> | <u>< 1</u> | <u>95</u> | |
| <u>1104</u> | <u>76.75</u> | <u>6.48</u> | <u>0.276</u> | <u>11.61</u> | <u>13.42</u> | <u>< 1</u> | <u>92</u> | <u>3.5 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>3/24/11</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good

Remarks _____

Signature L. Linde Page 1 of 1

Parametrix, Inc.

Well #: MW-6
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-6

| | | | |
|-----------------|--|------------|-----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>3/24/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW/RIFSSampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>J Bennett / M Baxter</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>same</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|-----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>17.93</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | <u>-</u> | Date Purged | <u>3/24/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1102 - 1157</u> |
| Date/Time Sampled | <u>3/24/2011 1130</u> | Flow = | <u>250 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 4.5

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ±10% if >10NTU | ORP (mV) ±10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|--------------|----------------------------|----------------------|--------------------|
| <u>1105</u> | <u>18.00</u> | <u>6.49</u> | <u>0.294</u> | <u>0.19</u> | <u>12.93</u> | <u>42.0</u> | <u>-14</u> | |
| <u>1109</u> | <u>18.02</u> | <u>6.42</u> | <u>0.291</u> | <u>0.00</u> | <u>12.65</u> | <u>45.6</u> | <u>-23</u> | |
| <u>1113</u> | <u>18.01</u> | <u>6.42</u> | <u>0.289</u> | <u>0.00</u> | <u>12.72</u> | <u>43.4</u> | <u>-27</u> | |
| <u>1117</u> | <u>18.02</u> | <u>6.42</u> | <u>0.290</u> | <u>0.00</u> | <u>12.67</u> | <u>46.4</u> | <u>-31</u> | |
| <u>1121</u> | <u>18.01</u> | <u>6.44</u> | <u>0.291</u> | <u>0.00</u> | <u>12.67</u> | <u>37.1</u> | <u>-32</u> | |
| <u>1125</u> | | | | | | | | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>3/25/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good

Remarks _____

Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-7
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-7

| | | | |
|-----------------|--|------------|------------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>3/24/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW/RIFSSampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>J Bennett / M. Keator</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>same</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|-----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>23.49</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | <u>—</u> | Date Purged | <u>3/24/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>0816-0934</u> |
| Date/Time Sampled | <u>3/24/2011 0915</u> | Flow = | <u>390 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 4

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10NTU | ORP (mV) ± 10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|--------------|------------------------------|-----------------------|--------------------|
| <u>0842</u> | <u>24.49</u> | <u>5.71</u> | <u>0.232</u> | <u>3.70</u> | <u>13.43</u> | <u>0.0</u> | <u>141</u> | |
| <u>0846</u> | <u>24.47</u> | <u>5.92</u> | <u>0.204</u> | <u>3.92</u> | <u>12.12</u> | <u>> 1.0</u> | <u>108</u> | |
| <u>0850</u> | <u>24.47</u> | <u>6.06</u> | <u>0.297</u> | <u>4.04</u> | <u>12.18</u> | <u>> 1.0</u> | <u>99</u> | |
| <u>0854</u> | <u>24.50</u> | <u>6.14</u> | <u>0.289</u> | <u>6.28</u> | <u>12.76</u> | <u>> 1.0</u> | <u>81</u> | |
| <u>0858</u> | <u>24.48</u> | <u>6.24</u> | <u>0.295</u> | <u>6.08</u> | <u>12.69</u> | <u>> 1.0</u> | <u>94</u> | |
| <u>0902</u> | <u>24.48</u> | <u>6.29</u> | <u>0.297</u> | <u>5.85</u> | <u>12.7</u> | <u>> 1.0</u> | <u>100</u> | |
| <u>0906</u> | <u>24.47</u> | <u>6.24</u> | <u>0.298</u> | <u>5.90</u> | <u>11.83</u> | <u>> 1.0</u> | <u>104</u> | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>3/25/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks Collected field duplicate at this location (OL-MW-11)
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-8
Sample #: _____

Groundwater Sampling Field Data Sheet

DL-MW-8

| | | | |
|-----------------|--|------------|-----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>3/24/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW/RIFSSampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>J Bennett / M Kestor</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>same</u> |

Casing Diameter: 2" 4" 6" Other

| | | | |
|---------------------------------------|---------------------|------------------------|-----------------------------|
| Depth to Water (feet) | <u>18.75</u> | Purge Vol. Meas.Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>3/24/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>09:00 - 10:10</u> |
| Date/Time Sampled | <u>3/24/11 1030</u> | Flow = | <u>410 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 4

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|--------------|-------------------------------|------------------------|--------------------|
| <u>1003</u> | <u>19.67</u> | <u>6.66</u> | <u>0.149</u> | <u>5.87</u> | <u>13.21</u> | <u>54.7</u> | <u>38</u> | |
| <u>1007</u> | <u>19.69</u> | <u>6.54</u> | <u>0.156</u> | <u>6.03</u> | <u>11.28</u> | <u>59.3</u> | <u>35</u> | |
| <u>1011</u> | <u>19.66</u> | <u>6.47</u> | <u>0.159</u> | <u>5.72</u> | <u>11.47</u> | <u>44.0</u> | <u>30</u> | |
| <u>1015</u> | <u>19.68</u> | <u>6.48</u> | <u>0.160</u> | <u>5.93</u> | <u>11.56</u> | <u>34.0</u> | <u>34</u> | |
| <u>1019</u> | <u>19.68</u> | <u>6.47</u> | <u>0.166</u> | <u>5.66</u> | <u>11.60</u> | <u>38.0</u> | <u>37</u> | |
| <u>1023</u> | <u>19.67</u> | <u>6.49</u> | <u>0.150</u> | <u>5.13</u> | <u>11.69</u> | <u>31.4</u> | <u>36</u> | |
| <u>1027</u> | <u>19.67</u> | <u>6.46</u> | <u>0.140</u> | <u>4.88</u> | <u>11.74</u> | <u>32.3</u> | <u>39</u> | |

| | | | |
|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>3/24/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

* Comminance depth.

Parametrix, Inc.

Well #: MW-10
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-10

| | | | |
|-----------------|--|------------|-----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>3/24/2011</u> |
| Project Name | <u>Olalla LF Qtrly GW/RIFSSampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>J Bennett / M Bunker</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>same</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|-----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>26.89</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | <u>—</u> | Date Purged | <u>3/24/2011</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1302 - 1420</u> |
| Date/Time Sampled | <u>3/24/2011 1335</u> | Flow = | <u>340 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) 5

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|--------------|-------------------------------|------------------------|--------------------|
| <u>1306</u> | <u>27.74*</u> | <u>6.51</u> | <u>0.910</u> | <u>0.16</u> | <u>13.41</u> | <u>61.9</u> | <u>33</u> | |
| <u>1310</u> | <u>27.75</u> | <u>6.29</u> | <u>0.990</u> | <u>0.00</u> | <u>11.70</u> | <u>66.3</u> | <u>67</u> | |
| <u>1314</u> | <u>27.73</u> | <u>6.24</u> | <u>0.997</u> | <u>0.00</u> | <u>12.08</u> | <u>57.8</u> | <u>74</u> | |
| <u>1318</u> | <u>27.72</u> | <u>6.25</u> | <u>0.999</u> | <u>0.00</u> | <u>12.18</u> | <u>54.3</u> | <u>83</u> | |
| <u>1322</u> | <u>27.73</u> | <u>6.27</u> | <u>0.999</u> | <u>0.00</u> | <u>12.40</u> | <u>52.8</u> | <u>91</u> | |
| <u>1326</u> | <u>—</u> | <u>6.27</u> | <u>0.999</u> | <u>0.00</u> | <u>12.41</u> | <u>50.5</u> | <u>93</u> | |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>3/24/2011</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

| | | | | | |
|----------------|---|------|----------|----|----------|
| Well Integrity | <u>Good</u> | Page | <u>1</u> | of | <u>1</u> |
| Remarks | <u>Collected, M6/M60 at this location</u> | | | | |
| Signature | <u>[Signature]</u> | | | | |

* Convenience Depth

Sampling Field Data Sheet

Station #: SW-2

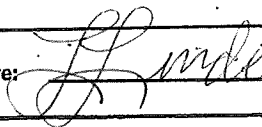
| | | | |
|------------------|------------------------|--------------|-----------------------------|
| Project Number: | <u>215-1578-121</u> | Date: | <u>3/23/11</u> |
| Project Name: | <u>Olalla LF R1/FS</u> | Client Name: | <u>Kitsap Cty SWD</u> |
| Project Address: | <u>Bandix Rd</u> | Sampled By: | <u>L. Linde / M. Baxter</u> |

| TIME (2400 hr) | pH (units) (± 0.1) | Ec (µmhos/cm 25°C) (± 3%) | TEMPERATURE °C | DO (mg/L) | COLOR (visual) | TURBIDITY (visual) |
|--------------------------------|--------------------|---------------------------|----------------|-----------|----------------|--------------------|
| <u>No parameters collected</u> | | | | | | |

| | | | |
|----------------------------|------------|-------------------------------------|------------|
| Sampling Equipment: | <u>N/A</u> | | |
| Laboratory: | <u>N/A</u> | Date Sent to Lab: | <u>N/A</u> |
| Chain-of-Custody (yes/no): | <u>N/A</u> | Field QC Sample Number: | <u>N/A</u> |
| Shipment Method: | <u>N/A</u> | Split With (names[s]/organization): | <u>N/A</u> |

Remarks:

No flow observed @ this location therefore no sample was collected

Signature: 

Sampling Field Data Sheet

Station #: SW-3

| | |
|--------------------------------------|--|
| Project Number: <u>215-2578-121</u> | Date: <u>3/23/11</u> |
| Project Name: <u>Olalla LF RI/FS</u> | Client Name: <u>Kitsap City SWD</u> |
| Project Address: <u>Bendix Rd</u> | Sampled By: <u>L. Lindel / M. Baxter</u> |

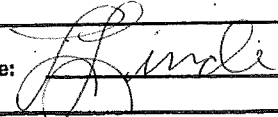
| TIME (2400 hr) | pH (units) (± 0.1) | Ec (µmhos/cm 25°C) (± 3%) | TEMPERATURE °C | DO (mg/L) | ORP COLOR (visual)(mV) | TURBIDITY (visual) (NTU) |
|-------------------|-----------------------|------------------------------|----------------|--------------|------------------------------|-----------------------------|
| <u>1139</u> | <u>6.79</u> | <u>0.061</u> | <u>11.12</u> | <u>13.60</u> | <u>133</u> | <u><1</u> |

| | |
|--|--|
| Sampling Equipment: <u>Grab</u> | |
| Laboratory: <u>ARI</u> | Date Sent to Lab: <u>3/24/11</u> |
| Chain-of-Custody (yes/no): <u>Yes</u> | Field QC Sample Number: <u>N/A</u> |
| Shipment Method: <u>Hand delivered</u> | Split With (names[s]/organization): <u>N/A</u> |

Remarks:

collected @ 1140

Flow rate: ~ 1 gal/min

Signature: 

Sampling Field Data Sheet

Station #: SW-4

| | |
|-------------------------------------|---|
| Project Number: <u>215-1578-121</u> | Date: <u>3/23/11</u> |
| Project Name: <u>Ball Lake SWD</u> | Client Name: <u>Kitap Co SWD</u> |
| Project Address: <u>Bardix Rd</u> | Sampled By: <u>L. Linde / M. Baxter</u> |

| TIME (2400 hr) | pH (units) (± 0.1) | Ec (µmhos/cm 25°C) (± 3%) | TEMPERATURE °C | DO (mg/L) | CRP COLOR (visual) (mv) | TURBIDITY (visual) (NTU) |
|-------------------|-----------------------|------------------------------|----------------|--------------|-------------------------------|-----------------------------|
| <u>1147</u> | <u>6.83</u> | <u>0.077</u> | <u>19.41</u> | <u>13.23</u> | <u>129</u> | <u>41</u> |

| | |
|--|--|
| Sampling Equipment: <u>Grab</u> | |
| Laboratory: <u>ARI</u> | Date Sent to Lab: <u>3/24/11</u> |
| Chain-of-Custody (yes/no): <u>Yes</u> | Field QC Sample Number: <u>N/A</u> |
| Shipment Method: <u>Hand Delivered</u> | Split With (names[s]/organization): <u>N/A</u> |

Remarks:

Collected @ 1150, OL-SW-4

Flow rate ~ 300 ml/min

Steady flow, no pulsing exhibited

Signature: Linde

Landfill Gas Sampling Field Data
Olalla Landfill Gas Sampling

Flare:

Date:

Field Team:

Field Conditions:
 Ambient Temperature:

| Vent | Time | CH ₄ | O ₂ | CO ₂ | Temp °C | Pressure (LSP) | Balance |
|---------|------|-----------------|----------------|-----------------|---------|----------------|---------|
| Flare 3 | 1233 | 29.9 | 1.3 | 7.8 | - | 10.5 | 61.0 |
| 2 | 1241 | 33.4 | 1.7 | 7.9 | - | 10.2 | 57.0 |
| 1 | 1245 | 27.5 | 4.7 | 6.8 | - | 10.2 | 61.0 |
| | | | | | | | |
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Note: surface holes dug in cap between flares 1 & 2

Dealla LF R/FS

3/23/11

Activities: Collect ^{new} samples

Weather: Sunny, ^{in 45°F}

Onsite: L. Lindell, M. Baxter

Author: L. Lindell

0715 Depart to get explosives

0730 Stop @ gas station to fill up generator & get ice

0755 Arrive @ Park

0815 Mike B. arrives - load up rest of gear

0830 Depart for site

0900 Arrive @ site - set up
Calibrate Hanna - good

0941 Begin purging mw-1

Late Entry 0920 Call lab to clarify
IL amber bottle count - thought
there were supposed to be 2 per
well but lab said only 1

0950 Kelli Warren arrives to hand
off tubing

1010 Collect sample @ mw-1, OL-mw-1

1020 Cleanup, move to mw-SA &
set up

1047 Begin purging mw-SA

Olalla LF R/FS

3/23/11

- 1110 Finish purging, collect sample @ MW-5A, OL-MW-5A
- 1120 Cleanup & move locations
- 1140 Collect sample @ SW-3
OL-SW-3
- 1150 Collect sample @ SW-4
OL-SW-4
- 1200 Break for lunch
- 1215 Call Jesse Bennett to bring extra coolers tomorrow & back up gear
- 1230 Decon pump & prep for landfill wells
- 1300 Cart gear to MW-2 & set up
- 1330 Begin purging MW-2
- Late Entry 1245 Leave coolers in office (Purgallup) & Mike W. will P/n in AM to deliver to lab
- 1400 Complete purge, collect sample @ MW-2, OL-MW-2
- 1415 Decon pump, move gear to next location
- 1443 Begin purging MW-4
- 1510 Collect sample @ MW-4, OL-MW-4
- 1520 Cleanup, cart everything to car

Olalla LF R/FS

3/23/11

- 1530 Decon pump, cleanup
- 1600 Collect rinsate on pump, OL-EP
Collect ~~the~~ trip block for tomorrow w/ time of 0000
- 1630 Depart site, secure gate
- 1700 Arrive @ P&R, Mike will leave samples @ office for Mike W. to P/n & take to lab in the morning

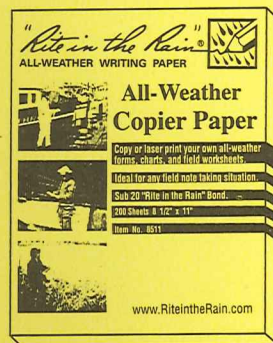
James 3/23/11

Thurs - March 24, 2011

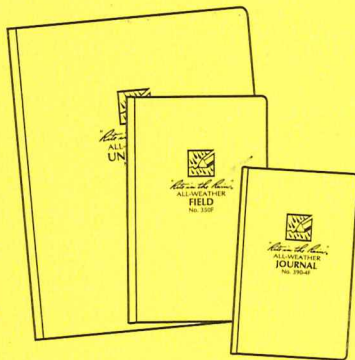
Samplers: Jesse Bennett, Mike Baxter
 Forecast: Hi 54 Lo 41 Chance of Shower
 Current Cond.: Overcast, light showers, no breeze

- 0800 Met Jesse at transfer station
 0815 Calibrate WQ Meter
 0825 Arrive at MW-7
 0840 Start purge
 0915 Collect MW-7
 0925 Collect Duplicate
 0940 Leave MW-7, return to Jesse's truck
 0950 Arrive at MW-8
 0958 Start purge
 1030 Collect MW-8
 1040 Stop purge
 1047 Depart MW-8
 1052 Arrive at MW-6
 1102 Start purge
 1130 Collect MW-6
 1137 End purge
 1145 Depart MW-6 and site to get packing
 tape in G.H.
 1240 Arrive at MW-10
 1302 Start purge

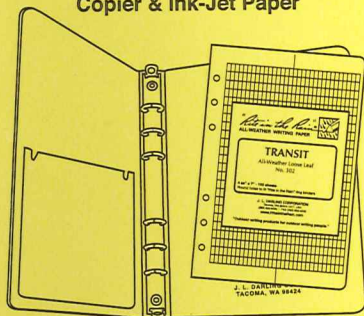
"Rite in the Rain"[®]
 ALL-WEATHER WRITING PAPER



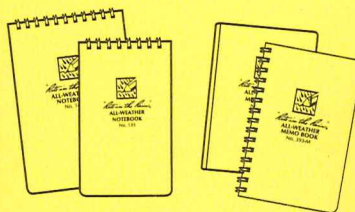
Copier & Ink-Jet Paper



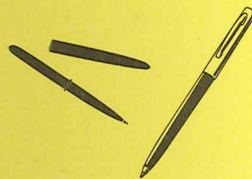
Bound Books



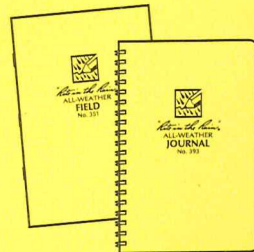
Loose Leaf / Ring Binder



Memo Books



All-Weather Pens

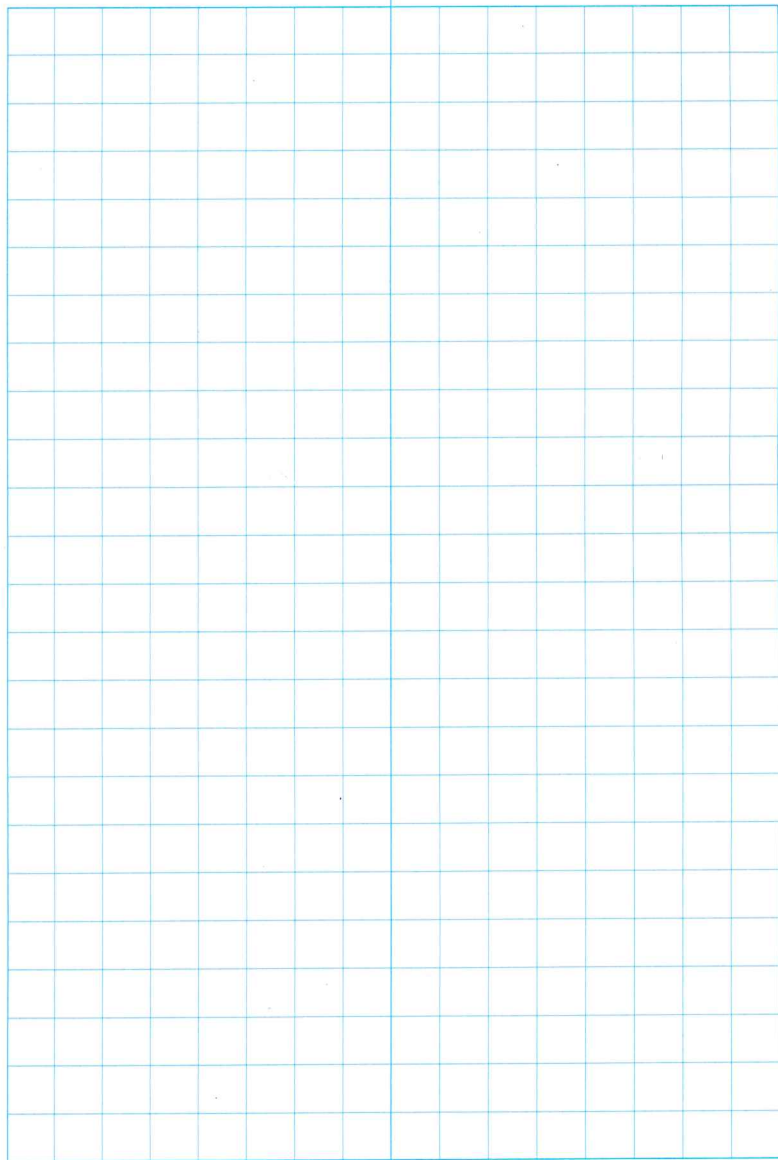
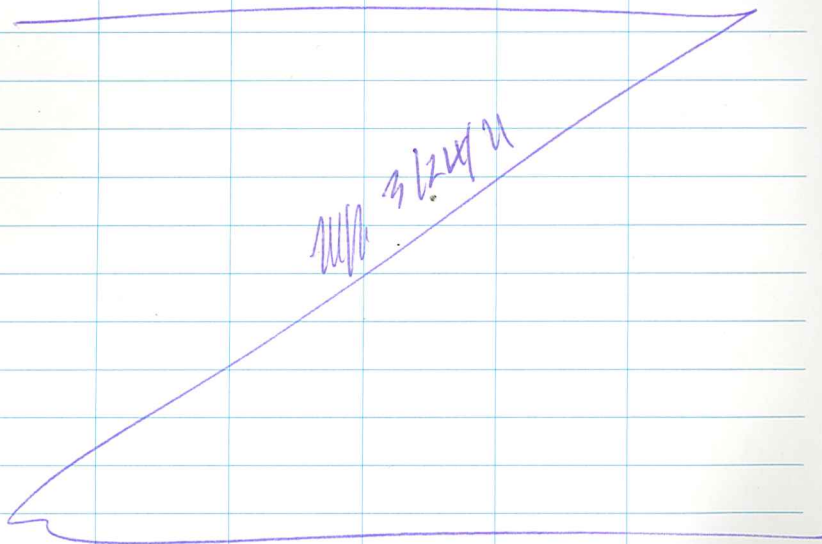


Notebooks

www.RiteintheRain.com

Thursday, March 24, 2011

- 1335 Collect MW-10, plus M6/M6B
- 1420 End purge
- 1435 Depart MW-10
- 1440 Arrive at MW-3
- 1445 Start purge
- 1430 Collect MW-3 sample
- 1540 End purge
- 1545 Depart MW-3
- 1550 Transfer samples and equipment to Jesse's truck, complete COC
- 1600 Depart project site, lock gate behind us.





Kitsap County Public Works

Solid Waste Division

614 Division Street (MS-27)
 Port Orchard, WA 98366
 Phone: 360-337-5777
 FAX: 360-337-4867

CHAIN OF CUSTODY RECORD

FOR LAB USE ONLY

| | |
|-----------------|---|
| LOGIN COMMENTS: | Sample(s) checked/logged by: _____ |
| | Sample temp on receipt at lab: _____ COMPLIANCE 4 +/- 4° C |
| | Are custody seals intact? <input type="checkbox"/> YES <input type="checkbox"/> NO |
| | Sample receipt (day/time): _____ / _____ |
| | Ice? <input type="checkbox"/> NONE <input type="checkbox"/> BLUE <input type="checkbox"/> WET |

TO BE COMPLETED BY THE SAMPLER:

TAT REQUESTED: STANDARD: OTHER: _____ PROJECT: Ball Lake ^{Individual} Lara Lindley / Mike Buxton SAMPLER PRINTED NAME AND SIGNATURE

| SAMPLE DATE | SAMPLE TIME | SITE NAME OR LOCATION | SAMPLE IDENTIFIER | MATRIX * | GRAB | COMP | ANALYSES REQUIRED — MARK AN X IN ALL TESTS REQUIRED FOR EACH SAMPLE LINE | | | | | | | | | | | |
|-------------|-------------|-----------------------|-------------------|----------|------|------|--|-------------|------------|----------|-------------|-------------|-----|----------|----------|---------|---------|------------------------------|
| | | | | | | | VOCs, VC | SVOCs, PCBs | T. Cyanide | TOC, COD | T. Coliform | F. Coliform | TDS | T. NH4-N | D. NH4-N | Ammonia | NITRATE | |
| 3/23 | 1010 | | DL-MW-1 | W | X | | X | X | X | X | X | X | X | X | X | X | X | ① Hg, Ni |
| | 1110 | | DL-MW-5A | | | | | | | | | | | | | | | ② Se, As, Ba, |
| | 1400 | | DL-MW-2 | | | | | | | | | | | | | | | Be, Cd, Cr, |
| | 1510 | | DL-MW-4 | | | | | | | | | | | | | | | Co, Cu, |
| | 1140 | | DL-SW-3 | | | | | | | | X | | | | | | X | Fe, Pb, Mg |
| | 1150 | | DL-SW-4 | | | | | | | | X | | | | | | X | Mn, K, Se, |
| | 1600 | | DL-ER | | | | | | | | X | X | | | | X | | Ag, Na, Thallium, |
| 3/15 | 0000 | | DL-TB | | | | | | | | X | | | | | | | Sn, V, Zn, field filtered |

* MATRIX TYPES: SW = Surface Water GW = Ground Water SO = Soil ^{③ Alkalinity, ammonia, carbonate, bicarbonate, chloride, nitrate, nitrite, sulfate, pH}

| SIGNATURE | PRINTED NAME | COMPANY AND TITLE | DATE | TIME |
|--------------------|--------------|-------------------------|-----------|------|
| <i>[Signature]</i> | Mike Buxton | Parametrix / Env Tech I | 3/24/2011 | 1800 |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |



Kitsap County Public Works
Solid Waste Division
 614 Division Street (MS-27)
 Port Orchard, WA 98366
 Phone: 360-337-5777
 FAX: 360-337-4867

CHAIN OF CUSTODY RECORD

FOR LAB USE ONLY

| | |
|-----------------|---|
| LOGIN COMMENTS: | Sample(s) checked/logged by: _____ |
| | Sample temp on receipt at lab: _____ COMPLIANCE 4 +/- 4° C |
| | Are custody seals intact? <input type="checkbox"/> YES <input type="checkbox"/> NO |
| | Sample receipt (day/time): _____ / _____ |
| | Ice? <input type="checkbox"/> NONE <input type="checkbox"/> BLUE <input type="checkbox"/> WET |

TO BE COMPLETED BY THE SAMPLER:

| | | |
|---|----------------------------|--|
| TAT REQUESTED: STANDARD: <input checked="" type="checkbox"/> OTHER: _____ | PROJECT: <u>Walla Wall</u> | SAMPLER PRINTED NAME AND SIGNATURE: <u>Mike Baxter / Jesse Bennett</u> |
|---|----------------------------|--|

| SAMPLE DATE | SAMPLE TIME | SITE NAME OR LOCATION | SAMPLE IDENTIFIER | MATRIX * | GRAB | COMP | ANALYSES REQUIRED — MARK AN X IN ALL TESTS REQUIRED FOR EACH SAMPLE LINE | | | | | | | | | | |
|-------------|-------------|-----------------------|-------------------|----------|------|------|--|-----------|----|----------|--------------|-----|-----------|-----------|--------------|--|--------------------------------------|
| | | | | | | | VOL.VC | ALC. 12.3 | CN | TOT. COP | T. COLIFORMS | TDS | T. Metals | P. Metals | Conductivity | | |
| 3/24 | 0915 | | OL MW-7 | GW | X | | X | X | X | X | X | X | X | X | | | |
| | 1030 | | OL MW-8 | | | | | | | | | | | | | | |
| | 1130 | | OL MW-6 | | | | | | | | | | | | | | |
| | 1330 | | OL MW-10 | | | | | | | | | | | | | | MS/MEO |
| | 1530 | | OL MW-3 | | | | | | | | | | | | | | |
| | 0800 | | OL TB | BIW | | | | | | | | | | | | | Disc. Metals samples still fill out. |

* MATRIX TYPES: SW = Surface Water GW = Ground Water SO = Soil

| SIGNATURE | PRINTED NAME | COMPANY AND TITLE | DATE | TIME |
|---------------------------------------|-------------------------|------------------------------------|----------------|-------------|
| RELINQUISHED BY: <u>Jesse Bennett</u> | <u>H. Jesse Bennett</u> | <u>Pro-Analysis, Sr. Env. Dir.</u> | <u>3/25/11</u> | <u>0843</u> |
| RECEIVED BY: <u>[Signature]</u> | <u>S. HOFFMAN</u> | <u>KCPW</u> | <u>3/25/11</u> | <u>0843</u> |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |
| RELINQUISHED BY: | | | | |
| RECEIVED BY: | | | | |

Olalla Landfill

OL-MW-##-W-0 ^{primary}

1 dup
2 trip
3 equip

0945 onsite

Met Kel: Means McKay
Jan Brower, KCHD
Eric Caddy, EPS
Mike Warfel, parametric

Glacier
Cascade
~~Paul~~
Noel Phillips
↓
Ecology
Well
construction
coordinator

10/7/10

MW-2 water level @ 11:12 am 66.89' toC

OL-MW-02-W-0 collected @ "11:00"

MW-4 water level @ 11:59 64.09' toC

OL-MW-04-W-0 collected @ "12:00"

Sampled by

Mike & Mary
Warfel Holder

with bailers
put on ice

Samples taken by Mike to
Aquatic Research

Approximately 13:30 started digging Test Pit 1

~ 4:00

Test Pit 2

~ 14:15

Test Pit 3

~ 14:35

Test Pit 4

~ 14:50 dug out drum & additional area @ TP 4 (North end)

~ 15:20

Test Pit 5

N. Phillips off site

~ 15:45 started backfilling

M. Warfel & J. Brower off site

reburied all waste

K. McKay-Means off site

16:15 M. Holder off site only E. Caddy & Ben Bohren (Glacier)
onsite - finishing the backfilling of TP 2 & TP 1.

Multi Rae (Oil and Safety) used to measure PID readings.

Parametrix, Inc.

Well #: MW-1
 Sample #: _____
 DL-MW-1-0

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/28/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>79.60</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/28/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>10:55 - 1111</u> |
| Date/Time Sampled | <u>12/28/10 1115</u> | | <u>500 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|-------------------------------|------------------------|--------------------|
| <u>1055</u> | <u>79.61</u> | <u>6.69</u> | <u>13.0</u> | <u>12.00</u> | <u>9.9</u> | <u>0.9</u> | <u>145</u> | |
| <u>1059</u> | <u>76.61</u> | <u>6.53</u> | <u>13.2</u> | <u>11.06</u> | <u>10.4</u> | <u>1.8</u> | <u>137</u> | |
| <u>1103</u> | <u>76.81</u> | <u>6.48</u> | <u>13.2</u> | <u>10.87</u> | <u>11.5</u> | <u>0.4</u> | <u>128</u> | |
| <u>1107</u> | — | <u>6.47</u> | <u>13.2</u> | <u>10.45</u> | <u>12.0</u> | <u>< 0.1</u> | <u>128</u> | |
| <u>1111</u> | <u>76.89</u> | <u>6.41</u> | <u>13.0</u> | <u>10.45</u> | <u>12.6</u> | <u>0.4</u> | <u>129</u> | <u>5 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-2
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-2-0

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/28/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>64.84</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/28/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1352 - 1408</u> |
| Date/Time Sampled | <u>12/28/10 1410</u> | | <u>420 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ±10% if >10NTU | ORP (mV) ±10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|----------------------------|----------------------|--------------------|
| <u>1352</u> | <u>64.96</u> | <u>7.33</u> | <u>11.4</u> | <u>10.26</u> | <u>12.0</u> | <u>6.8</u> | <u>68</u> | |
| <u>1356</u> | <u>64.88</u> | <u>7.20</u> | <u>11.4</u> | <u>10.13</u> | <u>13.1</u> | <u>4.5</u> | <u>65</u> | |
| <u>1400</u> | <u>64.92</u> | <u>7.13</u> | <u>11.4</u> | <u>10.23</u> | <u>13.3</u> | <u>4.7</u> | <u>65</u> | |
| <u>1404</u> | <u>64.90</u> | <u>7.09</u> | <u>11.3</u> | <u>10.08</u> | <u>13.9</u> | <u>7.2</u> | <u>64</u> | |
| <u>1408</u> | <u>64.92</u> | <u>7.08</u> | <u>11.4</u> | <u>10.15</u> | <u>13.7</u> | <u>8.6</u> | <u>64</u> | <u>4 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good

Remarks _____

Signature L. Linde Page 1 of 1

Parametrix, Inc.

Well #: 1NW-3
 Sample #: _____
 OL-MW-30

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|------------------------------|
| Depth to Water (feet) | <u>42.32</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/29/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>0842-0904</u> |
| Date/Time Sampled | <u>12/29/10 0910</u> | | <u>490 m²/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|------------|-------------------------------|------------------------|--------------------|
| 0842 | 42.38 | 5.83 | 35.5 | 7.17 | 9.9 | * | 211 | |
| 0846 | 42.36 | 5.93 | 35.9 | 5.65 | 10.7 | * | 201 | |
| 0850 | — | 6.03 | 35.8 | 5.20 | 11.4 | * | 191 | |
| 0854 | — | 6.11 | 35.8 | 4.91 | 12.2 | * | 182 | |
| 0856 | 42.34 | 6.14 | 35.9 | 4.81 | 12.5 | * | 177 | |
| 0900 | 42.36 | 6.15 | 35.8 | 4.70 | 12.5 | + | 173 | |
| 0904 | 42.35 | 6.16 | 35.8 | 4.68 | 12.5 | + | 170 | 45 gal |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/30/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

* turbidity sensor not recording

Parametrix, Inc.

Well #: MW-4
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-4-2

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/28/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" X 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>62.19</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/28/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1448 - 1504</u> |
| Date/Time Sampled | <u>12/28/10 1510</u> | | <u>420 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10NTU | ORP (mV) ± 10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|------------------------------|-----------------------|--------------------|
| <u>1418</u> | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| <u>1448</u> | _____ | _____ | _____ | _____ | _____ | _____ | _____ | _____ |
| <u>1452</u> | <u>62.23</u> | <u>7.09</u> | <u>12.7</u> | <u>9.29</u> | <u>12.2</u> | <u>6.0</u> | <u>66</u> | _____ |
| <u>1456</u> | <u>62.23</u> | <u>7.12</u> | <u>12.7</u> | <u>9.14</u> | <u>13.8</u> | <u>2.2</u> | <u>56</u> | _____ |
| <u>1500</u> | <u>62.23</u> | <u>7.13</u> | <u>12.7</u> | <u>9.29</u> | <u>13.9</u> | <u>2.0</u> | <u>55</u> | _____ |
| <u>1504</u> | <u>62.23</u> | <u>7.11</u> | <u>12.7</u> | <u>9.36</u> | <u>13.9</u> | <u>0.3</u> | <u>55</u> | <u>4 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good

Remarks _____

Signature L. Linde Page 1 of 1

Parametrix, Inc.

Well #: MW-5A
 Sample #: _____
OL-MW-5A-D

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/20/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>76.50</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/28/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1146-1202</u> |
| Date/Time Sampled | <u>12/20/10 1205</u> | | <u>440 mL/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ±10% if >10NTU | ORP (mV) ±10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|----------------------------|----------------------|--------------------|
| <u>1146</u> | <u>76.57</u> | <u>6.75</u> | <u>20.7</u> | <u>11.46</u> | <u>11.1</u> | <u>1.8</u> | <u>136</u> | |
| <u>1150</u> | <u>76.57</u> | <u>6.63</u> | <u>20.7</u> | <u>9.72</u> | <u>11.9</u> | <u><0.1</u> | <u>127</u> | |
| <u>1154</u> | <u>76.56</u> | <u>6.63</u> | <u>20.1</u> | <u>9.41</u> | <u>13.0</u> | <u><0.1</u> | <u>122</u> | |
| <u>1158</u> | <u>76.56</u> | <u>6.63</u> | <u>19.8</u> | <u>9.43</u> | <u>13.4</u> | <u><0.1</u> | <u>121</u> | |
| <u>1202</u> | <u>76.56</u> | <u>6.62</u> | <u>19.4</u> | <u>9.52</u> | <u>13.5</u> | <u><0.1</u> | <u>121</u> | <u>45 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/20/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: mw-6
 Sample #: _____
OL-MW-6-0

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>17.74</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/29/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>0934-0959</u> |
| Date/Time Sampled | <u>12/29/10 1005</u> | | <u>500 inl/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|-------------------------------|-----------------------|--------------------|
| <u>0934</u> | <u>17.83</u> | <u>6.49</u> | <u>35.1</u> | <u>8.11</u> | <u>10.0</u> | <u>*</u> | <u>33</u> | |
| <u>0938</u> | <u>17.84</u> | <u>6.53</u> | <u>27.5</u> | <u>5.33</u> | <u>11.0</u> | <u>*</u> | <u>7</u> | |
| <u>0943</u> | <u>-</u> | <u>6.64</u> | <u>23.9</u> | <u>9.08</u> | <u>11.4</u> | <u>*</u> | <u>-3</u> | |
| <u>0947</u> | <u>17.88</u> | <u>6.62</u> | <u>13.0</u> | <u>4.58</u> | <u>12.5</u> | <u>*</u> | <u>-11</u> | |
| <u>0951</u> | <u>17.91</u> | <u>6.62</u> | <u>16.7</u> | <u>4.35</u> | <u>12.7</u> | <u>*</u> | <u>-18</u> | |
| <u>0955</u> | <u>17.90</u> | <u>6.62</u> | <u>16.2</u> | <u>4.29</u> | <u>12.8</u> | <u>*</u> | <u>-22</u> | |
| <u>0959</u> | <u>17.90</u> | <u>6.63</u> | <u>16.3</u> | <u>4.28</u> | <u>12.7</u> | <u>*</u> | <u>-24</u> | <u>45 gal</u> |
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lost power →

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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/30/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>OL-MW-6-1</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature L. Linde Page 1 of 1

Duplicate collected @ 1015

Parametrix, Inc.

Well #: MW-7
 Sample #: _____
 OL MW-7-8

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>24.38</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/29/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1126-1136</u> |
| Date/Time Sampled | <u>12/29/10 1140</u> | | <u>500 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ±10% if >10NTU | ORP (mV) ±10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|----------------------------|----------------------|--------------------|
| <u>1120</u> | <u>24.44</u> | <u>6.77</u> | <u>13.0</u> | <u>11.41</u> | <u>10.3</u> | <u><0.1</u> | <u>75</u> | |
| <u>1124</u> | <u>24.44</u> | <u>6.78</u> | <u>13.0</u> | <u>8.92</u> | <u>11.2</u> | <u>*</u> | <u>75</u> | |
| <u>1128</u> | <u>24.44</u> | <u>6.78</u> | <u>12.9</u> | <u>8.58</u> | <u>11.8</u> | <u>*</u> | <u>75</u> | |
| <u>1132</u> | <u>24.44</u> | <u>6.77</u> | <u>12.9</u> | <u>8.43</u> | <u>12.1</u> | <u>*</u> | <u>78</u> | |
| <u>1136</u> | <u>24.43</u> | <u>6.77</u> | <u>12.9</u> | <u>8.41</u> | <u>12.1</u> | <u>*</u> | <u>81</u> | <u>5gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/30/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature L. Linde Page 1 of 1

* turbidity sensor not working

Parametrix, Inc.

Well #: MW-8
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-8-D

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

Depth to Water (feet) 18.79 Purge Vol. Meas. Method Meas. Cup/Stop Watch
 Depth of Well (feet) _____ Date Purged 12/29/10
 Reference Point (surveyors notch/etc) TOC Purge Time (from/to) 1027-1043
 Date/Time Sampled 12/29/10 1045 440 ml/min

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|-------------------------------|------------------------|--------------------|
| <u>1027</u> | <u>18.89</u> | <u>6.91</u> | <u>10.2</u> | <u>11.10</u> | <u>10.8</u> | <u>13.4</u> | <u>29</u> | |
| <u>1031</u> | <u>-</u> | <u>6.77</u> | <u>9.9</u> | <u>9.88</u> | <u>11.7</u> | <u>22.3</u> | <u>14</u> | |
| <u>1035</u> | <u>18.92</u> | <u>6.76</u> | <u>10.0</u> | <u>8.18</u> | <u>12.4</u> | <u>1.5</u> | <u>8</u> | |
| <u>1039</u> | <u>18.92</u> | <u>6.75</u> | <u>10.1</u> | <u>8.18</u> | <u>12.7</u> | <u>20.1</u> | <u>6</u> | |
| <u>1043</u> | <u>18.92</u> | <u>6.75</u> | <u>10.1</u> | <u>8.10</u> | <u>12.8</u> | <u>20.1</u> | <u>6</u> | <u>4 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/30/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature L. Linde Page 1 of 1

Parametrix, Inc.

Well #: MW-10
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-10-0

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/28/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" X 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | <u>27.70</u> | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/28/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1006 - 1022</u> |
| Date/Time Sampled | <u>12/28/10 1025</u> | | <u>300 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|-------------------------------|------------------------|--------------------|
| <u>1006</u> | <u>27.43</u> | <u>6.44</u> | <u>36.1</u> | <u>5.43</u> | <u>11.2</u> | <u>25.1</u> | <u>141</u> | |
| <u>1010</u> | <u>27.43</u> | <u>6.51</u> | <u>35.5</u> | <u>4.22</u> | <u>12.1</u> | <u>2.9</u> | <u>141</u> | |
| <u>1014</u> | <u>27.42</u> | <u>6.52</u> | <u>35.4</u> | <u>4.14</u> | <u>12.3</u> | <u>0.9</u> | <u>138</u> | |
| <u>1018</u> | <u>27.40</u> | <u>6.54</u> | <u>35.5</u> | <u>4.12</u> | <u>12.4</u> | <u>0.6</u> | <u>136</u> | |
| <u>1022</u> | <u>27.40</u> | <u>6.54</u> | <u>35.2</u> | <u>4.11</u> | <u>12.4</u> | <u>0.3</u> | <u>135</u> | <u>4.5 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks Sump repaired last qtr & working ok
 Signature [Signature] Page 1 of 1

Sampling Field Data Sheet

Station #: SW2

| | |
|-------------------------------------|---|
| Project Number: <u>215-1578-121</u> | Date: <u>12/28/10</u> |
| Project Name: <u>Oalla LF R/FS</u> | Client Name: <u>KCSWD</u> |
| Project Address: <u>Bandix Rd</u> | Sampled By: <u>L. Linde, C. Wiggins</u> |

| TIME (2400 hr) | pH (units) (± 0.1) | Ec (µmhos/cm 25°C) ^{ms/cm} (± 3%) | TEMPERATURE °C | DO (mg/L) | Turbidity COLOR (visual) (NTU) | TURBIDITY (visual) |
|-------------------|-----------------------|--|----------------|--------------|--------------------------------------|-----------------------|
| <u>0855</u> | <u>5.73</u> | <u>5.3</u> | <u>16.6</u> | <u>13.24</u> | <u>115</u> | <u>196</u> |

Sampling Equipment: None required except Horiba U-22 for WR

Laboratory: Aquatic Research Date Sent to Lab: 12/29/10

Chain-of-Custody (yes/no): Yes Field QC Sample Number: N/A

Shipment Method: Hand delivered Split With (names[s]/organization): N/A

Remarks:

Flow rate: 1 gal/sec

DL-SW-2-0 collected C 0900

Signature: L. Linde 12/28/10

Sampling Field Data Sheet

Station #: SW3

| | |
|--------------------------------------|---|
| Project Number: <u>215-1578-121</u> | Date: <u>12/28/10</u> |
| Project Name: <u>Dialla LF RI/FS</u> | Client Name: <u>KCSWD</u> |
| Project Address: <u>Bandix Rd</u> | Sampled By: <u>L. Linde, C. Wiggins</u> |

| TIME (2400 hr) | pH (units) (± 0.1) | Ec (µmhos/cm 25°C) (± 3%) | TEMPERATURE °C | DO (mg/L) | Turbidity COLOR (visual) NTA | Turbidity ORP (mV) TURBIDITY (visual) |
|-------------------|-----------------------|------------------------------|----------------|--------------|------------------------------------|--|
| <u>0934</u> | <u>6.05</u> | <u>3.1</u> | <u>6.7</u> | <u>11.43</u> | <u>127</u> | <u>168</u> |

| | |
|--|--|
| Sampling Equipment: <u>None Required except Horiba U-22 for WA</u> | |
| Laboratory: <u>Aquatic Research</u> | Date Sent to Lab: <u>12/29/10</u> |
| Chain-of-Custody (yes/no): <u>Yes</u> | Field QC Sample Number: <u>N/A</u> |
| Shipment Method: <u>Hand delivered</u> | Split With (names[s]/organization): <u>N/A</u> |

Remarks:

flow rate: 1 gal/sec

DE-SW-3-0 collected C0935

Signature: L. Linde 12/28/10

Sampling Field Data Sheet

Station #: SW4

| | |
|--------------------------------------|---|
| Project Number: <u>215-1578-121</u> | Date: <u>12/28/10</u> |
| Project Name: <u>Owalla LF RI/FS</u> | Client Name: <u>KCSWD</u> |
| Project Address: <u>Bandix Rd</u> | Sampled By: <u>L. Linde, C. Wiggins</u> |

| TIME (2400 hr) | pH (units) (± 0.1) | ms/m Ec (µmhos/cm 25°C) (± 3%) | TEMPERATURE °C | DO (mg/L) | Turbidity GOLOR (visual) | ORP TURBIDITY (visual) (m V) |
|----------------|--------------------|--------------------------------|----------------|--------------|--------------------------|------------------------------|
| <u>0917</u> | <u>5.97</u> | <u>4.9</u> | <u>6.7</u> | <u>12.64</u> | <u>20.0</u> | <u>178</u> |

Sampling Equipment: None required except Horiba U-22 for ORP

Laboratory: Aquatic Research Date Sent to Lab: 12/29/10

Chain-of-Custody (yes/no): Yes Field QC Sample Number: N/A

Shipment Method: Hand delivered Split With (names[s]/organization): N/A

Remarks:

discharge pulsing ~ 1 pulse/sec

flow rate : 500 ml/sec

OL-SW-4-0 collected @ 0920

Signature: L. Linde 12/28/10

Landfill Gas Sampling Field Data

Olalla Landfill Gas Sampling

Flare:

Date: 10/28/10

Field Team: LL, CW

Field Conditions: Cloudy, 40°F, Wind ~ 5 mph, snow in forecast

Ambient Temperature:

| Vent | Time | CH ₄ ^{10/28/10} _{LL} | O ₂ | CO ₂ | Temp °C | SP Pressure | Balance |
|--------------------|------|---|----------------|-----------------|---------|----------------|---------|
| Flare 3 | 1554 | 19.1 | 0.9 | 3.6 | | 0.2 | 71.4 |
| Flare 1 | 1601 | 19.0 | 0.5 | 9.3 | | 10.3 | 71.2 |
| Flare 2 | | 0 | 20.1 | 0 | | 10.6 | |
| Flare 2 | 1608 | 20.9 | 0.4 | 9.6 | | 0.3 | 109.1 |
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10/28/10
LL

Parametrix, Inc.

Well #: 0W-1
 Sample #: _____

Groundwater Sampling Field Data Sheet

0L-0W-1-0

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/28/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Transfer Station</u> |
| Project Address | _____ | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD/KCHD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" _____ 6" _____ Other 8"

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | _____ | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/28/10</u> |
| Reference Point (surveyors notch/etc) | _____ | Purge Time (from/to) | <u>1234-1240</u> |
| Date/Time Sampled | <u>12/28/10 1245</u> | | |

Cannot measure

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|------------|-------------------------------|-----------------------|--------------------|
| <u>1234</u> | <u>-</u> | <u>7.32</u> | <u>16.2</u> | <u>11.89</u> | <u>9.2</u> | <u>156.0</u> | <u>111</u> | |
| <u>1237</u> | <u>-</u> | <u>7.31</u> | <u>16.4</u> | <u>10.07</u> | <u>8.2</u> | <u>158.0</u> | <u>113</u> | |
| <u>1240</u> | <u>-</u> | <u>7.28</u> | <u>16.5</u> | <u>9.51</u> | <u>6.1</u> | <u>154.0</u> | <u>112</u> | <u>Unknown</u> |
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|---------------------------|------------------------------|-----------------------------------|----------------------|
| Purge Equipment | <u>Sampling Port/BR SINK</u> | Sampling Equipment | <u>Sampling Port</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

Attendants note rust in water
 Sample collected @ bathroom sink

no way to capture purge water
 or flow rate due to sink
 configuration

Parametrix, Inc.

Well #: OW-3
 Sample #: _____

Groundwater Sampling Field Data Sheet

~~OW-3-0~~ ^{HW} 12/29/10

| | | | |
|-----------------|------------------------------------|------------|------------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>2050 Purkey-Olalla Rd</u> |
| Project Address | <u>2650 Burien-Olalla Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD/KCHD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" _____ 6" _____ Other 8"

Depth to Water (feet) _____ Purge Vol. Meas. Method Meas. Cup/Stop Watch
 Depth of Well (feet) _____ Date Purged 12/29/10
 Reference Point (surveyors notch/etc) _____ Purge Time (from/to) _____
 Date/Time Sampled No sample collected No flow rate collected

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

tank water
 from unknown
 well source

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|------------|-------------------------------|------------------------|--------------------|
| <u>1347</u> | <u>-</u> | <u>6.32</u> | <u>16.7</u> | <u>7.56</u> | <u>5.3</u> | <u>57.5</u> | <u>-141</u> | <u>500 mL</u> |
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|---------------------------|-------------------------|-----------------------------------|----------------------------|
| Purge Equipment | Sampling Port | Sampling Equipment | Sampling Port |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>No sample collected</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Tank part good, well not pumping
 Remarks _____
 Signature L. Linde Page 1 of 1

1350 Cannot confirm water source fr. pressure tank

Parametrix, Inc.

Well #: OW-4
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-OW-4-C

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>OW-4</u> |
| Project Address | <u>2590 Barley Olalla Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD/KCHD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" _____ 6" _____ Other 8"

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------|
| Depth to Water (feet) | _____ | Purge Vol. Meas. Method | <u>Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/29/10</u> |
| Reference Point (surveyors notch/etc) | _____ | Purge Time (from/to) | <u>1407 - 1416</u> |
| Date/Time Sampled | <u>12/29/10 1420</u> | Flow Rate | <u>2L/min</u> |
| | | Purge Volume | <u>7.5 gal purged</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|-------------------------------|-----------------------|--------------------|
| <u>1407</u> | <u>-</u> | <u>7.42</u> | <u>12.2</u> | <u>5.45</u> | <u>10.2</u> | <u>*</u> | <u>-225</u> | |
| <u>1410</u> | <u>-</u> | <u>8.13</u> | <u>12.2</u> | <u>4.12</u> | <u>10.4</u> | <u>*</u> | <u>-248</u> | |
| <u>1413</u> | <u>-</u> | <u>8.27</u> | <u>12.2</u> | <u>3.97</u> | <u>10.6</u> | <u>*</u> | <u>-257</u> | |
| <u>1416</u> | <u>-</u> | <u>8.33</u> | <u>12.2</u> | <u>3.83</u> | <u>10.5</u> | <u>*</u> | <u>-244</u> | <u>7.5 gal</u> |
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|---------------------------|-------------------------|------------------|-----------------------|------------------------|------------|-----------------------------------|------------|
| Purge Equipment | _____ | Sampling Port | _____ | Sampling Equipment | _____ | Sampling Port | _____ |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/30/10</u> | Field QC Sample Number | <u>N/A</u> | Split with (name(s)/organization) | <u>N/A</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Shipment Method | <u>Hand Delivered</u> | | | | |

Well Integrity: Good
 Remarks: Spill water to purge water, 80 gal tank, pump turned on/off twice
 Signature: [Signature] Page _____ of _____

* turbidity sensor not working

Parametrix, Inc.

Well #: OW-9
 Sample #: _____

Groundwater Sampling Field Data Sheet

OW-9-0

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>12/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>OW-9</u> |
| Project Address | <u>13320 Annaple Dr SE</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD/KCHD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" _____ 6" _____ Other 8"

| | | | |
|---------------------------------------|----------------------|-------------------------|-----------------------------|
| Depth to Water (feet) | _____ | Purge Vol. Meas. Method | <u>Meas. Cup/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>12/29/10</u> |
| Reference Point (surveyors notch/etc) | _____ | Purge Time (from/to) | <u>1438 - 1447</u> |
| Date/Time Sampled | <u>12/29/10 1450</u> | | <u>6L/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) ± 0.1 | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% if > 10 NTU | ORP (mV) ± 10mV | CUM. VOL. (gal) |
|-------------------|--------------------------------|------------------------|-------------------------|-----------------------|-------------|-------------------------------|-----------------------|--------------------|
| <u>1438</u> | <u>-</u> | <u>7.68</u> | <u>18.0</u> | <u>6.71</u> | <u>11.0</u> | <u>*</u> | <u>-35</u> | |
| <u>1441</u> | <u>-</u> | <u>7.27</u> | <u>18.0</u> | <u>4.78</u> | <u>11.2</u> | <u>*</u> | <u>-28</u> | |
| <u>1444</u> | <u>-</u> | <u>7.15</u> | <u>18.0</u> | <u>4.57</u> | <u>11.1</u> | <u>*</u> | <u>-24</u> | |
| <u>1447</u> | <u>-</u> | <u>7.05</u> | <u>18.0</u> | <u>4.50</u> | <u>11.1</u> | <u>*</u> | <u>-28</u> | <u>12 gal</u> |
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|---------------------------|-------------------------|-----------------------------------|-----------------|
| Purge Equipment | Sampling Port | Sampling Equipment | Sampling Port |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>12/30/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks No odor
 Signature [Signature] Page 1 of 1

Olalla LF R/FS

12/28/10

Activities: Collect girdsw samples

Weather: Cloudy, 40°F, 5-10 mph wind

Onsite: L. Lindle, C. Wiggins

Author: L. Lindle

0710 Depart for TORO, P/n ice

0730 Meet w/ Mike W (PMX) to get containers for KCHD samples. He informs me to collect all MFS parameters for surface water samples instead of just f. coliform & nitrate

0740 Depart for landfill

0800 Arrive @ Landfill, meet Keli (RESW) to get well & gate keys and Grandfos control box; she will be back w/ Lucas (KCHD) @ 1500 for landfill inspection

0805 Keli departs

0810 Chad W (PMX) arrives, dummies lock note so Keli & Lucas can get in later.

0820 Setup for sampling surface water first - all 3 locations have discharge

0830 Calibrate Horiba - good

0845 Unlock stormwater pond gate

Olalla LF R/FS

12/28/10

0900 Collect sample OL-SW-2-0

0910 Gather gear & move to next location, secure stormwater pond gate

0920 Collect sample OL-SW-4-0

0925 Gather gear, in transit to next location note new ss on installed C access road to MW runoff flowing under access & not through culvert

0935 Collect sample OL-SW-3-0

0945 Go back to MW-10, Mike W requested it be sampled first since it went in for repairs last set up @ MW-10

1000 Begin purging MW-10

1025 Complete purge & collect sample OL-MW-10-0

1030 Cleanup, put purge water staged drum - completely no more water will fit.

1040 Go to MW-1, setup

1055 Begin purging MW-1

1115 ~~Complete~~ Complete purge & collect sample OL-MW-1-0

Olalla LF R/FS

12/28/10

1120 Clean up, purge water not retained

1130 Go to MW-5A, setup, talk to transfer station attendant - he will remove wooden cover so we can inspect pressure tanks in the attendant station bathroom

1146 Begin purging MW-5A

1205 Complete purge, collect sample OL-MW-5A-D, purge water not retained

1210 Cleanup, park C transfer station, inspect pressure tanks (2) (+) 1 sm. hot water heater, no access to sample from a part C tanks, inspect well C entrance to transfer station (batted cover water supply) no parts for sampling, attendants tell us there's 2 exterior spigots but the one to the south is broken, they are unsure if the spigots pull from the same source but give us plans to look at, plans are unhelpful, only source that is confirmed is the bathroom sink, no way to capture flow

Olalla LF R/FS

12/28/10

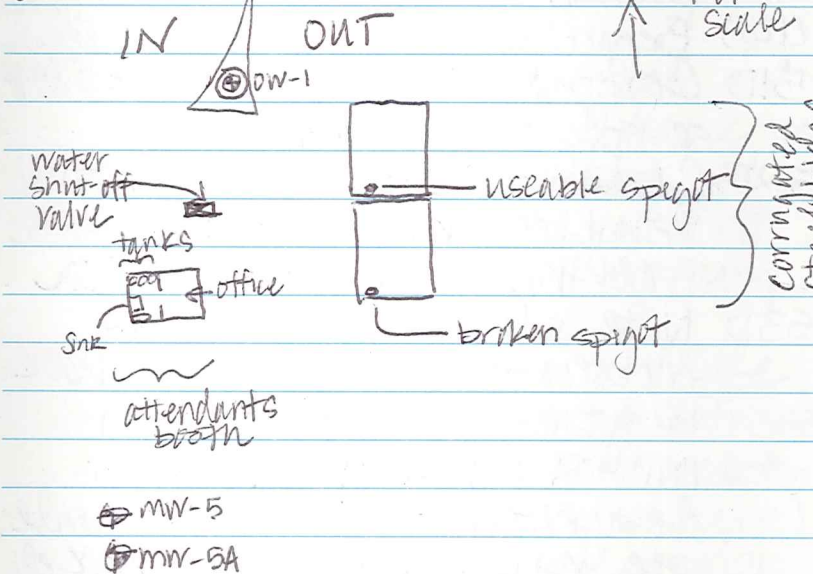
rate or total volume purged due to awkwardness of setup

1234 Begin purging OW-1

1245 Complete purge, collect sample ~~1238~~ at OL-OW-1-D, purge water down sink drain

Burley Olalla Rd

OW-1 Sketch.



1310 Cleanup, go to MW-2 & 4 access on south side of landfill, setup

1320 Decon portable Grundfos

1340 move gear by hand out to MW-2 to setup

1352 Begin purging MW-2

Olalla LF R/FS

12/28/10

1410 Complete purge, collect sample

C OL-mw-2-0, retain ^{purge water}1420 Cleanup, note ^{clear to tan} gelatin-like globs adhering to pump & discharge hose (no odor)

1435 Move gear to mw-4, decon Pump

1448 Begin purging mw-4, ^{purge water} ^{odor like gas}1510 Complete purge, collect ^{large} sample OL-mw-4-0

1520 Cleanup, carry gear back to vehicles, Keli & Lucas on site walking fr. transfer station

1530 Note similar gelatin-like globs on pump & discharge hose

1535 Purge water retained from mw-2 & mw-4 placed in designated drum on roadside, see Keli & Lucas, inform Keli there is an area about the size of a grapefruit between mw-2 & drainage grate in which the liner is exposed

1540 Keli & Lucas continue inspection

1550 Finish cleaning up to start

Olalla LF R/FS

12/28/10

landfill gas readings, calibrate (zero) Landtec, rain imminent

1554 Begin landfill gas readings C Flares (3), call Mike W. (PMV) to meet C TORO @ 1700 to hand off sample, see field sheets for readings

1610 Complete readings

1620 Depart site, secure gate, meet w/ Chad in morning @ same time

1640 Arrive C TORO, fill out COCs

1725 Mike arrives to hand off samples for tomorrow morning delivery

1730 Sign out samples to Mike, inform him of culvert & confirm its for overflow & about gelatin-globs (could be Fe bacteria), sampled all 3 SW stations, wants us to call 1 hr before we're done w/ landfill wells so he can be present for domestic well sample collection

1740 Depart

Finale 12/28/10

Olalla LF R/FS 12/29/10

Activities: Collect gw & dom. well samples

Weather: Partly Cloudy, ~ 35°F

Onsite: L. Linde, C. Wiggins

Arthur: L. Linde

0730 Depart for site

0810 Arrive @ Landfill, Chad onsite

0815 Go to MW-3 & setup, calibrate
Horiba - good

0842 Begin purging MW-3

0910 Complete purge, collect sample
OL-MW-3, retain purge water

0915 Cleanup, move to MW-6, setup

0934 Begin purging MW-6

0940 Receive call from Mike W. (PMX)
NC Rentals will be coming by to
p/n excavator, Horiba turbidity
sensor still not working after rinsing
off w/ DI water

1005 Complete purge, collect sample
OL-MW-6-0

1015 Collect field duplicate OL-MW-6-1

1020 Cleanup, move to MW-8, retain
purge water, setup

1027 Begin purging MW-8

Olalla LF R/FS 12/29/10

1045 Complete purge, collect sample
at MW-8, OL-MW-8-0, retain water

1100 Cleanup, move to MW-7, setup

1120 Begin purging MW-7

1140 Complete purge, collect sample
OL-MW-7-0

1145 Cleanup, purge water from MW-3,
MW-8, & MW-6 in MW-8 dev.
water drum, MW-7 purge water
in empty drum + 1/2 bucket
from MW-8, all drums
full except extra drum
next to dev water drum
for MW-2 & MW-4

1205 Meet w/ Mike, break for lunch
depart & secure gate

1300 Arrive back @ transfer station,
reorganize gear

1320 Go to OW-3 (in red barn-type
shed, attempt to pump water
but have trouble with pressure
tank, turn on valve outside
shed, water begins flowing
but pump in wellhead not
on 12/29/10

Olalla IFR/FS

12/22/10

- 1347 Purge ~ 500 mL & collect one set of field parameters, decide to stop purge since we can't confirm water source
- 1350 Check in w/ owner's wife - will be back a little later when husband gets home
- 1355 Depart & go to OW-4, owner not home, well house open set up, pump is running
- 1407 Begin purging OW-4, water has sulfur odor
- 1420 Finish purge ~ 7.5 gal purged @ 2 l/min from spigot part inside well house, pump cycles twice while purging, well inside well house
- 1430 Depart & go to OW-5, well house locked & owner's not home, will return later
- 1435 Go to OW-9, gain access to well & pressure tank w/ spigot part, set up
- 1438 Begin purging OW-9, no odor observed

collect
OL-OW-4

Olalla IFR/FS

12/22/10

- 1450 Finish purge, ~ 12 gal purged @ 6 l/min from spigot inside shop/garage, pump running, well located in grass between shop/garage & house
- 1500 Go back to OW-3, owner back, control panel for pressure tank was borrowed by neighbor, he will run to neighbor's house to retrieve
- 1510 Depart back to transfer station to wait for owner @ OW-3 to return, fill out COCs & redistribute gear
- 1520 OW-3 owner returns, neighbor is out of town on vacation
- 1530 Depart back to OW-5, owner still not back
- 1545 Sign out samples to Chad to drop off tomorrow, Chad departs
- 1550 Depart site w/ Mike & will return next week to finish

collect
OL-OW-9

J. J. J. 12/22/10



Aquatic Research Incorporated

3927 Aurora Ave. N / Seattle, WA 98103 / (206) 632-2715

CHAIN-OF-CUSTODY RECORD

CLIENT: Kitsap City Health District
SAMPLING DATE: 12/28/10
SAMPLERS: Lara Lindel/Chad Wiggins, Parametrix

SHEET 1 OF 1
PROJECT ID: 01alla 1F
CASE FILE NO.: _____
DATA RECORDED BY: _____

SAMPLE INFORMATION

PARAMETERS

| SAMPLE ID | DATE/TIME COLLECTED | VOL | DIS | BOTT # | NOTES |
|-----------|---------------------|-----|-----|--------|-------|
| DL-OW-1-0 | 12/28/10 1345 | XXX | XXX | | |
| DL-OW-1-2 | 12/28/10 | XX | | | TB |
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|--------------|---------------------------------------|-----------------------------------|---|-------------------------------------|
| Printed Name | Relinquished By <u>Lara Lindel</u> | Date/Time <u>12/28/10 1730</u> | Received By <u>Michael R. Wurfel</u> | Date/Time <u>12/28/2010 1730</u> |
| Signature | | | | |
| Affiliation | <u>Parametrix</u> | | <u>Parametrix</u> | |

| | | | | |
|--------------|---|-------------------------------------|--------------------------------|-----------------------------------|
| Printed Name | Relinquished By <u>Michael K. Wurfel</u> | Date/Time <u>12/29/2010 1038</u> | Received By <u>S. NIGRO</u> | Date/Time <u>12/29/10 1038</u> |
| Signature | | | | |
| Affiliation | <u>Parametrix</u> | | <u>AR</u> | |

Miscellaneous Notes (Hazardous Materials, Quick turn-around time, etc.): * field filtered, metals = As, Fe, Mn



Aquatic Research Incorporated

3927 Aurora Ave. N / Seattle, WA 98103 / (206) 632-2715

CHAIN-OF-CUSTODY RECORD

CLIENT: Kitsap County Health Dept.
 SAMPLING DATE: 1/27/11
 SAMPLERS: L. Lindell (PMX)

SHEET 1 OF 1
 PROJECT ID: _____
 CASE FILE NO.: _____
 DATA RECORDED BY: _____

SAMPLE INFORMATION

PARAMETERS

| SAMPLE ID | DATE/TIME COLLECTED | PARAMETERS | | | | | | | | | | | | | | | | BOTT # | NOTES |
|-----------|---------------------|------------|-----|------------|--|--|--|--|--|--|--|--|--|--|--|--|--|------------------------|-------|
| | | VLS | VLS | Diss. MTLS | | | | | | | | | | | | | | | |
| DL-OW-2-0 | 1/27/11 1230 | XXX | XXX | | | | | | | | | | | | | | | | |
| DL-OW-3-0 | ↓ 1150 | ↓ | ↓ | ↓ | | | | | | | | | | | | | | Diss MTLS - As, Fe, Mn | |
| DL-OW-5-0 | ↓ 1105 | ↓ | ↓ | ↓ | | | | | | | | | | | | | | * field filtered | |
| DL-OW-2-2 | ↓ 0930 | ↓ | ↓ | ↓ | | | | | | | | | | | | | | | |

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|--------------|--------------------------------------|-----------|---------------------|---------------------------------|-----------|-----------------------|
| Printed Name | Relinquished By: <u>Lara Lindell</u> | Date/Time | <u>1/28/11 0730</u> | Received By: <u>[Signature]</u> | Date/Time | <u>7:34 1/28/2011</u> |
| Signature | <u>[Signature]</u> | | | | | |
| Affiliation | <u>PMX</u> | | | | | |

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|--------------|-------------------------------------|-----------|-----------------------|---------------------------------|-----------|------------------|
| Printed Name | Relinquished By: <u>[Signature]</u> | Date/Time | <u>1/28/2011 7:43</u> | Received By: <u>[Signature]</u> | Date/Time | <u>1-28-2011</u> |
| Signature | <u>[Signature]</u> | | | | | <u>7:43 PM</u> |
| Affiliation | | | | | | |

Miscellaneous Notes (Hazardous Materials, Quick turn-around time, etc.):

S. NELSON 1/28/11
[Signature]
AVC 0910

Parametrix, Inc.

Well #: mw-1
 Sample #: 02-mw-1

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|-----------------------------|------------|---------------------|
| Project Number | 215-1578-121 | Date | 11/28/10 |
| Project Name | Olalla LF Qtrly GW Sampling | Location | Olalla LF |
| Project Address | Bandix Rd | Sampled By | L. Linde/C. Wiggins |
| Client Name | KC SWD | Purged By | L. Linde/C. Wiggins |

Casing Diameter: 2" X 4" 6" Other _____

| | | | |
|---------------------------------------|---------------|-------------------------|---------------------|
| Depth to Water (feet) | 79.59 | Purge Vol. Meas. Method | Grad Cyl/Stop Watch |
| Depth of Well (feet) | | Date Purged | 10/28/10 |
| Reference Point (surveyors notch/etc) | TOC | Purge Time (from/to) | 0850-0920 |
| Date/Time Sampled | 0930 11/28/10 | | 450 ml/min |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|------------|----------------|------------------------|--------------------|
| 0850 | 79.81 | 6.30 | 16.9 | 9.95 | 10.1 | 26.1 | 247 | |
| 0854 | 79.70 | 6.28 | 17.2 | 9.90 | 10.1 | 42.6 | 235 | |
| 0858 | 79.71 | 6.26 | 17.3 | 9.90 | 10.0 | 34.8 | 204 | |
| 0902 | 79.72 | 6.25 | 17.2 | 9.91 | 9.9 | 37.4 | 189 | |
| 0906 | 79.71 | 6.24 | 16.3 | 9.85 | 10.2 | 28.4 | 178 | |
| 0910 | | | | | | | | |
| 0912 | 79.72 | 6.41 | 15.0 | 10.51 | 10.3 | 43.8 | 171 | |
| 0916 | 79.63 | 6.36 | 15.0 | 9.23 | 10.5 | 31.7 | 173 | |
| 0920 | 79.66 | 6.33 | 15.0 | 9.76 | 10.5 | 34.3 | 163 | 2.5 gal |
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|---------------------------|----------------------|-----------------------------------|----------------------|
| Purge Equipment | Grundfos Ready-flo 2 | Sampling Equipment | Grundfos Ready-flo 2 |
| Laboratory | Aquatic Research | Date Sent to Lab | 10/29/10 |
| Chain-of-Custody (yes/no) | Yes | Field QC Sample Number | N/A |
| Shipment Method | Hand Delivered | Split with (name(s)/organization) | N/A |

Well Integrity Good
 Remarks having trouble maintaining flow
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-2
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-2

| | | | |
|-----------------|-----------------------------|------------|---------------------|
| Project Number | 215-1578-121 | Date | <u>10/28/10</u> |
| Project Name | Olalla LF Qtrly GW Sampling | Location | Olalla LF |
| Project Address | Bandix Rd | Sampled By | L. Linde/C. Wiggins |
| Client Name | KC SWD | Purged By | L. Linde/C. Wiggins |

Casing Diameter: 2" X 4" 6" Other _____

Depth to Water (feet) 66.63 Purge Vol. Meas. Method Grad. Cyl/Stop Watch
 Depth of Well (feet) _____ Date Purged 10/28/10
 Reference Point (surveyors notch/etc) TOC Purge Time (from/to) 1102 - 1130
 Date/Time Sampled 10/28/10 1145 300 ml/min

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) ± mV | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|-------------|----------------|---------------------|--------------------|
| <u>1102</u> | <u>—</u> | <u>7.18</u> | <u>13.8</u> | <u>9.40</u> | <u>11.5</u> | <u>05.8</u> | <u>102</u> | |
| <u>1110</u> | <u>66.76</u> | <u>7.08</u> | <u>13.3</u> | <u>9.16</u> | <u>13.0</u> | <u>89.1</u> | <u>109</u> | |
| <u>1114</u> | <u>66.78</u> | <u>7.03</u> | <u>13.1</u> | <u>8.86</u> | <u>13.9</u> | <u>66.3</u> | <u>97</u> | |
| <u>1118</u> | <u>66.75</u> | <u>7.02</u> | <u>13.1</u> | <u>8.86</u> | <u>14.2</u> | <u>54.4</u> | <u>106</u> | |
| <u>1122</u> | <u>66.80</u> | <u>7.02</u> | <u>13.1</u> | <u>8.79</u> | <u>14.2</u> | <u>45.9</u> | <u>114</u> | |
| <u>1126</u> | <u>66.81</u> | <u>7.01</u> | <u>13.1</u> | <u>8.76</u> | <u>14.1</u> | <u>45.1</u> | <u>123</u> | |
| <u>1130</u> | <u>—</u> | <u>7.00</u> | <u>13.2</u> | <u>8.75</u> | <u>14.1</u> | <u>22.1</u> | <u>138</u> | <u>55 gal</u> |
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Purge Equipment Grundfos Ready-flo 2 Sampling Equipment Grundfos Ready-flo 2
 Laboratory Aquatic Research Date Sent to Lab 10/29/10
 Chain-of-Custody (yes/no) Yes Field QC Sample Number N/A
 Shipment Method Hand Delivered Split with (name(s)/organization) N/A

Well Integrity Good
 Remarks Difficulty maintaining flow
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-3
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-3

| | | | |
|-----------------|-----------------------------|------------|---------------------|
| Project Number | 215-1578-121 | Date | <u>10/28/10</u> |
| Project Name | Olalla LF Qtrly GW Sampling | Location | Olalla LF |
| Project Address | Bandix Rd | Sampled By | L. Linde/C. Wiggins |
| Client Name | KC SWD | Purged By | L. Linde/C. Wiggins |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|---------------------|
| Depth to Water (feet) | <u>46.80</u> | Purge Vol. Meas. Method | Grad Cyl/Stop Watch |
| Depth of Well (feet) | _____ | Date Purged | <u>10/28/10</u> |
| Reference Point (surveyors notch/etc) | TOC | Purge Time (from/to) | <u>1533-1553</u> |
| Date/Time Sampled | <u>10/28/10 1600</u> | | <u>440 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|------------|----------------|------------------------|--------------------|
| 1533 | 46.79 | 6.22 | 43.1 | 0.62 | 11.0 | 169.0 | 244 | |
| 1537 | 46.80 | 6.03 | 42.5 | 0.27 | 12.1 | 170.0 | 233 | |
| 1541 | 46.80 | 6.07 | 42.5 | 0.38 | 13.1 | 173.0 | 222 | |
| 1545 | 46.80 | 6.14 | 42.5 | 0.34 | 13.3 | 174.0 | 208 | |
| 1549 | 46.80 | 6.20 | 42.5 | 0.36 | 13.4 | 175.0 | 200 | |
| 1553 | 46.81 | 6.28 | 42.7 | 0.40 | 13.5 | 176.0 | 199 | 5 gal |
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|---------------------------|----------------------|-----------------------------------|----------------------|
| Purge Equipment | Grundfos Ready-flo 2 | Sampling Equipment | Grundfos Ready-flo 2 |
| Laboratory | Aquatic Research | Date Sent to Lab | <u>10/29/10</u> |
| Chain-of-Custody (yes/no) | Yes | Field QC Sample Number | <u>OL-MW-9C 1615</u> |
| Shipment Method | Hand Delivered | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

Field duplicate collected OL-MW-9C 1615

Parametrix, Inc.

Well #: MW-4
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-4

| | | | |
|-----------------|-----------------------------|------------|---------------------|
| Project Number | 215-1578-121 | Date | <u>10/28/10</u> |
| Project Name | Olalla LF Qtrly GW Sampling | Location | Olalla LF |
| Project Address | Bandix Rd | Sampled By | L. Linde/C. Wiggins |
| Client Name | KC SWD | Purged By | L. Linde/C. Wiggins |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---|----------------------|-------------------------|---------------------|
| Depth to Water (feet) | <u>64.09</u> | Purge Vol. Meas. Method | Grad Cyl/Stop Watch |
| Depth of Well (feet) | _____ | Date Purged | <u>10/28/10</u> |
| Reference Point (surveyors notch/etc) TOC | _____ | Purge Time (from/to) | <u>1245-1313</u> |
| Date/Time Sampled | <u>10/28/10 1325</u> | | <u>300 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|-------------|----------------|------------------------|--------------------|
| <u>1245</u> | — | <u>7.22</u> | <u>15.5</u> | <u>9.09</u> | <u>11.0</u> | <u>248</u> | <u>236</u> | |
| <u>1249</u> | <u>63.96</u> | <u>7.11</u> | <u>15.3</u> | <u>8.86</u> | <u>11.1</u> | <u>236</u> | <u>238</u> | |
| <u>1253</u> | <u>63.94</u> | <u>7.09</u> | <u>14.7</u> | <u>8.44</u> | <u>11.4</u> | <u>232</u> | <u>202</u> | |
| <u>1357</u> | <u>63.94</u> | <u>7.03</u> | <u>14.5</u> | <u>8.32</u> | <u>12.0</u> | <u>310</u> | <u>177</u> | |
| <u>1361</u> | <u>63.98</u> | <u>7.05</u> | <u>14.4</u> | <u>8.34</u> | <u>12.3</u> | <u>265</u> | <u>164</u> | |
| <u>1365</u> | <u>63.97</u> | <u>7.05</u> | <u>13.5</u> | <u>8.15</u> | <u>15.3</u> | <u>17.7</u> | <u>101</u> | |
| <u>1309</u> | <u>63.98</u> | <u>7.03</u> | <u>14.3</u> | <u>8.16</u> | <u>16.3</u> | <u>44</u> | <u>116</u> | |
| <u>1313</u> | <u>63.98</u> | <u>7.03</u> | <u>14.3</u> | <u>8.23</u> | <u>15.2</u> | <u>29</u> | <u>137</u> | <u>4 gal</u> |

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|---------------------------|----------------------|-----------------------------------|----------------------|
| Purge Equipment | Grundfos Ready-flo 2 | Sampling Equipment | Grundfos Ready-flo 2 |
| Laboratory | Aquatic Research | Date Sent to Lab | <u>10/29/10</u> |
| Chain-of-Custody (yes/no) | Yes | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | Hand Delivered | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity: GOOD
 Remarks: 2" Hard to maintain flow
 Signature: _____ Page 1 of 1

Parametrix, Inc.

Well #: MW-5A
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-5A

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>10/28/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>77.83</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>10/28/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1432-1500</u> |
| Date/Time Sampled | <u>10/28/10 1510</u> | | <u>500 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) ± 10 mV | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|-------------|----------------|------------------------|--------------------|
| <u>1432</u> | <u>77.87</u> | <u>6.59</u> | <u>18.0</u> | <u>9.14</u> | <u>11.5</u> | <u>60.5</u> | <u>239</u> | |
| <u>1438</u> | <u>77.87</u> | <u>6.45</u> | <u>17.6</u> | <u>8.54</u> | <u>11.8</u> | <u>14.9</u> | <u>193</u> | |
| <u>1442</u> | <u>77.88</u> | <u>6.46</u> | <u>17.6</u> | <u>8.58</u> | <u>13.8</u> | <u>10.8</u> | <u>177</u> | |
| <u>1446</u> | <u>77.89</u> | <u>6.52</u> | <u>17.4</u> | <u>8.58</u> | <u>13.6</u> | <u>6.6</u> | <u>172</u> | |
| <u>1450</u> | <u>77.97</u> | <u>6.55</u> | <u>17.2</u> | <u>9.40</u> | <u>14.1</u> | <u>10.7</u> | <u>165</u> | |
| <u>1456</u> | <u>77.98</u> | <u>6.58</u> | <u>17.3</u> | <u>9.32</u> | <u>14.0</u> | <u>13.5</u> | <u>166</u> | |
| <u>1500</u> | <u>77.98</u> | <u>6.55</u> | <u>17.3</u> | <u>9.25</u> | <u>13.4</u> | <u>17.0</u> | <u>175</u> | <u>5.5 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>10/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-6
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-6

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>10/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>21.17</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>10/29/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>0915 - 0931</u> |
| Date/Time Sampled | <u>10/29/10 0940</u> | | <u>500 mL/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|-------------|----------------|-------------|--------------------|
| <u>0915</u> | <u>21.28</u> | <u>6.45</u> | <u>49.5</u> | <u>3.45</u> | <u>10.9</u> | <u>19.8</u> | <u>43</u> | |
| <u>0919</u> | <u>21.28</u> | <u>6.43</u> | <u>50.8</u> | <u>1.25</u> | <u>11.8</u> | <u>41.3</u> | <u>24</u> | |
| <u>0923</u> | <u>21.30</u> | <u>6.45</u> | <u>51.2</u> | <u>0.64</u> | <u>12.3</u> | <u>12.0</u> | <u>17</u> | |
| <u>0927</u> | <u>21.30</u> | <u>6.48</u> | <u>51.2</u> | <u>0.39</u> | <u>12.4</u> | <u>5.9</u> | <u>12</u> | |
| <u>0931</u> | <u>21.30</u> | <u>6.44</u> | <u>51.5</u> | <u>0.30</u> | <u>12.3</u> | <u>6.4</u> | <u>9</u> | <u>5 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>10/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-7
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-7

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>10/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>26.56</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>10/29/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1131-1139</u> |
| Date/Time Sampled | <u>10/29/10 1145</u> | | <u>480 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|-------------|----------------|-------------|--------------------|
| <u>1121</u> | <u>26.61</u> | <u>6.92</u> | <u>16.8</u> | <u>6.16</u> | <u>10.9</u> | <u>64.0</u> | <u>219</u> | |
| <u>1127</u> | <u>26.59</u> | <u>6.79</u> | <u>15.5</u> | <u>5.67</u> | <u>11.6</u> | <u>13.0</u> | <u>164</u> | |
| <u>1131</u> | <u>26.59</u> | <u>6.68</u> | <u>15.4</u> | <u>5.57</u> | <u>12.3</u> | <u>10.4</u> | <u>142</u> | |
| <u>1135</u> | <u>26.59</u> | <u>6.73</u> | <u>15.3</u> | <u>5.54</u> | <u>12.6</u> | <u>8.1</u> | <u>143</u> | |
| <u>1139</u> | <u>26.59</u> | <u>6.76</u> | <u>15.3</u> | <u>5.55</u> | <u>12.9</u> | <u>8.5</u> | <u>148</u> | <u>3 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>10/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature L. Linde Page _____ of _____

Parametrix, Inc.

Well #: MW-8
 Sample #: _____
OL-MW-8

Groundwater Sampling Field Data Sheet

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>10/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>21.12</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>10/29/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1015 - 1035</u> |
| Date/Time Sampled | <u>10/29/10 1040</u> | | <u>400 ml/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|-------------|----------------|-------------|--------------------|
| <u>1015</u> | <u>21.53</u> | <u>6.54</u> | <u>25.0</u> | <u>2.36</u> | <u>11.0</u> | <u>137</u> | <u>106</u> | |
| <u>1019</u> | <u>21.53</u> | <u>6.44</u> | <u>25.7</u> | <u>0.73</u> | <u>11.3</u> | <u>31.4</u> | <u>84</u> | |
| <u>1023</u> | <u>21.53</u> | <u>6.42</u> | <u>33.6</u> | <u>0.30</u> | <u>12.1</u> | <u>74.7</u> | <u>77</u> | |
| <u>1027</u> | <u>21.53</u> | <u>6.46</u> | <u>35.7</u> | <u>0.22</u> | <u>12.5</u> | <u>48.3</u> | <u>69</u> | |
| <u>1031</u> | <u>21.53</u> | <u>6.48</u> | <u>37.0</u> | <u>0.15</u> | <u>12.7</u> | <u>36.4</u> | <u>65</u> | |
| <u>1035</u> | <u>21.53</u> | <u>6.48</u> | <u>37.4</u> | <u>0.15</u> | <u>12.7</u> | <u>22.3</u> | <u>63</u> | <u>4 gal</u> |
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|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>10/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity Good
 Remarks _____
 Signature [Signature] Page 1 of 1

Parametrix, Inc.

Well #: MW-10
 Sample #: _____

Groundwater Sampling Field Data Sheet

OL-MW-10

| | | | |
|-----------------|------------------------------------|------------|----------------------------|
| Project Number | <u>215-1578-121</u> | Date | <u>10/29/10</u> |
| Project Name | <u>Olalla LF Qtrly GW Sampling</u> | Location | <u>Olalla LF</u> |
| Project Address | <u>Bandix Rd</u> | Sampled By | <u>L. Linde/C. Wiggins</u> |
| Client Name | <u>KC SWD</u> | Purged By | <u>L. Linde/C. Wiggins</u> |

Casing Diameter: 2" 4" 6" Other _____

| | | | |
|---------------------------------------|----------------------|-------------------------|----------------------------|
| Depth to Water (feet) | <u>31.26</u> | Purge Vol. Meas. Method | <u>Grad Cyl/Stop Watch</u> |
| Depth of Well (feet) | _____ | Date Purged | <u>10/29/10</u> |
| Reference Point (surveyors notch/etc) | <u>TOC</u> | Purge Time (from/to) | <u>1326 - 1946</u> |
| Date/Time Sampled | <u>10/29/10 1355</u> | | <u>450 mL/min</u> |

Purge Volume Calculation: $(\pi r^2 h)(7.48 \text{ gal/ft}^3)(\# \text{ Casing volumes})$
 Purge Volume (gallons) for 2" = $(0.16)(h)(\#Cv)$; 4" = $(0.653)(h)(\#Cv)$; 6" = $(1.48)(h)(\#Cv)$
 Calculated Purge Volume (gallons) _____ Actual Purge Volume (gallons) _____

| TIME (2400 hr) | WATER LEVEL (feet) | pH (units) ± 0.1 | COND (mS/cm) ± 3% | DO (mg/L) ± 10% | TEMP °C | TURB. ± 10% | ORP (mV) | CUM. VOL. (gal) |
|-------------------|--------------------------|------------------------|-------------------------|-----------------------|-------------|----------------|-------------|--------------------|
| <u>1326</u> | <u>31.33</u> | <u>6.52</u> | <u>49.7</u> | <u>2.19</u> | <u>10.8</u> | <u>24.9</u> | <u>234</u> | |
| <u>1330</u> | <u>31.33</u> | <u>6.46</u> | <u>50.0</u> | <u>0.80</u> | <u>11.7</u> | <u>16.4</u> | <u>170</u> | |
| <u>1334</u> | <u>31.33</u> | <u>6.47</u> | <u>49.8</u> | <u>2.12</u> | <u>12.0</u> | <u>10.2</u> | <u>153</u> | |
| <u>1338</u> | <u>31.34</u> | <u>6.51</u> | <u>49.7</u> | <u>0.69</u> | <u>12.1</u> | <u>9.5</u> | <u>145</u> | |
| <u>1942</u> | <u>31.33</u> | <u>6.53</u> | <u>49.8</u> | <u>0.53</u> | <u>12.1</u> | <u>10.0</u> | <u>140</u> | |
| <u>1946</u> | <u>31.33</u> | <u>6.52</u> | <u>49.9</u> | <u>0.46</u> | <u>12.1</u> | <u>9.5</u> | <u>135</u> | <u>4.5 gal</u> |
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| | | | |
|---------------------------|-----------------------------|-----------------------------------|-----------------------------|
| Purge Equipment | <u>Grundfos Ready-flo 2</u> | Sampling Equipment | <u>Grundfos Ready-flo 2</u> |
| Laboratory | <u>Aquatic Research</u> | Date Sent to Lab | <u>10/29/10</u> |
| Chain-of-Custody (yes/no) | <u>Yes</u> | Field QC Sample Number | <u>N/A</u> |
| Shipment Method | <u>Hand Delivered</u> | Split with (name(s)/organization) | <u>N/A</u> |

Well Integrity: Good
 Remarks: Dedicated Grundfos not working, pull for repair
 Signature: [Signature] Page _____ of _____

10/15/10 52°F

17

OLUCA

Overcast then P.C

| Loc | P | CH ₄ /O ₂ | CO ₂ | T | I/E |
|------|---|---------------------------------|-----------------|----|--------|
| FL-1 | φ | φ/20.1 | 0.8 | 58 | -29/φ |
| FL-2 | φ | 12/0.8 | 13.4 | 55 | 730°/0 |
| FL-3 | φ | 0.4/17.8 | 2.7 | 57 | 730°/ |

- Verified canister vacuums
- Monitored static pressure until stable
- Pumped gas stream until gas concentrations stabilized
- Connected canister and opened until flow stopped
- Rebuilt canister pressure
- Labeled canister + packaged
- Fed ex sample

Shallaf RIFC

10/25/10 ²³

Activities: Collect gw samples

Weather: Rainy, 48°F

Organize: L. Linde, C. Wiggins

Author: L. Linde

0705 Depart

0735 Stop to get ice & paper towels

0745 Back on the road

0800 Arrive C dump box gate, wait

0810 meet w/ Chad, go to bagg gate

0840 Calibrate Horiba - good

0850 Begin purging MW-1

0930 Collect sample @ MW-1, take
30 min to collect enough volume

1000 Pack up, go to MW-2, decon
rental pump, carry gear to MW-2

1102 Begin purging MW-2

1145 Collect sample @ MW-2

1200 Decon pump

1215 Pack gear to MW-4, setup

1245 Begin purging MW-4

1325 Collect sample @ MW-4

1345 Pack gear out back to cars

1350 Decon pump

1415 Go to MW-3 & setup

Olalla LF RI/FS 10/28/10

- 1432 Begin purging MW-5A
 1510 Collect sample @ MW-5A
 1520 Pack up gear, move to MW-3
 1530 Setup @ MW-3
 1533 Begin purging MW-3
 1600 Collect sample @ MW-5A
 1615 Collect field duplicate @ MW-5A, labeled as MW-9
 1630 Rearrange gear, will meet w/ Mike in the morning to hand off samples, containerize ^{Purge} water
 1647 Finish collecting water levels @ remainder of wells - far potentiometric surface

| Water Level Table | Well ID | DTW |
|-------------------|---------|-------|
| | MW-1 | 79.59 |
| | MW-2 | 66.63 |
| | MW-3 | 46.80 |
| | MW-4 | 64.09 |
| | MW-5A | 77.83 |
| | MW-6 | 21.07 |
| | MW-7 | 26.58 |
| | MW-8 | 21.37 |
| | MW-10 | 31.21 |

Olalla LF RI/FS 10/28/10

- Late Entry 0900 Cannot maintain a constant flow
 Late Entry 0915 Receive call from Damien @ Aquatic Research, confirm hold time, will call back tomorrow w/ estimated time to finish
 Late Entry 0920 Call Mike W., Chad can take samples for delivery today but might not get enough wells done, will call back @ 1400 to discuss progress
 Late Entry 1405 Call Mike W. back to update status - completed 3 wells, can do 2 more if Chad stays, Mike will meet w/ me in AM to hand off samples, containerize ^{Purge} water in drum
 1650 Put coolers w/ my gear, Chad to take portable Grundfos to Bramerton office for shipping & P/U an empty cooler
 1715 Will meet @ site @ 0800 tomorrow morning w/ Chad

Alalla LF R/FS 10/28/10
 1717 Depart site, secure gate
 1905 Call Mike, will meet
 @ TORO @ 0730 tomorrow
 morning to hand off samples
 to satisfy confirm holding
 time requirement

Scott (Pmk) 10/28/10

Alalla LF R/FS 10/29/10
 Activities: Finish gw sample collection
 Weather: Sunny 55°F
 Onsite: L. Linde, C. Wiggins
 Author: L. Linde

0700 Depart for TORO
 0725 Arrive @ TORO, drop off
 samples w/ Mike to take
 up this morning
 0740 Depart TORO
 0750 Stop for ice
 0755 Depart for site
 0810 Arrive @ site, Chad (Pmk)
 just arrived before me
 0820 At MW-10, Set up
 0830 Portable Grundfos ^{10/29/10}
 Dedicated Grundfos (new)
 not working, keeps giving
 "HW Ground Fault", trouble
 shoot
 0845 Go to MW-6 to test
 control box, pump & box
 works, move gear to MW-6
 0900 Call Scott (Pmk), ask to
 bring portable Grundfos

- Olalla LF R/FS 10/29/10
- 0915 Calibrate Horiba - good,
begin purging MW-6
- 0940 Collect sample @ MW-6
- 0945 Scott (PMX) onsite, deliver
pump & extra tubing to
sample MW-10 in case
we still can't get control
box to work
- 0955 Scott offsite
- 1000 Dump purge water from
MW-6 in MW-8 develop.
water drum
- 1005 Setup @ MW-8
- 1015 Begin purging MW-8
- 1040 Collect sample @ MW-8
Discard purge water in
MW-8 development water
drum
- 1100 Setup @ MW-7
- ~~1050~~ ~~10/29/10~~
- Late Entry 1050 Ford Ranger
battery dead, Highlander
provides jump start
- 1121 Begin purging MW-7
- 1145 Collect sample @ MW-7

- Olalla LF R/FS 10/29/10
- 1150 Dump purge water from
MW-7 in drainage ditch
beside well, not required
to containerize @ this
location
- 1210 Back @ MW-10, setup
again, attempt to reconnect
control box to dedicated
pump w/o "ground fault"
mechanism. Attempts failed
- 1215 Contact INW to field
troubleshoot connections,
talk to Tyler - will get
his tech guy to call back
right away, take quick
break for lunch
- 1235 Call Tyler back, says tech
guy is on lunch & will
call back as soon as he
gets back
- 1240 Receive vm from Shawn
@ INW to call back, was
on the phone w/ Tyler
when he called
- 1250 Leave vm for Shawn to

- Alalla LF RI/FS 10/29/10
 call back ASAP, waiting in
 the field
- 1255 Receive call from Shawn @
 INW, field troubleshooting,
 appears that connection
 cannot be fixed in the field
- 1305 Pull dedicated pump in MW-10
 & coil for repair/service
- 1310 Deploy rental pump in
 well, was previously decon
 yesterday after last use
- 1326 Begin phasing MW-10
- 1355 Collect sample @ MW-10
- 1410 Decon pump & clean up,
 dump purge water in
 develop. water drum for
 MW-10
- 1430 Redistribute rear
- 1445 Call Mike W. (PMX), inform
 him of completion & difficulty
 with MW-10, approves rental
 pump return & dedicated
 pump return to INW on
 Mon (11/1) w/ Chad.
- 1450 Call Aquatic Research &

- Alalla LF RI/FS 10/29/10
 inform them samples will
 be arriving a little before
 1700 due to field difficulties
 Give Chad their # to call
 in case he will be late
- 1500 Call Keli (Kc) to see
 if I can drop off pump
 controller, no answer,
 will go to office & return it
 on Mon (11/1)
- 1505 Secure gate, Chad &
 I depart from site.
- 1525 Arrive in Bremerton
- 1530 Unload gear

Handwritten signature and date: 10/29/10



Kitsap County Public Works
Solid Waste Division
 614 Division Street (MS-27)
 Port Orchard, WA 98366
 Phone: 360-337-5777
 FAX: 360-337-4867

CHAIN OF CUSTODY RECORD

FOR LAB USE ONLY

| | |
|----------------------------------|---|
| LOGIN COMMENTS: | Sample(s) checked/logged by: _____ |
| | Sample temp on receipt at lab: _____ COMPLIANCE 4 +/- 4° C |
| | Are custody seals intact? <input type="checkbox"/> YES <input type="checkbox"/> NO |
| Sample receipt (day/time): _____ | Ice? <input type="checkbox"/> NONE <input type="checkbox"/> BLUE <input type="checkbox"/> WET |

TO BE COMPLETED BY THE SAMPLER:

TAT REQUESTED: STANDARD: OTHER: _____ PROJECT: Olalla LF Lara Linde SAMPLER PRINTED NAME AND SIGNATURE

| SAMPLE DATE | SAMPLE TIME | SITE NAME OR LOCATION | SAMPLE IDENTIFIER | MATRIX * | GRAB | COMP | ANALYSES REQUIRED — MARK AN X IN ALL TESTS REQUIRED FOR EACH SAMPLE LINE | | | | | | | |
|-------------|-------------|-----------------------|-------------------|----------|------|------|--|------|-----|--------------|----------|----|---------------|---------|
| | | | | | | | VOCs, VC | PEBs | ON- | Conventional | CO2, TOC | pH | T.F. Coliform | T. M115 |
| 10/29 | 1355 | MW-10 | DL-MW-10 | GW | X | | X | X | X | X | X | X | X | X |
| 10/29 | 0740 | MW-6 | DL-MW-6 | | | | | | | | | | | |
| 10/29 | 1145 | MW-7 | DL-MW-7 | | | | | | | | | | | |
| 10/29 | 1040 | MW-8 | DL-MW-8 | | | | | | | | | | | |
| 10/29 | 0900 | Trip Blank | TB-2 | | | | | | | | | | | |

* MATRIX TYPES: SW = Surface Water GW = Ground Water SO = Soil * Field Filtered

| SIGNATURE | PRINTED NAME | COMPANY AND TITLE | DATE | TIME |
|---------------------|--------------|-------------------|----------|------|
| <i>Lara Linde</i> | Lara Linde | Parametric | 10/29/10 | 1450 |
| <i>Chad Wiggins</i> | Chad Wiggins | Parametric | 10/29/10 | 1450 |
| <i>Chad Wiggins</i> | Chad Wiggins | Parametric | 10/29/10 | 1640 |
| <i>S. Hillson</i> | S. HILLSON | Anal | 10/29/10 | 1640 |
| | | | | |
| | | | | |

DISTRIBUTION: WHITE-With lab report YELLOW-KCPW PINK-Lab

KCPW 3053-REV 05/02

Conventional = Alkalinity, Ammonia, Carbonate, Bicarbonate, Chloride, Nitrate, Nitrite, Sulfate

Appendix G

Appendix II and Appendix III List of Constituents
from WAC 173-351-990

ested parties including the county or city having jurisdiction over the site and the department. Within thirty days after the hearing the health officer shall notify the applicant or the holder of the permit in writing of his determination thereof. Any party aggrieved by such determination may appeal to the pollution control hearings board by filing with the hearings board a notice of appeal within thirty days after receipt of notice of the determination of the health officer. The hearings board shall hold a hearing in accordance with the provisions of the Administrative Procedure Act, chapter 34.05 RCW, as now or hereafter amended.

[Statutory Authority: Chapter 70.95 RCW and 40 CFR 258.93-22-016, § 173-351-760, filed 10/26/93, effective 11/26/93.]

WAC 173-351-990 Appendices.

APPENDIX I¹

Appendix I - Constituents for Detection Monitoring

| COMMON NAME ² | CAS RN ³ |
|---|---------------------|
| Inorganic Constituents | |
| 1) Antimony | (Dissolved) |
| 2) Arsenic | (Dissolved) |
| 3) Barium | (Dissolved) |
| 4) Beryllium | (Dissolved) |
| 5) Cadmium | (Dissolved) |
| 6) Chromium | (Dissolved) |
| 7) Cobalt | (Dissolved) |
| 8) Copper | (Dissolved) |
| 9) Lead | (Dissolved) |
| 10) Nickel | (Dissolved) |
| 11) Selenium | (Dissolved) |
| 12) Silver | (Dissolved) |
| 13) Thallium | (Dissolved) |
| 14) Vanadium | (Dissolved) |
| 15) Zinc | (Dissolved) |
| 16) Nitrate | |
| Organic Constituents | |
| 17) Acetone | 67-64-1 |
| 18) Acrylonitrile | 107-13-1 |
| 19) Benzene | 71-43-2 |
| 20) Bromochloromethane | 74-97-5 |
| 21) Bromodichloromethane | 75-27-4 |
| 22) Bromoform; Tribromomethane | 75-25-2 |
| 23) Carbon disulfide | 75-15-0 |
| 24) Carbon tetrachloride | 56-23-5 |
| 25) Chlorobenzene | 108-90-7 |
| 26) Chloroethane; Ethyl chloride | 75-00-3 |
| 27) Chloroform; Trichloromethane | 67-66-3 |
| 28) Dibromochloromethane; Chlorodibromomethane | 124-48-1 |
| 29) 1,2-Dibromo-3-chloropropane; DBCP | 96-12-8 |
| 30) 1,2-Dibromoethane; Ethylene dibromide; EDB | 106-93-4 |
| 31) o-Dichlorobenzene; 1,2-Dichlorobenzene | 95-50-1 |
| 32) p-Dichlorobenzene; 1,4-Dichlorobenzene | 106-46-7 |

| COMMON NAME ² | CAS RN ³ |
|---|---------------------|
| Inorganic Constituents | |
| 33) trans-1,4-Dichloro-2-butene | 110-57-6 |
| 34) 1,1-Dichloroethane; Ethylidene chloride | 75-34-3 |
| 35) 1,2-Dichloroethane; Ethylene dichloride | 107-06-2 |
| 36) 1,1-Dichloroethylene; 1,1-Dichloroethene; Vinylidene chloride | 75-35-4 |
| 37) cis-1,2-Dichloroethylene; cis-1,2-Dichloroethene | 156-59-2 |
| 38) trans-1,2-Dichloroethylene; trans-1,2-Dichloroethene | 156-60-5 |
| 39) 1,2-Dichloropropane; Propylene dichloride | 78-87-5 |
| 40) cis-1,3-Dichloropropene | 10061-01-5 |
| 41) trans-1,3-Dichloropropene | 10061-02-6 |
| 42) Ethylbenzene | 100-41-4 |
| 43) 2-Hexanone; Methyl butyl ketone | 591-73-6 |
| 44) Methyl bromide; Bromomethane | 74-83-9 |
| 45) Methyl chloride; Chloromethane | 74-87-3 |
| 46) Methylene bromide; Dibromomethane | 74-95-3 |
| 47) Methylene chloride; Dichloromethane | 75-09-2 |
| 48) Methyl ethyl ketone; MEK; 2-Butanone | 78-93-3 |
| 49) Methyl iodide; Iodomethane | 74-88-4 |
| 50) 4-Methyl-2-pentanone; Methyl isobutyl ketone | 108-10-1 |
| 51) Styrene | 100-42-5 |
| 52) 1,1,1,2-Tetrachloroethane | 630-20-6 |
| 53) 1,1,2,2-Tetrachloroethane | 79-34-5 |
| 54) Tetrachloroethylene; Tetrachloroethene; Perchloroethylene | 127-18-4 |
| 55) Toluene | 108-88-3 |
| 56) 1,1,1-Trichloroethane; Methyl chloroform | 71-55-6 |
| 57) 1,1,2-Trichloroethane | 79-00-5 |
| 58) Trichloroethylene; Trichloroethene | 79-01-6 |
| 59) Trichlorofluoromethane; CFC-11 | 75-69-4 |
| 60) 1,2,3-Trichloropropane | 96-18-4 |
| 61) Vinyl acetate | 108-05-4 |
| 62) vinyl chloride | 75-01-4 |
| 63) Xylenes | 1330-20-7 |

¹This list contains 47 volatile organics for which possible analytical procedures provided in EPA Report SW-846 "Test Methods for Evaluating Solid Waste," third edition, November 1986, as revised December 1987, includes Method 8260; and 15 metals for which SW-846 provides either Method 6010 or a method from the 7000 series of methods.

² Common names are those widely used in government regulations, scientific publications, and commerce; synonyms exist for many chemicals.

³ Chemical Abstracts Service registry number.

APPENDIX II

GROUND WATER QUALITY PARAMETERS

Field Parameters

pH

specific conductance
 temperature
 static water level

Geochemical Indicator Parameters

Calcium (Ca) Sodium (Na)
 Bicarbonate (HCO₃) Chloride (Cl)
 Magnesium (Mg) Potassium (K)
 Sulfate (SO₄) Alkalinity (as Ca CO₃)

Iron (Fe)
 Manganese (Mn)

Leachate Indicators

Ammonia (NH₃-N)
 Total Organic Carbon (TOC)
 Total Dissolved Solids (TDS)

APPENDIX III**List of Hazardous Inorganic and Organic Constituents.¹**

| Common Name² (mg/L)⁶ | CAS RN³ | Chemical abstracts service index name⁴ | Suggested methods⁵ | PQL |
|---|---------------------------|---|--|-------------------|
| Acenaphthene | 83-32-9 | Acenaphthylene, 1,2-dihydro- | 8100 8270 | 200 10 |
| Acenaphthylene | 208-96-8 | Acenaphthylene | 8100 8270 | 200 10 |
| Acetone | 67-64-1 | 2-Propanone | 8260 | 100 |
| Acetonitrile; Methyl cyanide | 75-05-8 | Acetonitrile | 8015 | 100 |
| Acetophenone | 98-86-2 | Ethanone, 1-phenyl- | 8270 | 10 |
| 2-Acetylaminofluorene; 2-AAF | 53-96-3 | Acetamide, N-9H-fluoren-2-yl- | 8270 | 20 |
| Acrolein | 107-02-8 | 2-Propenal | 8030 8260 | 5 100 |
| Acrylonitrile | 107-13-1 | 2-Propenenitrile | 8030 8260 | 5 200 |
| Aldrin | 309-00-2 | 1,4:5,8-Dimethanonaphthalene, 1,2,3,4,10,10-hexachloro-1,4, 4a,5,8,8a-hexahydro- (1 α ,4 α , 4a β ,5 α ,8 α ,8a β)- | 8080 8270 | 0.05 10 |
| Allyl chloride | 107-05-1 | 1-Propene, 3-chloro- | 8010 8260 | 5 10 |
| 4-Aminobiphenyl | 92-67-1 | [1,1 1 -Biphenyl]-4-amine | 8270 | 20 |
| Anthracene | 120-12-7 | Anthracene | 8100 8270 | 200 10 |
| Antimony | (Dissolved) | Antimony | 6010 7040 7041 | 300 2000 30 |
| Arsenic | (Dissolved) | Arsenic | 6010 7060 7061 | 500 10 20 |
| Barium | (Dissolved) | Barium | 6010 7080 | 20 1000 |
| Benzene | 71-43-2 | Benzene | 8020 8021 8260 | 2 0.1 5 |
| Benzo[a]anthracene; Benzanthracene | 56-55-3 | Benz[a]anthracene | 8100 8270 | 200 10 |

Municipal Solid Waste Landfills

173-351-990

| Common Name ² (mg/L) ⁶ | CAS RN ³ | Chemical abstracts service index name ⁴ | Suggested methods ⁵ | PQL |
|--|---------------------|--|-----------------------------------|----------------|
| Benzo[b]fluoranthene | 205-99-2 | Benz[e]acephenanthrylene | 8100 8270 | 200 10 |
| Benzo[k]fluoranthene | 207-08-9 | Benzo[k]fluoranthene | 8100 8270 | 200 10 |
| Benzo[ghi]perylene | 191-24-2 | Benzo[ghi]perylene | 8100 8270 | 200 10 |
| Benzo[a]pyrene | 50-32-8 | Benzo[a]pyrene | 8100 8270 | 200 10 |
| Benzyl alcohol | 100-51-6 | Benzenemethanol | 8270 | 20 |
| Beryllium | (Dissolved) | Beryllium | 6010 7090 7091 | 3 50 2 |
| alpha-BHC | 319-84-6 | Cyclohexane, 1,2,3,4,5,6- hexachloro-, (1 α ,2 α ,3 β ,4 α ,5 β ,6 β)- | 8080 8270 | 0.05 10 |
| beta-BHC | 319-85-7 | Cyclohexane, 1,2,3,4,5,6- hexachloro-, (1 α ,2 β ,3 α ,4 β ,5 α ,6 β)- | 8080 8270 | 0.05 20 |
| delta-BHC | 319-86-8 | Cyclohexane, 1,2,3,4,5,6- hexachloro-, (1 α ,2 α ,3 α ,4 β ,5 α ,6 β)- | 8080 8270 | 0.1 20 |
| gamma-BHC; Lindane | 58-89-9 | Cyclohexane, 1,2,3,4,5,6- hexachloro-, (1 α ,2 α ,3 β ,4 α ,5 α ,6 β)- | 8080 8270 | 0.05 20 |
| Bis(2-chloroethoxy)methane | 111-91-1 | Ethane, 1,1 1 - [methylenebis(oxy)]bis[2-chloro- | 8110 8270 | 5 10 |
| Bis(2-chloroethyl) ether; Dichloroethyl ether | 111-44-4 | Ethane, 1,1 1 -oxybis[2-chloro- | 8110 8270 | 3 10 |
| Bis-(2-chloro-1-methylethyl) ether; 2,2 1 - Dichlorodiisopropyl ether; DCIP, See note 7 | 108-60-1 | Propane, 2,2 1 -oxybis[1-chloro- | 8110 8270 | 10 10 |
| Bis(2-ethylhexyl) phthalate | 117-81-7 | 1,2-Benzenedicarboxylic acid, bis(2-ethylhexyl) ester | 8060 | 20 |
| Bromochloromethane; Chlorobromomethane | 74-97-5 | Methane, bromochloro- | 8021 8260 | 0.1 5 |
| Bromodichloromethane; Dibromochloromethane | 75-27-4 | Methane, bromodichloro- | 8010 8021 8260 | 1 0.2 5 |
| Bromoform; Tribromomethane | 75-25-2 | Methane, tribromo- | 8010 8021 8260 | 2 15 5 |
| 4-Bromophenyl phenyl ether | 101-55-3 | Benzene, 1-bromo-4-phenoxy- | 8110 8270 | 25 10 |
| Butyl benzyl phthalate; Benzyl butyl phthalate | 85-68-7 | 1,2-Benzenedicarboxylic acid, butyl phenylmethyl ester | 8060 8270 | 5 10 |
| Cadmium | (Dissolved) | Cadmium | 6010 7130 7131 | 40 50 1 |
| Carbon disulfide | 75-15-0 | Carbon disulfide | 8260 | 100 |
| Carbon tetrachloride | 56-23-5 | Methane, tetrachloro- | 8010 8021 8260 | 1 0.1 10 |
| Chlordane | See Note 8 | 4,7-Methano-1H-indene, 1,2,4,5, 6,7,8,8-octachloro-2,3,3a,4,7, 7a-hexahydro- | 8080 8270 | 0.1 50 |
| p-Chloroaniline | 106-47-8 | Benzenamine, 4-chloro- | 8270 | 20 |

| Common Name ² (mg/L) ⁶ | CAS RN ³ | Chemical abstracts service index name ⁴ | Suggested methods ⁵ | PQL |
|---|---------------------|--|-----------------------------------|------|
| Chlorobenzene | 108-90-7 | Benzene, chloro- | 8010 | 2 |
| | | | 8020 | 2 |
| | | | 8021 | 0.1 |
| | | | 8260 | 5 |
| Chlorobenzilate | 510-15-6 | Benzeneacetic acid, 4-chloro- α - (4-chlorophenyl)- α -hydroxy-, ethyl ester | 8270 | 10 |
| p-Chloro-m-cresol; 4-Chloro-3- methylphenol | 59-50-7 | Phenol, 4-chloro-3-methyl- | 8040 | 5 |
| | | | 8270 | 20 |
| Chloroethane; Ethyl chloride | 75-00-3 | Ethane, chloro- | 8010 | 5 |
| | | | 8021 | 1 |
| | | | 8260 | 10 |
| Chloroform; Trichloromethane | 67-66-3 | Methane, trichloro- | 8010 | 0.5 |
| | | | 8021 | 0.2 |
| | | | 8260 | 5 |
| 2-Chloronaphthalene | 91-58-7 | Naphthalene, 2-chloro- | 8120 | 10 |
| | | | 8270 | 10 |
| 2-Chlorophenol | 95-57-8 | Phenol, 2-chloro- | 8040 | 5 |
| | | | 8270 | 10 |
| 4-Chlorophenyl phenyl ether | 7005-72-3 | Benzene, 1-chloro-4-phenoxy- | 8110 | 40 |
| | | | 8270 | 10 |
| Chloroprene | 126-99-8 | 1,3-Butadiene, 2-chloro- | 8010 | 50 |
| | | | 8260 | 20 |
| Chromium | (Dissolved) | Chromium | 6010 | 70 |
| | | | 7190 | 500 |
| | | | 7191 | 10 |
| Chrysene | 218-01-9 | Chrysene | 8100 | 200 |
| | | | 8270 | 10 |
| Cobalt | (Dissolved) | Cobalt | 6010 | 70 |
| | | | 7200 | 500 |
| | | | 7201 | 10 |
| Copper | (Dissolved) | Copper | 6010 | 60 |
| | | | 7210 | 200 |
| | | | 7211 | 10 |
| | | | 8270 | 10 |
| m-Cresol; 3-methylphenol | 108-39-4 | Phenol, 3-methyl- | 8270 | 10 |
| o-Cresol; 2-methylphenol | 95-48-7 | Phenol, 2-methyl- | 8270 | 10 |
| p-Cresol; 4-methylphenol | 106-44-5 | Phenol, 4-methyl- | 8270 | 10 |
| Cyanide | 57-12-5 | Cyanide | 9010 | 200 |
| 2,4-D; 2,4- Dichlorophenoxyacetic acid | 94-75-7 | Acetic acid, (2,4- dichlorophenoxy)- | 8150 | 10 |
| 4,4 1 -DDD | 72-54-8 | Benzene 1,1 1 -(2,2- dichloroethylidene)bis[4- chloro- | 8080 | 0.1 |
| | | | 8270 | 10 |
| 4,4 1 -DDE | 72-55-9 | Benzene, 1,1 1 - (dichloroethylenylidene)bis[4- chloro- | 8080 | 0.05 |
| | | | 8270 | 10 |
| 4,4 1 -DDT | 50-29-3 | Benzene, 1,1 1 -(2,2,2- trichloroethylidene)bis[4- chloro- | 8080 | 0.1 |
| | | | 8270 | 10 |
| Diallate | 2303-16-4 | Carbamothioic acid, bis(1- methylethyl)-,S-(2,3-dichloro- 2-propenyl) ester | 8270 | 10 |
| Dibenz[a,h]anthracene | 53-70-3 | Dibenz[a,h]anthracene | 8100 | 200 |
| | | | 8270 | 10 |
| Dibenzofuran | 132-64-9 | Dibenzofuran | 8270 | 10 |

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| Common Name ² (mg/L) ⁶ | CAS RN ³ | Chemical abstracts service index name ⁴ | Suggested methods ⁵ | PQL |
|--|---------------------|---|-----------------------------------|------|
| Dibromochloromethane; Chlorodibromomethane | 124-48-1 | Methane, dibromochloro- | 8010 | 1 |
| | | | 8021 | 0.3 |
| | | | 8260 | 5 |
| 1,2-Dibromo-3-chloropropane; DBCP | 96-12-8 | Propane, 1,2-dibromo-3-chloro- | 8011 | 0.1 |
| | | | 8021 | 30 |
| | | | 8260 | 25 |
| | | | 8011 | 0.1 |
| 1,2-Dibromoethane; Ethylene dibromide; EDB | 106-93-4 | Ethane, 1,2-dibromo- | 8021 | 10 |
| | | | 8260 | 5 |
| | | | 8060 | 5 |
| Di-n-butyl phthalate | 84-74-2 | 1,2-Benzenedicarboxylic acid, dibutyl ester | 8270 | 10 |
| | | | 8010 | 2 |
| o-Dichlorobenzene; 1,2- Dichlorobenzene | 95-50-1 | Benzene, 1,2-dichloro- | 8020 | 5 |
| | | | 8021 | 0.5 |
| | | | 8120 | 10 |
| | | | 8260 | 5 |
| | | | 8270 | 10 |
| | | | 8010 | 5 |
| | | | 8020 | 5 |
| m-Dichlorobenzene; 1,3- Dichlorobenzene | 541-73-1 | Benzene, 1,3-Dichloro- | 8021 | 0.2 |
| | | | 8120 | 10 |
| | | | 8260 | 5 |
| | | | 8270 | 10 |
| | | | 8010 | 2 |
| p-Dichlorobenzene; 1,4- Dichlorobenzene | 106-46-7 | Benzene, 1,4-dichloro- | 8020 | 5 |
| | | | 8021 | 0.1 |
| | | | 8120 | 15 |
| | | | 8260 | 5 |
| | | | 8270 | 10 |
| 3,3'-Dichlorobenzidine | 91-94-1 | [1,1'-Biphenyl]-4,4'-diamine, 3,3'- dichloro- | 8270 | 20 |
| | | | 8260 | 100 |
| trans-1,4-Dichloro-2-butene Dichlorodifluoromethane; CFC 12; | 110-57-6 | 2-Butene, 1,4-dichloro-, (E)- | 8021 | 0.5 |
| | 75-71-8 | Methane, dichlorodifluoro- | 8260 | 5 |
| | 75-34-3 | Ethane, 1,1-dichloro- | 8010 | 1 |
| 1,1-Dichloroethane; Ethylidene chloride | 107-06-2 | Ethane, 1,1-dichloro- | 8021 | 0.5 |
| | | | 8260 | 5 |
| | | | 8010 | 0.5 |
| 1,1-Dichloroethylene; 1,1- Dichloroethene; Vinylidene chloride | 75-35-4 | Ethene, 1,1-dichloro- | 8021 | 0.5 |
| | | | 8260 | 5 |
| | | | 8010 | 1 |
| cis-1,2-Dichloroethylene; cis- 1,2-Dichloroethene | 156-59-2 | Ethene, 1,2-dichloro-, (Z)- | 8021 | 0.2 |
| | | | 8260 | 5 |
| trans-1,2-Dichloroethylene trans-1,2-Dichloroethene | 156-60-5 | Ethene, 1,2-dichloro-, (E)- | 8010 | 1 |
| | | | 8021 | 0.5 |
| | | | 8260 | 5 |
| 2,4-Dichlorophenol | 120-83-2 | Phenol, 2,4-dichloro- | 8040 | 5 |
| | | | 8270 | 10 |
| 2,6-Dichlorophenol 1,2-Dichloropropane; Propylene dichloride | 87-65-0 | Phenol, 2,6-dichloro- Propane, 1,2-dichloro- | 8270 | 10 |
| | | | 8010 | 0.5 |
| | | | 8021 | 0.05 |
| 1,3-Dichloropropane; Trimethylene dichloride | 142-28-9 | Propane, 1,3-dichloro- | 8260 | 5 |
| | | | 8021 | 0.3 |
| | | | 8260 | 5 |

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|--|---------------------|--|-----------------------------------|------|
| 2,2-Dichloropropane; Isopropylidene chloride | 594-20-7 | Propane, 2,2-dichloro- | 8021 | 0.5 |
| | | | 8260 | 15 |
| 1,1-Dichloropropene | 563-58-6 | 1-Propene, 1,1-dichloro- | 8021 | 0.2 |
| | | | 8260 | 5 |
| cis-1,3-Dichloropropene | 10061-01-5 | 1-Propene, 1,3-dichloro-, (Z)- | 8010 | 20 |
| | | | 8260 | 10 |
| trans-1,3-Dichloropropene | 10061-02-6 | 1-Propene, 1,3-dichloro-, (E)- | 8010 | 5 |
| | | | 8260 | 10 |
| Dieldrin | 60-57-1 | 2,7:3,6-Dimethanonaphth[2,3-b]oxirene, 3,4,5,6,9,9-hexa, chloro-1a,2,2a,3,6,6a,7,7a-octahydro-, (1 α ,2 β ,2 α ,3 β ,6 β ,6 α ,7 β ,7 α)- | 8080 | 0.05 |
| | | | 8270 | 10 |
| Diethyl phthalate | 84-66-2 | 1,2-Benzenedicarboxylic acid, diethyl ester | 8060 | 5 |
| | | | 8270 | 10 |
| 0,0-Diethyl 0-2-pyrazinyl phosphorothioate; Thionazin Dimethoate | 297-97-2 | Phosphorothioic acid, 0,0-diethyl 0-pyrazinyl ester | 8141 | 5 |
| | | | 8270 | 20 |
| p-(Dimethylamino)azobenzene | 60-51-5 | Phosphorodithioic acid, 0,0-dimethyl S-[2-(methylamino)-2-oxoethyl] ester | 8141 | 3 |
| | | | 8270 | 20 |
| 7,12-Dimethylbenz[a]anthracene | 60-11-7 | Benzenamine, N,N-dimethyl-4-(phenylazo)- | 8270 | 10 |
| | | | 8270 | 10 |
| 3,3'-Dimethylbenzidine | 57-97-6 | Benz[a]anthracene, 7,12-dimethyl-[1,1'-biphenyl]-4,4'-diamine, 3,3'-dimethyl- | 8270 | 10 |
| | | | 8270 | 10 |
| 2,4-Dimethylphenol; m-Xylenol | 119-93-7 | Phenol, 2,4-dimethyl- | 8040 | 5 |
| | | | 8270 | 10 |
| Dimethyl phthalate | 105-67-9 | 1,2-Benzenedicarboxylic acid, dimethyl ester | 8060 | 5 |
| | | | 8270 | 10 |
| m-Dinitrobenzene | 99-65-0 | Benzene, 1,3-dinitro- | 8270 | 20 |
| | | | 8270 | 20 |
| 4,6-Dinitro-o-cresol 4,6-Dinitro-2-methylphenol | 534-52-1 | Phenol, 2-methyl-4,6-dinitro | 8040 | 150 |
| | | | 8270 | 50 |
| 2,4-Dinitrophenol; | 51-28-5 | Phenol, 2,4-dinitro- | 8040 | 150 |
| | | | 8270 | 50 |
| 2,4-Dinitrotoluene | 121-14-2 | Benzene, 1-methyl-2,4-dinitro- | 8090 | 0.2 |
| | | | 8270 | 10 |
| 2,6-Dinitrotoluene | 606-20-2 | Benzene, 2-methyl-1,3-dinitro- | 8090 | 0.1 |
| | | | 8270 | 10 |
| Dinoseb; DNBP; 2-sec-Butyl-4,6-dinitrophenol | 88-85-7 | Phenol, 2-(1-methylpropyl)-4,6-dinitro- | 8150 | 1 |
| | | | 8270 | 20 |
| Di-n-octyl phthalate | 117-84-0 | 1,2-Benzenedicarboxylic acid, dioctyl ester | 8060 | 30 |
| | | | 8270 | 10 |
| Diphenylamine | 122-39-4 | Benzenamine, N-phenyl- | 8270 | 10 |
| | | | 8270 | 10 |
| Disulfoton | 298-04-4 | Phosphorodithioic acid, 0,0-diethyl S-[2-(ethylthio)ethyl] ester | 8140 | 2 |
| | | | 8141 | 0.5 |
| Endosulfan I | 959-98-8 | 6,9-Methano-2,4,3-benzodioxathiepin, 6,7,8,9,10,10-hexachloro-1,5,5a,6,9,9a-hexahydro-, 3-oxide, | 8270 | 10 |
| | | | 8080 | 0.1 |
| Endosulfan II | 33213-65-9 | 6,9-Methano-2,4,3-benzodioxathiepin, 6,7,8,9,10,10-hexachloro-1,5,5a,6,9,9a-hexahydro-, 3-oxide, (3 α ,5 α ,6 β ,9 β ,9 α)- | 8270 | 20 |
| | | | 8080 | 0.05 |
| | | | 8270 | 20 |

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|---|------------------------|--|-----------------------------------|------|-----|
| Endosulfan sulfate | 1031-07-8 | 6,9-Methano-2,4,3-benzodioxathiepin, 6,7,8,9,10,10-hexa-chloro-1,5,5a,6,9,9a-hexahydro-,3-3-dioxide | 8080 | 0.5 | |
| | | | 8270 | 10 | |
| Endrin | 72-20-8 | 2,7:3,6-Dimethanonaphth[2,3-b]oxirene, 3,4,5,6,9,9-hexachloro-1a,2,2a,3,6,6a,7,7a-octahydro-, (1 α , 2 β ,2a β ,3 α ,6 α ,6a β ,7 β ,7a α)- | 8080 | 0.1 | |
| | | | 8270 | 20 | |
| Endrin aldehyde | 7421-93-4 | 1,2,4-Methenocyclopenta[cd]pentalene-5-carboxaldehyde, 2,2a,3,3,4,7-hexachlorodecahydro-, (1 α ,2 β ,2a β ,4 β ,4a β ,5 β ,6a β ,6b β ,7R*)- | 8080 | 0.2 | |
| | | | 8270 | 10 | |
| Ethylbenzene | 100-41-4 | Benzene, ethyl- | 8020 | 2 | |
| | | | 8221 | 0.05 | |
| | | | 8260 | 5 | |
| | | | 8015 | 5 | |
| Ethyl methacrylate | 97-63-2 | 2-Propenoic acid, 2-methyl-, ethyl ester | 8260 | 10 | |
| | | | 8270 | 10 | |
| | | | 8270 | 20 | |
| | | | 8270 | 20 | |
| Ethyl methanesulfonate | 62-50-0 | Methanesulfonic acid, ethyl ester | 8270 | 20 | |
| | Famphur 52-85-7 | Phosphorothioic acid, 0-[4-[(dimethylamino)sulfonyl]phenyl] 0,0-dimethyl ester | 8270 | 20 | |
| Fluoranthene | 206-44-0 | Fluoranthene | 8100 | 200 | |
| | | | 8270 | 10 | |
| Fluorene | 86-73-7 | 9H-Fluorene | 8100 | 200 | |
| | | | 8270 | 10 | |
| Heptachlor | 76-44-8 | 4,7-Methano-1H-indene, 1,4,5,6,7,8,8-heptachloro-3a,4,7,7a-tetrahydro- | 8080 | 0.05 | |
| | | | 8270 | 10 | |
| Heptachlor epoxide | 1024-57-3 | 2,5-Methano-2H-indeno[1,2-b]oxirene, 2,3,4,5,6,7,7-heptachloro-1a,1b,5,5a,6,6a-hexahydro-, (1 α , 1b β , 2 α , 5 α ,5a β , 6 β , 6a α) | 8080 | 1 | |
| | | | 8270 | 10 | |
| Hexachlorobenzene | 118-74-1 | Benzene, hexachloro- | 8120 | 0.5 | |
| | | | 8270 | 10 | |
| Hexachlorobutadiene | 87-68-3 | 1,3-Butadiene, 1,1,2,3,4,4-hexachloro- | 8021 | 0.5 | |
| | | | 8120 | 5 | |
| | | | 8260 | 10 | |
| | | | 8270 | 10 | |
| | | | 8270 | 10 | |
| Hexachlorocyclopentadiene | 77-47-4 | 1,3-Cyclopentadiene, 1,2,3,4,5,5-hexachloro- | 8120 | 5 | |
| | | | 8270 | 10 | |
| Hexachloroethane | 67-72-1 | Ethane, hexachloro- | 8120 | 0.5 | |
| | | | 8260 | 10 | |
| | | | 8270 | 10 | |
| Hexachloropropene | 1888-71-7 | 1-Propene, 1,1,2,3,3,3-hexachloro- | 8270 | 10 | |
| | | | 8270 | 10 | |
| 2-Hexanone; Methyl butyl ketone | 591-78-6 | 2-Hexanone | 8260 | 50 | |
| | Indeno(1,2,3-cd)pyrene | 193-39-5 | Indeno(1,2,3-cd)pyrene | 8100 | 200 |
| | | 8270 | 10 | | |
| Isobutyl alcohol | 78-83-1 | 1-Propanol, 2-methyl- | 8015 | 50 | |
| | | | 8240 | 100 | |

| Common Name ² (mg/L) ⁶ | CAS RN ³ | Chemical abstracts service index name ⁴ | Suggested methods ⁵ | PQL |
|---|---------------------|--|-----------------------------------|------|
| Isodrin | 465-73-6 | 1,4,5,8-Dimethanonaphthalene,1, 2,3,4,10,10- hexachloro-1,4,4a, 5,8,8a hexahydro- (1 α ,4 α ,4a β , 5 β ,8 β ,8a β)- | 8270 | 20 |
| | | | 8260 | 10 |
| Isophorone | 78-59-1 | 2-Cyclohexen-1-one, 3,5,5- trimethyl- | 8090 | 60 |
| Isosafrole | 120-58-1 | 1,3-Benzodioxole, 5-(1-propenyl)- | 8270 | 10 |
| Kepon | 143-50-0 | 1,3,4-Metheno-2H- cyclobuta[cd]pentalen-2-one, 1, 1a,3,3a,4,5,5a,5b,6- decachlorooctahydro- | 8270 | 20 |
| Lead | (Dissolved) | Lead | 6010 | 400 |
| | | | 7420 | 1000 |
| | | | 7421 | 10 |
| | | | 7470 | 2 |
| Mercury | (Total) | Mercury | 7470 | 2 |
| Methacrylonitrile | 126-98-7 | 2-Propenenitrile, 2-methyl- | 8015 | 5 |
| | | | 8260 | 100 |
| Methapyrilene | 91-80-5 | 1,2-Ethanediamine, N,N- dimethyl-N 1 -2-pyridinyl-N1/2- thienylmethyl)- | 8270 | 100 |
| | | | | |
| Methoxychlor | 72-43-5 | Benzene,1,1 1 -(2,2,2, trichloroethylidene)bis[4-methoxy- | 8080 | 2 |
| | | | 8270 | 10 |
| Methyl bromide; Bromomethane | 74-83-9 | Methane, bromo- | 8010 | 20 |
| | | | 8021 | 10 |
| Methyl chloride; Chloromethane | 74-87-3 | Methane, chloro- | 8010 | 1 |
| | | | 8021 | 0.3 |
| 3-Methylcholanthrene | 56-49-5 | Benz[j]aceanthrylene, 1,2-dihydro-3- methyl- | 8270 | 10 |
| | | | | |
| Methyl ethyl ketone; MEK; 2-Butanone | 78-93-3 | 2-Butanone | 8015 | 10 |
| | | | 8260 | 100 |
| Methyl iodide; Iodomethane | 74-88-4 | Methane, iodo- | 8010 | 40 |
| | | | 8260 | 10 |
| Methyl methacrylate | 80-62-6 | 2-Propenoic acid, 2-methyl-, methyl ester | 8015 | 2 |
| | | | 8260 | 30 |
| Methyl methanesulfonate | 66-27-3 | Methanesulfonic acid, methyl ester | 8270 | 10 |
| 2-Methylnaphthalene | 91-57-6 | Naphthalene, 2-methyl- | 8270 | 10 |
| Methyl parathion; Parathion methyl | 298-00-0 | Phosphorothioic acid, 0,0-dimethyl | 8140 | 0.5 |
| | | | 8141 | 1 |
| | | | 8270 | 10 |
| | | | 8015 | 5 |
| 4-Methyl-2-pentanone; Methyl isobutyl ketone | 108-10-1 | 2-Pentanone, 4-methyl- | 8260 | 100 |
| | | | 8010 | 15 |
| Methylene bromide; Dibromomethane | 74-95-3 | Methane, dibromo- | 8021 | 20 |
| | | | 8260 | 10 |
| Methylene chloride; Dichloromethane | 75-09-2 | Methane, dichloro- | 8010 | 5 |
| | | | 8021 | 0.2 |
| Naphthalene | 91-20-3 | Naphthalene | 8260 | 10 |
| | | | 8021 | 0.5 |
| 1,4-Naphthoquinone | 130-15-4 | 1,4-Naphthalenedione | 8100 | 200 |
| | | | 8260 | 5 |
| 1-Naphthylamine | 134-32-7 | 1-Naphthalenamine | 8270 | 10 |
| 2-Naphthylamine | 91-59-8 | 2-Naphthalenamine | 8270 | 10 |
| Nickel | (Total) | Nickel | 6010 | 150 |
| | | | 7520 | 400 |

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|---|---------------------|---|-----------------------------------|-----|
| o-Nitroaniline; 2-Nitroaniline | 88-74-4 | Benzenamine, 2-nitro- | 8270 | 50 |
| m-Nitroaniline; 3-Nitroaniline | 99-09-2 | Benzenamine, 3-nitro- | 8270 | 50 |
| p-Nitroaniline; 4-Nitroaniline | 100-01-6 | Benzenamine, 4-nitro- | 8270 | 20 |
| Nitrobenzene | 98-95-3 | Benzene, nitro- | 8090 | 40 |
| | | | 8270 | 10 |
| o-Nitrophenol; 2-Nitrophenol | 88-75-5 | Phenol, 2-nitro- | 8040 | 5 |
| | | | 8270 | 10 |
| p-Nitrophenol; 4-Nitrophenol | 100-02-7 | Phenol, 4-nitro- | 8040 | 10 |
| | | | 8270 | 50 |
| N-Nitrosodi-n-butylamine | 924-16-3 | 1-Butanamine, N-butyl-N-nitroso- | 8270 | 10 |
| N-Nitrosodiethylamine | 55-18-5 | Ethanamine, N-ethyl-N-nitroso- | 8270 | 20 |
| N-Nitrosodimethylamine | 62-75-9 | Methanamine, N-methyl-N-nitroso- | 8070 | 2 |
| N-Nitrosodiphenylamine | 86-30-6 | Benzenamine, N-nitroso-N-phenyl- | 8070 | 5 |
| N-Nitrosodipropylamine; N-Nitroso-N-dipropylamine; Di-n-propylnitrosamine | 621-64-7 | 1-Propanamine, N-nitroso-N-propyl- | 8070 | 10 |
| N-Nitrosomethylethylamine | 10595-95-6 | Ethanamine, N-methyl-N-nitroso- | 8270 | 10 |
| N-Nitrosopiperidine | 100-75-4 | Piperidine, 1-nitroso- | 8270 | 20 |
| N-Nitrosopyrrolidine | 930-55-2 | Pyrrolidine, 1-nitroso- | 8270 | 40 |
| 5-Nitro-o-toluidine | 99-55-8 | Benzenamine, 2-methyl-5-nitro- | 8270 | 10 |
| Parathion | 56-38-2 | Phosphorothioic acid, 0,0-diethyl 0-(4-nitrophenyl) ester | 8141 | 0.5 |
| | | | 8270 | 10 |
| Pentachlorobenzene | 608-93-5 | Benzene, pentachloro- | 8270 | 10 |
| Pentachloronitrobenzene | 82-68-8 | Benzene, pentachloronitro- | 8270 | 20 |
| Pentachlorophenol | 87-86-5 | Phenol, pentachloro- | 8040 | 5 |
| | | | 8270 | 50 |
| Phenacetin | 62-44-2 | Acetamide, N-(4-ethoxyphenyl) | 8270 | 20 |
| Phenanthrene | 85-01-8 | Phenanthrene | 8100 | 200 |
| | | | 8270 | 10 |
| Phenol | 108-95-2 | Phenol | 8040 | 1 |
| p-Phenylenediamine | 106-50-3 | 1,4-Benzenediamine | 8270 | 10 |
| Phorate | 298-02-2 | Phosphorodithioic acid, 0,0-diethyl S-[(ethylthio)methyl] ester | 8140 | 2 |
| | | | 8141 | 0.5 |
| | | | 8270 | 10 |
| Polychlorinated biphenyls; PCBs; Aroclors | See Note 9 | 1,1'-Biphenyl, chloro derivatives | 8080 | 50 |
| | | | 8270 | 200 |
| Pronamide | 23950-58-5 | Benzamide, 3,5-dichloro-N-(1,1-dimethyl-2-propynyl)- | 8270 | 10 |
| Propionitrile; Ethyl cyanide | 107-12-0 | Propanenitrile | 8015 | 60 |
| | | | 8260 | 150 |
| Pyrene | 129-00-0 | Pyrene | 8100 | 200 |
| | | | 8270 | 10 |
| Safrole | 94-59-7 | 1,3-Benzodioxole, 5-(2-propenyl)- | 8270 | 10 |
| Selenium | (Dissolved) | Selenium | 6010 | 750 |
| | | | 7740 | 20 |
| | | | 7741 | 20 |
| Silver | (Dissolved) | Silver | 6010 | 70 |
| | | | 7760 | 100 |
| | | | 7761 | 10 |
| Silvex; 2,4,5-TP | 93-72-1 | Propanoic acid, 2-(2,4,5-trichlorophenoxy)- | 8150 | 2 |
| Styrene | 100-42-5 | Benzene, ethenyl- | 8020 | 1 |
| | | | 8021 | 0.1 |
| | | | 8260 | 10 |

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|---|---------------------|---|-----------------------------------|------|
| Sulfide | 18496-25-8 | Sulfide | 9030 | 4000 |
| 2,4,5-T; 2,4,5- Trichlorophenoxyacetic acid | 93-76-5 | Acetic acid, (2,4,5-trichlorophenoxy)- | 8150 | 2 |
| 1,2,4,5-Tetrachlorobenzene | 95-94-3 | Benzene, 1,2,4,5-tetrachloro- | 8270 | 10 |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | Ethane, 1,1,1,2-tetrachloro- | 8010 | 5 |
| | | | 8021 | 0.05 |
| | | | 8260 | 5 |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | Ethane, 1,1,2,2-tetrachloro- | 8010 | 0.5 |
| | | | 8021 | 0.1 |
| | | | 8260 | 5 |
| Tetrachloroethylene; | 127-18-4 | Ethene, tetrachloro- | 8010 | 0.5 |
| Tetrachloroethene; | | | 8021 | 0.5 |
| Perchloroethylene | | | 8260 | 5 |
| 2,3,4,6-Tetrachlorophenol | 58-90-2 | Phenol, 2,3,4,6-tetrachloro- | 8270 | 10 |
| Thallium | (Dissolved) | Thallium | 6010 | 400 |
| | | | 7840 | 1000 |
| | | | 7841 | 10 |
| Tin | (Dissolved) | Tin | 6010 | 40 |
| Toluene | 108-88-3 | Benzene, methyl- | 8020 | 2 |
| | | | 8021 | 0.1 |
| | | | 8260 | 5 |
| o-Toluidine | 95-53-4 | Benzenamine, 2-methyl- | 8270 | 10 |
| Toxaphene | See Note 10 | Toxaphene | 8080 | 2 |
| 1,2,4-Trichlorobenzene | 120-82-1 | Benzene, 1,2,4-trichloro- | 8021 | 0.3 |
| | | | 8120 | 0.5 |
| | | | 8260 | 10 |
| | | | 8270 | 10 |
| 1,1,1-Trichloroethane; | 71-55-6 | Ethane, 1,1,1-trichloro- | 8010 | 0.3 |
| Methylchloroform | | | 8021 | 0.3 |
| | | | 8260 | 5 |
| 1,1,2-Trichloroethane | 79-00-5 | Ethane, 1,1,2-trichloro- | 8010 | 0.2 |
| | | | 8260 | 5 |
| Trichloroethylene; | 79-01-6 | Ethene, trichloro- | 8010 | 1 |
| Trichloroethene | | | 8021 | 0.2 |
| | | | 8260 | 5 |
| Trichlorofluoromethane; CFC-11 | 75-69-4 | Methane, trichlorofluoro- | 8010 | 10 |
| | | | 8021 | 0.3 |
| | | | 8260 | 5 |
| 2,4,5-Trichlorophenol | 95-95-4 | Phenol, 2,4,5-trichloro- | 8270 | 10 |
| 2,4,6-Trichlorophenol | 88-06-2 | Phenol, 2,4,6-trichloro- | 8040 | 5 |
| | | | 8270 | 10 |
| 1,2,3-Trichloropropane | 96-18-4 | Propane, 1,2,3-trichloro- | 8010 | 10 |
| | | | 8021 | 5 |
| | | | 8260 | 15 |
| 0,0,0-Triethyl phosphorothioate | 126-68-1 | Phosphorothioic acid, 0,0,0-triethyl- ester | 8270 | 10 |
| sym-Trinitrobenzene | 99-35-4 | Benzene, 1,3,5-trinitro- | 8270 | 10 |
| Vanadium | (Dissolved) | Vanadium | 6010 | 80 |
| | | | 7910 | 2000 |
| | | | 7911 | 40 |
| Vinyl acetate | 108-05-4 | Acetic acid, ethenyl ester | 8260 | 50 |
| Vinyl chloride; Chloroethene | 75-01-4 | Ethene, chloro- | 8010 | 2 |
| | | | 8021 | 0.4 |
| | | | 8260 | 10 |

| Common Name ² (mg/L) ⁶ | CAS RN ³ | Chemical abstracts service index name ⁴ | Suggested methods ⁵ | PQL |
|---|---------------------|---|-----------------------------------|-----|
| Xylene (total) | See Note 11 | Benzene, dimethyl- | 8020 | 5 |
| | | | 8021 | 0.2 |
| | | | 8260 | 5 |
| Zinc | (Dissolved) | Zinc | 6010 | 20 |
| | | | 7950 | 50 |
| | | | 7951 | 0.5 |

Notes:

- The regulatory requirements pertain only to the list of substances; the right hand columns (Methods and PQL) are given for informational purposes only. See also footnotes 5 and 6. Also, note that the state ground water quality criteria, chapter 173-200 WAC, takes precedence over these recommended PQL's.
- Common names are those widely used in government regulations, scientific publications, and commerce; synonyms exist for many chemicals.
- Chemical Abstracts Service registry number. Where "Total" is entered, all species in the ground water that contain this element are included.
- CAS index are those used in the 9th Collective Index.
- Suggested Methods refer to analytical procedure numbers used in EPA Report SW-846 "Test Methods for Evaluating Solid Waste", third edition, November 1986, as revised, December 1987. Analytical details can be found in SW-846 and in documentation on file at the agency. CAUTION: The methods listed are representative SW-846 procedures and may not always be the most suitable method(s) for monitoring an analyte under the regulations.
- Practical Quantitation Limits (PQLs) are the lowest concentrations of analytes in ground waters that can be reliably determined within specified limits of precision and accuracy by the indicated methods under routine laboratory operating conditions. The PQLs listed are generally stated to one significant figure. PQLs are based on 5 mL samples for volatile organics and 1 L samples for semivolatile organics. CAUTION: The PQL values in many cases are based only on a general estimate for the method and not on a determination for individual compounds; PQLs are not a part of the regulation.
- This substance is often called Bis(2-chloroisopropyl) ether, the name Chemical Abstracts Service applies to its noncommercial isomer, Propane, 2,2'-oxybis[2-chloro- (CAS RN 39638-32-9).
- Chlordane: This entry includes alpha-chlordane (CAS RN 5103-71-9), beta-chlordane (CAS RN 5103-74-2), gamma-chlordane (CAS RN 5566-34-7), and constituents of chlordane (CAS RN 57-74-9 and CAS RN 12789-03-6). PQL shown is for technical chlordane. PQLs of specific isomers are about 20 µg/L by method 8270.
- Polychlorinated biphenyls (CAS RN 1336-36-3); this category contains congener chemicals, including constituents of Aroclor 1016 (CAS RN 12674-11-2), Aroclor 1221 (CAS RN 11104-28-2), Aroclor 1232 (CAS RN 11141-16-5), Aroclor 1242 (CAS RN 53469-21-9), Aroclor 1248 (CAS RN 12672-29-6), Aroclor 1254 (CAS RN 11097-69-1), and Aroclor 1260 (CAS RN 11096-82-5). The PQL shown is an average value for PCB congeners.
- Toxaphene: This entry includes congener chemicals contained in technical toxaphene (CAS RN 8001-35-2), i.e., chlorinated camphene.
- Xylene (total): This entry includes o-xylene (CAS RN 96-47-6), m-xylene (CAS RN 108-38-3), p-xylene (CAS RN 106-42-3), and unspecified xylenes (dimethylbenzenes) (CAS RN 1330-20-7). PQLs for method 8021 are 0.2 for o-xylene and 0.1 for m-or p-xylene. The PQL for m-xylene is 2.0 µg/L by method 8020 or 8260.

APPENDIX IV

PARAMETERS FOR LEACHATE ANALYSIS

Appendix I¹ Parameters

Appendix II Parameters

Nitrite

Total Colliform

COD

BOD

Cyanide

- 1_ All metals analysis should be for total recoverable metals, for the leachate analysis only.

Important Note: All other appendices require dissolved metals (field-filtration for metals).

[Statutory Authority: Chapter 70.95 RCW and 40 CFR 258. 93-22-016, § 173-351-990, filed 10/26/93, effective 11/26/93.]

Reviser's note: The brackets and enclosed material in the text of the above section occurred in the copy filed by the agency.

Appendix H

Area-Wide Arsenic Data Discussion and Background Determination for Arsenic, Iron, and Manganese



TECHNICAL MEMORANDUM

Arsenic in Groundwater at the Olalla Landfill Kitsap County, WA

Introduction

Arsenic is a naturally occurring element in the environment and its presence in groundwater is largely the result of natural dissolution of arsenic-bearing minerals over time (USGS, 2008). Elevated concentrations of arsenic in groundwater affect many parts of the country where arsenic-bearing minerals are common, including the Puget Sound area of Washington State.

This Technical Memorandum has been prepared to provide supporting documentation to demonstrate that the occurrence of arsenic in groundwater samples from monitoring wells at the Olalla Landfill is partially caused by regional elevated arsenic concentrations in soil and groundwater.

Arsenic in Regional Groundwater

Higher levels of arsenic tend to be found more in groundwater sources than in surface water sources (i.e., lakes and rivers) of drinking water. Many states, including Washington, have a significant number of public water systems with arsenic concentrations greater than EPA's standard of 10 $\mu\text{g/L}$ (EPA, 2011).

The United States Geological Survey (USGS) has evaluated arsenic data and has published maps summarizing a national data set for arsenic in groundwater. These maps provide a summary view of patterns of naturally occurring arsenic in groundwater across the United States. The arsenic occurrence map presented in Figure 1 shows elevated arsenic concentrations greater than the drinking water standard of 10 $\mu\text{g/L}$ in western Washington, particularly in the Puget Sound, including Kitsap County.

The Washington State Department of Health (DOH) notes arsenic is found in well water throughout Washington State and recommends that water used for drinking or food preparation contain no more than 10 $\mu\text{g/L}$ arsenic (DOH, 2008). DOH tracks arsenic in samples from regulated public drinking water sources at concentrations greater than the drinking water standard of 10 $\mu\text{g/L}$. These data are summarized on a map of Washington State presented in Figure 2. This map shows that regulated water systems in Kitsap County, including locations near the Olalla Landfill, have arsenic concentrations greater than the Washington State Drinking Water Standard of 10 $\mu\text{g/L}$.

The Bremerton-Kitsap County Health District Environmental Health Division conducted a domestic well survey in 1995 where a background level of arsenic in southern Kitsap County was determined to be 9.69 $\mu\text{g/L}$ (KCHD 1995). The background arsenic concentration was taken from the Kitsap County Ground Water Management Plan, July 1989. The calculated concentration represents an average concentration measured in 84 wells in southern Kitsap County. According to the survey, the wells used in the comparison are all greater than one hundred feet deep; therefore, the wells are considered comparable to the conditions at the Olalla Landfill.

Arsenic in Kitsap County Groundwater

As groundwater flows through rocks and soil that contain naturally occurring arsenic bearing minerals some of the arsenic dissolves into the groundwater. Drinking water in Washington State typically contains less than 3 µg/L (DOH, 2008). However, levels of health concern (from 10 µg/L to 33,000 µg/L) have been detected in samples from some wells in Washington. These concentrations are commonly associated with aquifers located in rock or soil that has naturally high arsenic content (DOH, 2008).

Much of the Puget Sound area, including Kitsap County, has near surface geology consisting largely of glacially deposited volcanic rocks, which contain naturally occurring arsenic. In addition, the ASARCO smelter formerly operating in Ruston WA, approximately 10 miles from the Landfill, was one of the few smelters able to process ore containing high concentrations of arsenic. The ASARCO smelter released particulates containing high concentrations of arsenic and lead to the atmosphere, which was carried by the wind over a wide expanse of King, Pierce, Thurston, and Kitsap counties (Seattle and King County Public Health, 2011). For comparison to local groundwater arsenic concentrations, arsenic concentrations averaging 17 µg/L were detected in rain and snow collected downwind from the ASARCO smelter (Creclius 1975).

The DOH Office of Drinking Water database contains analytical data, which include arsenic data for Kitsap County Class A and B water supply wells. Samples from many of these wells have arsenic at concentrations greater than the Washington State Drinking Water Standard of 10 µg/L. Database queries in the DOH database indicate that arsenic concentrations in Kitsap County Class A water supply wells are as high as 32 µg/L and as high as 100 µg/L in Class B water supply wells. Most of the drinking water standard exceedances in Kitsap County drinking water wells for which DOH has data are in the 11-20 µg/L range. The arsenic data exceeding drinking water standards in Kitsap County water supply wells are summarized in Figure 2 and Figure 3 (DOH, 2011). Among the public water supply systems noted in Figure 3, West Sound Utility Water District #1 has water supply wells that are closest to the Landfill at a distance of approximately 7 miles.

Arsenic in Olalla Landfill Groundwater

Arsenic has been routinely detected from 1992 to present in groundwater samples from all monitoring wells at the Landfill, including samples from upgradient well MW-1. Detected arsenic concentrations in samples from upgradient well MW-1 range from a high of 2.0 µg/L in January 1994 to a low of 0.05 µg/L during several monitoring events, the most recent in December 2002. The mean arsenic concentration in samples from 1992 to 2011 from MW-1 is 0.171 µg/L. Arsenic concentrations in samples from downgradient offsite wells range from a high of 11 µg/L to non-detect at a reporting limit of 0.05 µg/L. The mean arsenic concentration in samples from downgradient offsite wells is 1.04 µg/L.

Arsenic was also detected in samples from all water supply wells that were sampled as part of the RI/FS at concentrations ranging from a high of 7.04 µg/L in the sample from OW-3 to a low of 0.215 µg/L in the sample from OW-2. The water supply well designated as OW-3, which had the greatest arsenic concentration, is completed in a deeper confined aquifer that is separated from the unconfined aquifer underlying the Landfill by approximately 145 feet of dense clay.

Arsenic Background Concentration in Olalla Landfill Groundwater

Water supply wells were sampled prior to and during the RI activities conducted at the Landfill. Specifically, water supply wells were sampled in April 1995, September 1997, December 2010, and January 2011. The sampling results from the water supply wells and Landfill upgradient monitoring well MW-1 were used to define the data set for background concentration determination. Parametrix provided background calculations for arsenic in a letter dated August 13, 2012 (Parametrix, 2012). The background concentration of arsenic in groundwater at the Olalla Landfill was determined to be 1.29 µg/L.

Conclusions

The following conclusions regarding the presence of arsenic in groundwater are based on information and data presented in the preceding sections.

- Arsenic is a naturally occurring element that is found in groundwater in areas containing arsenic-bearing minerals.
- The Puget Sound, including Kitsap County, contains arsenic-bearing minerals that affect groundwater quality throughout the region.
- The presence of arsenic in the deeper aquifer at a concentration significantly greater than the mean concentration in data from downgradient monitoring well MW-6 indicates that the presence of arsenic in groundwater is partially a regional issue related to arsenic-bearing minerals.
- An arsenic background concentration of 1.29 µg/L was calculated for the Landfill using procedures described within MTCA.

References:

United States Geological Survey (USGS), Fact Sheet 063-00, August 7, 2008.

United States Environmental Protection Agency (EPA, 2011.) Web site reference:
<http://water.epa.gov/lawsregs/rulesregs/sdwa/arsenic/Basic-Information.cfm>

Washington State Department of Health (DOH), Arsenic and Your Private Well. Publication # 334-156, June 2008.

Bremerton-Kitsap County Health District (KCHD), Environmental Health Division, Solid Waste Program. Olalla Landfill Domestic Well Survey Results. Prepared by Shawn Ultican, Environmental Health Specialist. October 23, 1995.

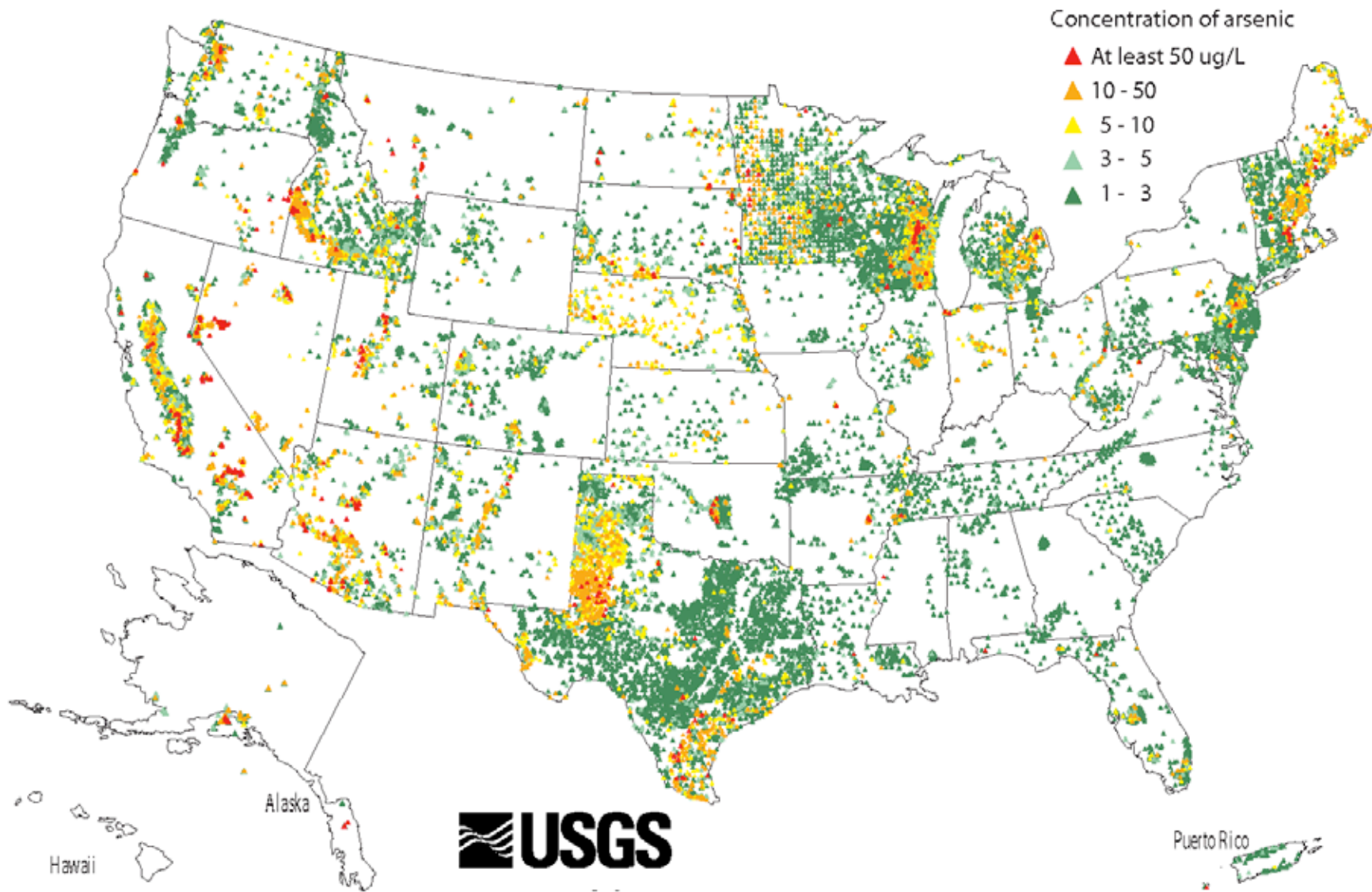
Seattle and King County Public Health website. Background on the Tacoma Smelter Plume. <http://www.kingcounty.gov/healthservices/health/ehs/toxic/TacomaSmelterPlume/background.aspx>. Updated September 9, 2011.

Crececius, E.A. 1975. The geochemical cycle of arsenic in Lake Washington and its relation to other elements. *Limnology and Oceanography* 20, no. 3: 441-451.

Washington State Department of Health DOH, Division of Environmental Health, Office of Drinking Water. <https://fortress.wa.gov/portal/odw/si/ListWaterQualityExceedences.aspx>. (2011)

Blakely, N. 1992. *Statistical Guidance for Ecology Site Managers*. Washington Department of Ecology. Olympia, Washington. (Ecology, 1992)

Parametrix, 2012. Re: Olalla RI/FS Background Calculation Scenarios for Arsenic in Groundwater. Blaine Hardy, Parametrix to Krystyna Kowalik, Department of Ecology. August 13.



Concentration of arsenic

- ▲ At least 50 ug/L
- ▲ 10 - 50
- ▲ 5 - 10
- ▲ 3 - 5
- ▲ 1 - 3



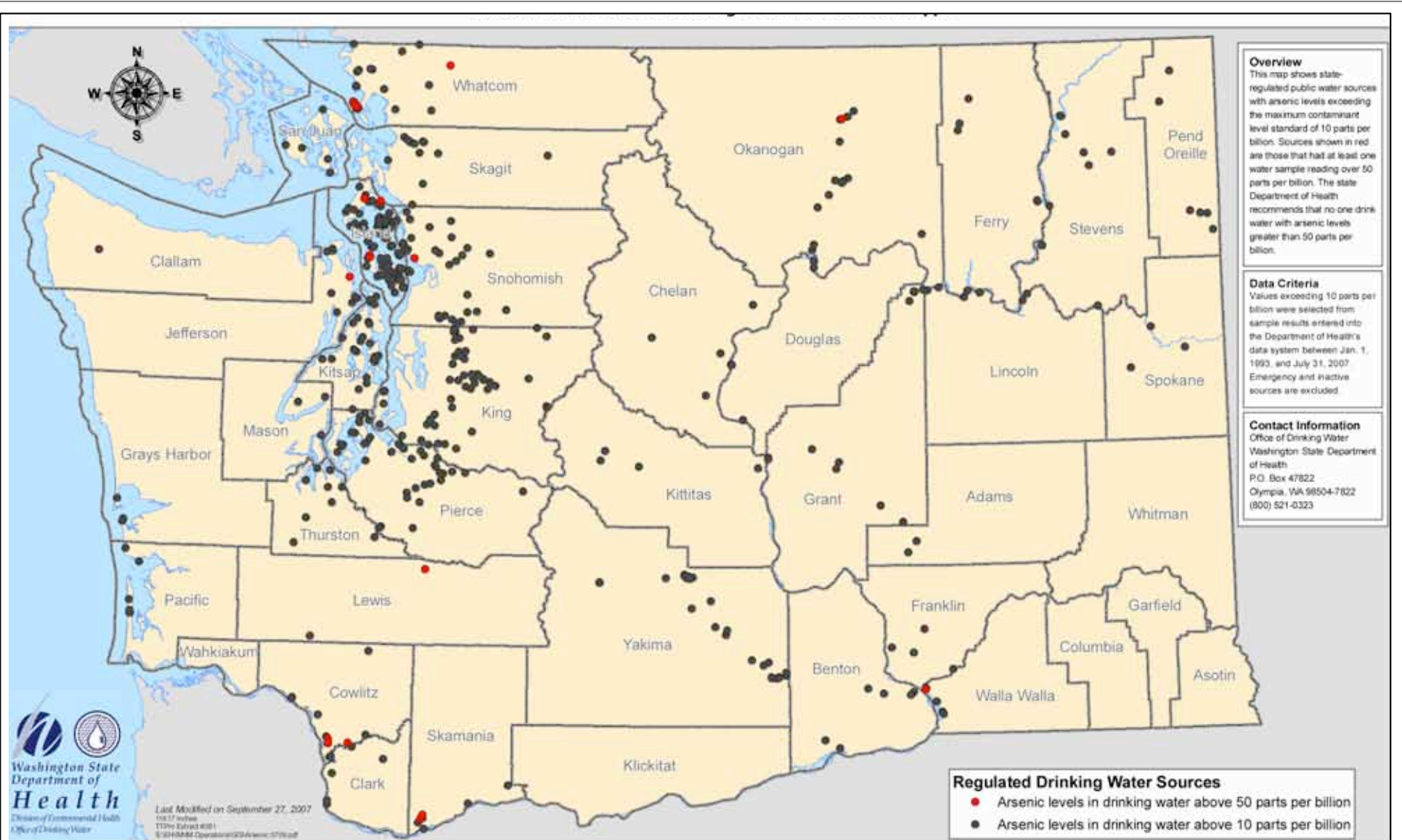
KEY:

SOURCE:
 U.S. Department of the Interior | U.S. Geological Survey
 URL: http://water.usgs.gov/nawqa/trace/pubs/geo_v46n11/fig1.html
 Page Contact Information: gs-w_nawqa_info@usgs.gov
 Page Last Modified: Tuesday, 20-Oct-2009 16:04:53 EDT

epl ENVIRONMENTAL PARTNERS INC
 295 NE Gilman Boulevard, Suite 201
 Issaquah, Washington 98027

FIGURE 1
 ARSENIC CONCENTRATIONS IN
 GROUNDWATER THROUGHOUT THE
 UNITED STATES

| | | | |
|------------------------|--|---------------------------|-------------------------|
| PROJECT | 60101.0 OLALLA, WASHINGTON | | |
| PREPARED FOR | KITSAP COUNTY | | |
| LOCATION | OLALLA LANDFILL KITSAP COUNTY, WASHINGTON | | |
| SHEET 1 of 1 | DRAWN BY ARM | REVIEWED BY DCK | DATE 02/14/12 |



KEY:

epl ENVIRONMENTAL PARTNERS INC
295 NE Gilman Boulevard, Suite 201
Issaquah, Washington 98027

FIGURE 2

ARSENIC DETECTIONS IN WASHINGTON
PUBLIC WATER SUPPLIES

PROJECT 60101.0 OLALLA, WASHINGTON

PREPARED FOR KITSAP COUNTY

LOCATION OLALLA LANDFILL
KITSAP COUNTY, WASHINGTON

| SHEET | DRAWN BY | REVIEWED BY | DATE |
|--------------|-----------------|--------------------|-------------|
| 1 of 1 | ARM | DCK | 02/14/12 |



Division of Environmental Health Office of Drinking Water

Help

Water Quality Exceedences View

| WSID | WS Name | County | Grp | Type | Status | Src # | DOE Src | Collect Date | Analyte Name | Result Quantity | UOM Code | Test Panel | Analyte Group | Sample Purpose | Sam # | Lab # |
|-----------------------|--------------------------------|--------|-----|------|--------|-------|---------|--------------|--------------|-----------------|----------|------------|---------------|----------------|-----------------------|-------|
| 67613 | AMBER WATER ASSOCIATION | KITSAP | B | | Act | 01 | | 4/15/1991 | ARSENIC | 0.100 | mg/L | ICHEM | IOC | RC | 04269 | 081 |
| 46117 | LARSON WATER | KITSAP | B | | Act | 01 | | 3/11/1980 | ARSENIC | 0.065 | mg/L | ICHEM | IOC | RC | 03470 | 052 |
| 46117 | LARSON WATER | KITSAP | B | | Act | 01 | | 2/19/1979 | ARSENIC | 0.061 | mg/L | ICHEM | IOC | RC | 03413 | 051 |
| 00700 | THORS WELL ASSN | KITSAP | B | | Act | 01 | | 1/17/2006 | ARSENIC | 0.036 | mg/L | IOC | IOC | RC | 43407 | 010 |
| AC483 | Breidablik Baptist Church | KITSAP | A | TNC | Act | 01 | | 7/22/2002 | ARSENIC | 0.032 | mg/L | AR | IOC | RC | 76120 | 010 |
| AC242 | YAQUINA | KITSAP | B | | PreAct | 01 | | 10/14/2010 | ARSENIC | 0.025 | mg/L | IOC_SHORT | IOC | RC | 04701 | 010 |
| AC242 | YAQUINA | KITSAP | B | | PreAct | 01 | | 7/15/2010 | ARSENIC | 0.020 | mg/L | IOC_SHORT | IOC | RC | 25601 | 010 |
| AC242 | YAQUINA | KITSAP | B | | PreAct | 01 | | 6/17/2010 | ARSENIC | 0.018 | mg/L | IOC_SHORT | IOC | RC | 37001 | 010 |
| 05122 | NORTH PENINSULA | KITSAP | A | Comm | Act | 13 | | 6/15/2006 | ARSENIC | 0.016 | mg/L | IOC | IOC | RC | 55988 | 010 |
| 05122 | NORTH PENINSULA | KITSAP | A | Comm | Act | 13 | | 6/13/2007 | ARSENIC | 0.016 | mg/L | IOC_SHORT | IOC | RC | 85739 | 010 |
| 05122 | NORTH PENINSULA | KITSAP | A | Comm | Act | 13 | | 10/6/2010 | ARSENIC | 0.016 | mg/L | IOC | IOC | RC | 82301 | 010 |
| 92400 | WALKER BEACH | KITSAP | B | | Act | 02 | | 7/6/1999 | ARSENIC | 0.016 | mg/L | IOC | IOC | RC | 35890 | 010 |
| AC242 | YAQUINA | KITSAP | B | | PreAct | 01 | | 7/15/2010 | ARSENIC | 0.016 | mg/L | IOC_SHORT | IOC | RC | 25602 | 010 |
| 02600 | WEST SOUND UTILITY DISTRICT #1 | KITSAP | A | Comm | Act | 14 | 15G014 | 4/19/1983 | ARSENIC | 0.015 | mg/L | ICHEM | IOC | RC | 04344 | 089 |
| 05122 | NORTH PENINSULA | KITSAP | A | Comm | Act | 13 | | 2/7/2007 | ARSENIC | 0.015 | mg/L | IOC_SHORT | IOC | RC | 75862 | 010 |
| 23740 | ERICKSON | KITSAP | B | | Act | 01 | 15G502 | 3/18/2007 | ARSENIC | 0.015 | mg/L | IOC | IOC | RC | 78974 | 010 |
| 66936 | BUCKLIN | KITSAP | A | Comm | Act | 01 | 15G493 | 2/14/2007 | ARSENIC | 0.014 | mg/L | AR | IOC | RC | 14983 | 089 |
| 66936 | BUCKLIN | KITSAP | A | Comm | Act | 01 | 15G493 | 4/24/2007 | ARSENIC | 0.014 | mg/L | AR | IOC | RC | 16015 | 089 |
| AB876 | TOAD HOLLER | KITSAP | A | NTNC | Act | 01 | | 1/6/2010 | ARSENIC | 0.014 | mg/L | AR | IOC | RC | 20203 | 010 |
| AB876 | TOAD HOLLER | KITSAP | A | NTNC | Act | 01 | | 4/12/2011 | ARSENIC | 0.014 | mg/L | AR | IOC | RC | 38401 | 010 |
| AB876 | TOAD HOLLER | KITSAP | A | NTNC | Act | 01 | | 7/12/2011 | ARSENIC | 0.014 | mg/L | AR | IOC | RC | 75301 | 010 |
| 05122 | NORTH PENINSULA | KITSAP | A | Comm | Act | 13 | | 6/15/2006 | ARSENIC | 0.013 | mg/L | IOC_SHORT | IOC | RC | 55987 | 010 |
| 05122 | NORTH PENINSULA | KITSAP | A | Comm | Act | 13 | | 7/27/2010 | ARSENIC | 0.013 | mg/L | AR | IOC | RC | 60002 | 010 |
| 24800 | FERNCLIFF | KITSAP | A | Comm | Act | 02 | | 9/5/2006 | ARSENIC | 0.013 | mg/L | IOC | IOC | RC | 64585 | 010 |
| 66936 | BUCKLIN | KITSAP | A | Comm | Act | 01 | 15G493 | 8/23/2006 | ARSENIC | 0.013 | mg/L | IOC_SHORT | IOC | RC | 11852 | 089 |

KEY:



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*295 NE Gilman Boulevard, Suite 201
Issaquah, Washington 98027*

FIGURE 3

ARSENIC CONCENTRATIONS EXCEEDING DRINKING WATER STANDARDS IN REGULATED WATER SUPPLY WELLS, KITSAP COUNTY, WASHINGTON

PROJECT

60101.0 OLALLA, WASHINGTON

PREPARED FOR

KITSAP COUNTY

LOCATION

OLALLA LANDFILL
KITSAP COUNTY, WASHINGTON

SHEET

1 of 1

DRAWN BY

ARM

REVIEWED BY

DCK

DATE

08/25/11

4660 KITSAP WAY, SUITE A
BREMERTON, WA 98312-2357
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www.parametrix.com

August 13, 2012
PMX No. 215-1578-121

Ms. Krystyna Kowalik
Regional Hydrogeologist
Department of Ecology
Northwest Regional Office
3190 160th Ave SE
Bellevue, WA 98008-5452

Re: Olalla RI/FS Background Calculation Scenarios for Arsenic in Groundwater

Dear Ms. Kowalik:

INTRODUCTION

The May 31, 2012 Parametrix letter regarding Kitsap County Olalla Landfill Remedial Investigation /Feasibility Study (RI/FS) Arsenic Background Calculation Scenarios contained several scenarios for selecting data to be included in the arsenic background calculations. The recommended Scenario 2 uses data from wells located in close proximity to the Landfill and known to be completed in the uppermost aquifer beneath the landfill, including on-site wells MW-1 (upgradient) and MW-5A (cross-gradient), and off-site wells OW-1, OW-2, OW-4, and OW-9. The July 20, 2012 Department of Ecology letter agreed with the overall approach for Scenario 2 but requested that the calculations be expanded using the last three years of groundwater quality data for wells MW-1 and MW-5A. The data set was reevaluated in consideration of this request, and the results are presented in this letter.

BACKGROUND

In accordance with WAC 173-340-200, "Natural background" means the concentration of a hazardous substance consistently present in the environment that has not been influenced by localized human activities. In accordance with WAC 173-340-709(c), for log normally distributed data sets, background shall be defined as the true upper 90th percentile or four times the true 50th percentile, whichever is lower.

Additional guidance for determining background levels is presented in the Unified Guidance, Statistical Analysis of Groundwater Monitoring Data at RCRA Facilities (EPA 2009), which states that sample data from the background population are assumed to be independent and identically distributed. A key implication of this assumption is that a series of sample measurements should be stationary. A *stationary* statistical distribution is one whose population characteristics do not change over time and/or space.

*Inspired people
Inspired solutions
making a difference*



ANALYSIS

A time-series plot of the historical data for MW-1 and MW-5A (Attachment A) indicates that the more recent data (since 2003) have lower variability than previous data, and do not contain obvious outliers or apparent trends. The lower variability is likely due to improved sampling procedures.

Trends in the data for MW-1 and MW-5A were evaluated using the Mann-Kendall and Sens Slope tests for trend. Three different sets of data were evaluated for each well: 1) all data since 1992; 2) data from 2003 to 2012; and 3) data from 2009 to 2012. Statistically significant downward trends in the data were observed using all data for both wells, and using the 2003-2012 data for well MW-5A. There were no upward or downward trends in the data for either well using the 2009-2012 data. Therefore using the past three years of data (2009-2012) for these wells, as requested by Ecology, is appropriate to represent the current water quality conditions at those locations.

Since the concentrations of arsenic in samples from wells in the vicinity of the landfill that are located within the same aquifer have a natural spatial variability due to variations in soil geochemistry, it is important to incorporate as many spatial locations as possible to calculate the arsenic background level. Therefore data for the four off-site wells not affected by the landfill will be used in addition to the two on-site wells in calculating the background concentration for arsenic, even though less data are available for the off-site wells. Since there are 13 or 14 points available in the 2009-2012 data set for wells MW-1 and MW-5A, it is necessary to calculate one representative concentration for each off-site well so that all six wells will be equally weighted spatially in the background calculation.

Probability plots were prepared using the 2009-2012 data for MW-1 and MW-5A and the results indicate the data are normally distributed (Attachment B). The distribution and summary statistics for each well were then verified using *MTCASat* and the recommended distribution was lognormal (see calculations in Attachment C). Therefore the lognormal mean of the data for each of the wells (0.106 ug/L at MW-1 and 0.160 ug/L at MW-5A) was selected to represent the current concentration at that well.

For off-site wells OW-2 and OW-4, either the most recent data point could be used or a representative concentration could be calculated that would use all three measurements (1995 through 2010/2011). However, since the data are not adequate to evaluate the presence of trends or to assess the distribution of the data (see distribution and summary statistics in Attachment C), it is recommended that only the most recent data point be used to represent those wells in the background calculations.

Background arsenic concentrations were calculated using *MTCASat* using the mean of the 2009-2012 data for MW-1 and MW-5A, and the individual data points collected in 2010 or 2011 for all the off-site wells (OW-1, OW-2, OW-4, and OW-9). This data set was determined to be log-normally distributed and the resulting background arsenic concentration is 1.29 ug/L. The data used are summarized in the table below and the calculations using *MTCASat* are presented in Attachment D.



Data Used to Calculate Arsenic Background Concentration Using MCTAStat

| Arsenic Concentration (ug/L) | Well | Data Used in Background Calculation | Calculated Arsenic Background Concentration (ug/L) |
|-------------------------------------|-------------|--|---|
| 0.719 | OW-1 | 12/28/2010 | |
| 0.215 | OW-2 | 1/27/2011 | |
| 1.68 | OW-4 | 12/28/2010 | |
| 0.253 | OW-9 | 12/28/2010 | |
| 0.106 | MW-1 | 2009-2012 Lognormal Mean | |
| 0.160 | MW-5A | 2009-2012 Lognormal Mean | 1.29 ug/L |

CONCLUSION

Based on arsenic groundwater data from 2009 through 2012 for wells located in close proximity to the Landfill and known to be completed in the uppermost aquifer beneath the landfill, the background arsenic concentration is 1.29 ug/L.

Sincerely,
PARAMETRIX, INC.

Blaine Hardy, PE

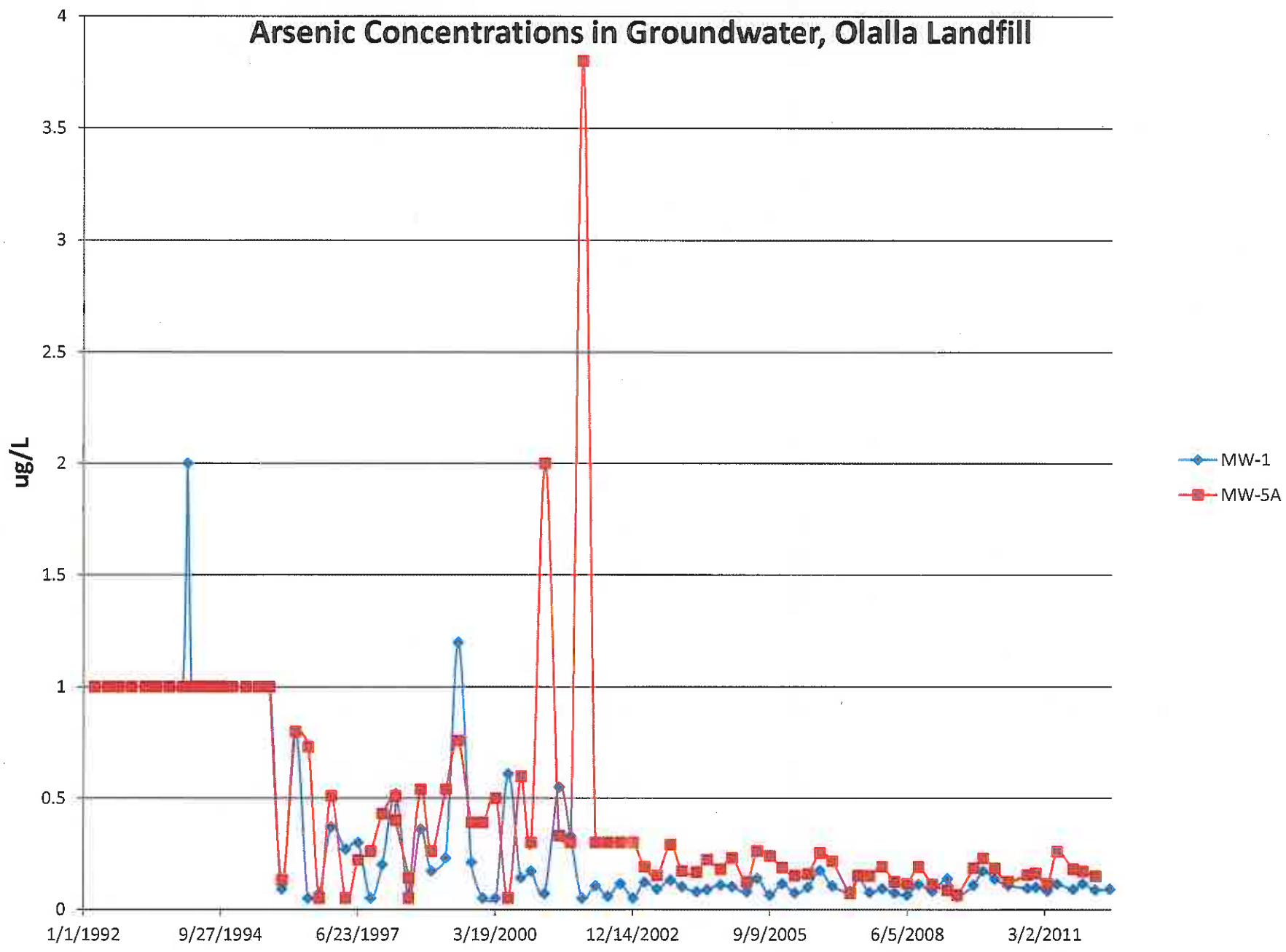
Attachments

- cc: Jan Brower, Kitsap Public Health District
- Madeline Wall, Department of Ecology
- Peter Christiansen, Department of Ecology
- Keli McKay-Means, Kitsap County Public Works
- Pat Campbell, Kitsap County Public Works
- Doug Kunkel, Environmental Partners, Inc.

ATTACHMENT A

Time Series Plot MW-1 and MW-5A

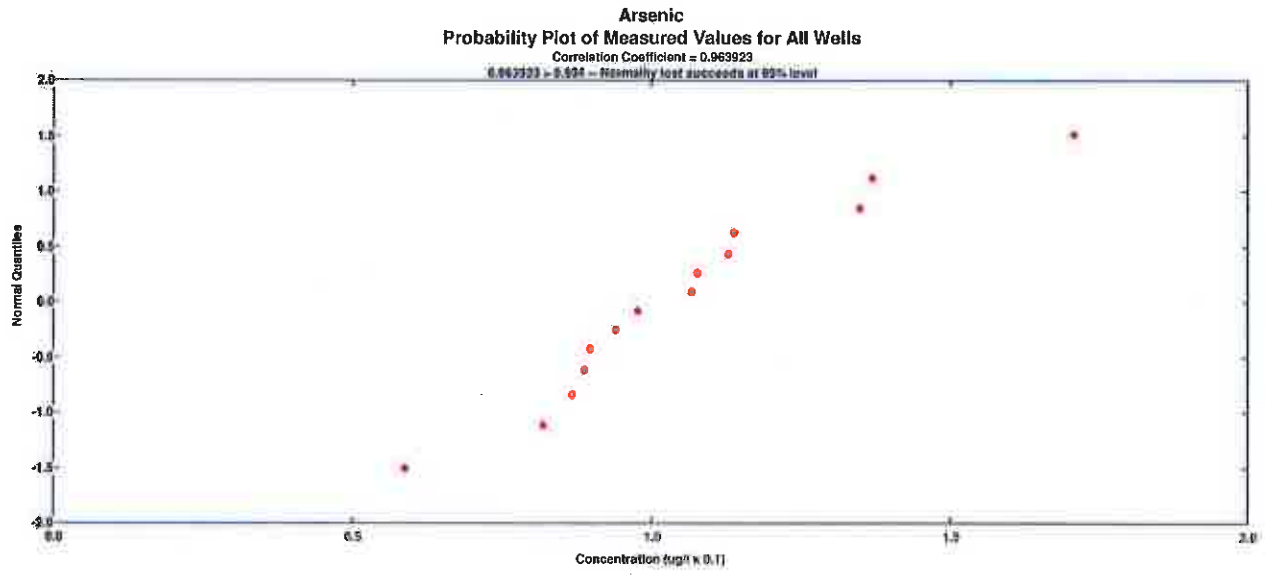
Arsenic Concentrations in Groundwater, Olalla Landfill



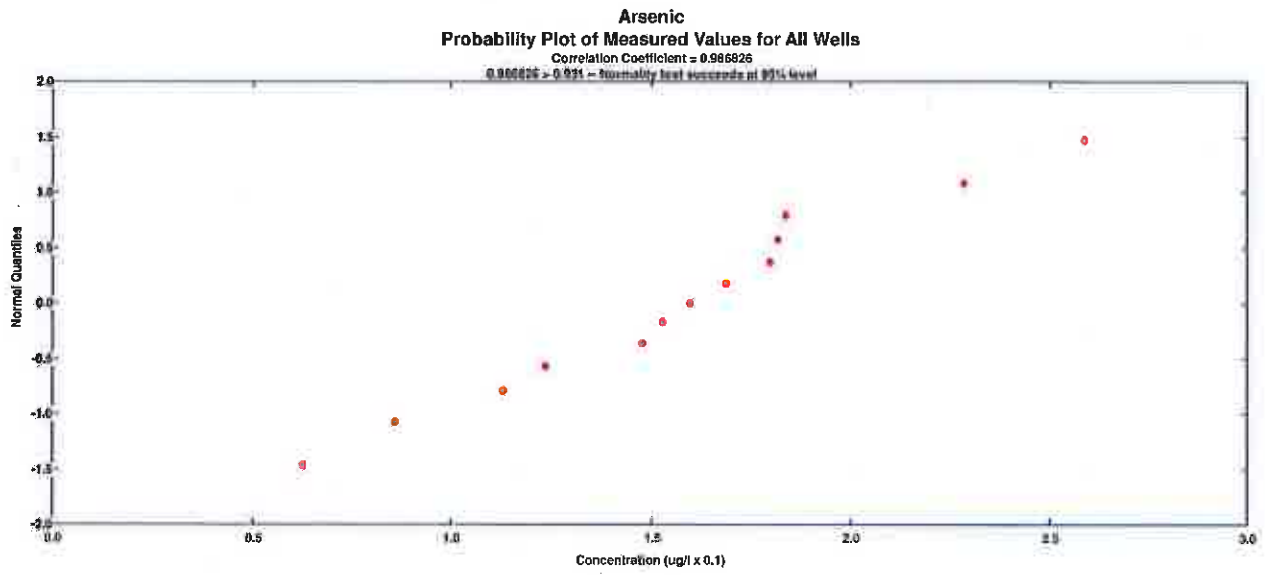
ATTACHMENT B

Probability Plots MW-1 and MW-5A

MW-1 2009-2012



MW-5A 2009-2012



ATTACHMENT C

Distribution and Summary Statistics

Compliance Calculations

MW-1 Arsenic Data (2009-2012)

| Concentration (ug/L) | Date (mm/dd/yy) |
|-------------------------|--------------------|
| 0.059 | 6/2/09 |
| 0.082 | 3/23/11 |
| 0.087 | 3/6/12 |
| 0.089 | 9/26/11 |
| 0.09 | 6/21/12 |
| 0.0944 | 10/28/10 |
| 0.098 | 12/28/10 |
| 0.107 | 6/9/10 |
| 0.108 | 9/30/09 |
| 0.113 | 6/1/11 |
| 0.114 | 12/5/11 |
| 0.135 | 3/3/10 |
| 0.137 | 3/23/09 |

Distribution and Summary Stats by the MCTAStat Method

| Number of samples | | Summary Statistics | |
|------------------------|----|--------------------|---------|
| Uncensored | 14 | Mean | 0.1060 |
| Censored | 0 | Lognormal mean | 0.1063 |
| Detection limit or PQL | | Std. devn. | 0.02778 |
| Method detection limit | | Median | 0.1025 |
| TOTAL | 14 | Min. | 0.059 |
| | | Max. | 0.171 |

| | |
|-------------------------|----------------------|
| Lognormal distribution? | Normal distribution? |
| r-squared is: 0.956 | r-squared is: 0.932 |

Recommendations:

Use lognormal distribution.

Compliance Calculations

MW-5A Arsenic Data (2009-2012)

| Concentration (ug/L) | Date (mm/dd/yy) |
|-------------------------|--------------------|
| 0.063 | 6/2/09 |
| 0.086 | 3/23/09 |
| 0.113 | 3/23/11 |
| 0.124 | 6/9/10 |
| 0.148 | 3/6/12 |
| 0.153 | 10/28/10 |
| 0.16 | 12/28/10 |
| 0.169 | 12/5/11 |
| 0.18 | 9/26/11 |
| 0.182 | 3/3/10 |
| 0.184 | 9/30/09 |
| 0.229 | 12/8/09 |
| 0.259 | 6/2/11 |

Distribution and Summary Stats by the MCTAStat Method

| Number of samples | | Summary Statistics | |
|------------------------------------|----|----------------------|---------|
| Uncensored | 13 | Mean | 0.1577 |
| Censored | 0 | Lognormal mean | 0.1596 |
| Detection limit or PQL | | Std. devn. | 0.05371 |
| Method detection limit | | Median | 0.1600 |
| TOTAL | 13 | Min. | 0.063 |
| | | Max. | 0.259 |
| Lognormal distribution? | | Normal distribution? | |
| r-squared is: 0.930 | | r-squared is: 0.975 | |
| Recommendations: | | | |
| Use lognormal distribution. | | | |

Compliance calculations

OW-2 Arsenic Data

| Concentration (ug/L) | Date (mm/dd/yy) |
|-------------------------|--------------------|
| 0.16 | 9/24/97 |
| 0.215 | 1/27/11 |
| 0.48 | 4/25/95 |

Distribution and Summary Stats by the MCTAStat Method

| | | | |
|--|---|----------------------|---------|
| Number of samples | | Uncensored values | |
| Uncensored | 3 | Mean | 0.2850 |
| Censored | 0 | Lognormal mean | 0.2993 |
| Detection limit or PQL | | Std. devn. | 0.17110 |
| Method detection limit | | Median | 0.2150 |
| TOTAL | 3 | Min. | 0.16 |
| | | Max. | 0.48 |
| Lognormal distribution? | | Normal distribution? | |
| r-squared is: | | r-squared is: | |
| Recommendations: | | | |
| Unable to analyze probability plot for lognormal case. | | | |
| Unable to analyze probability plot for normal case. | | | |
| Consult Statistical Guidance document | | | |

Compliance calculations

OW-4 Arsenic Data

| Concentration (ug/L) | Date (mm/dd/yy) |
|-------------------------|--------------------|
| 1.68 | 12/28/10 |
| 1.8 | 9/24/97 |
| 2.77 | 4/25/95 |

Distribution and Summary Stats by the MCTAStat Method

| | | | |
|--|---|----------------------|---------|
| Number of samples | | Uncensored values | |
| Uncensored | 3 | Mean | 2.0833 |
| Censored | 0 | Lognormal mean | 2.1069 |
| Detection limit or PQL | | Std. devn. | 0.59769 |
| Method detection limit | | Median | 1.8000 |
| TOTAL | 3 | Min. | 1.68 |
| | | Max. | 2.77 |
| Lognormal distribution? | | Normal distribution? | |
| r-squared is: | | r-squared is: | |
| Recommendations: | | | |
| Unable to analyze probability plot for lognormal case. | | | |
| Unable to analyze probability plot for normal case. | | | |
| Consult Statistical Guidance document | | | |

ATTACHMENT D

Background Calculations Using *MCTAStat*

Background calculations

0.106 MW-1 Calculation of Arsenic Background for Olalla Landfill
 0.16 MW-5A
 0.215 OW-2
 0.253 OW-9
 0.719 OW-1
 1.68 OW-4

| MTCASat 3.0 | | | | |
|---|---|---------------------------------|--|--|
| Number of samples | | Uncensored values | | |
| Uncensored | 6 | Mean | 0.52 | |
| Censored | 0 | Lognormal mean | 0.55 | |
| TOTAL | 6 | Std. devn. | 0.61 | |
| | | Median | 0.234 | |
| | | Min. | 0.106 | |
| | | Max. | 1.68 | |
| Lognormal distribution? | | Normal distribution? | | |
| r-squared is: 0.92 | | r-squared is: 0.73 | | |
| Recommendations: | | | | |
| <div style="border: 3px double black; padding: 5px; display: inline-block;"> Use lognormal distribution. </div> | | | | |
| Distribution selection | | Enter percentile | Value corresponding to that percentile is: | |
| 1 | | 90 | 1.44 | |
| 1 = Lognormal | | 50th | 0.32 | |
| 2 = Normal | | 4 X 50th | 1.29 | |
| 3 = Nonparametric method | | Coefficient of Variation = 1.72 | | |

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www.parametrix.com

TECHNICAL MEMORANDUM

Date: December 23, 2013
To: File
From: David Dinkuhn
Subject: Background Calculations for Iron and Manganese
Project Number: 215-1578-121
Project Name: Olalla Landfill Remedial Investigation/Feasibility Study

STATISTICAL METHODS

Manganese and iron data for up gradient well MW-1, cross-gradient well MW-5A, and four off-site drinking water wells (OW1, OW2, OW4, and OW9) were used to calculate the background concentrations at the Olalla Landfill using the data evaluation procedures for arsenic described in the Parametrix letter to the Department of Ecology dated August 13, 2012. The 2009 through 2012 data were selected to represent background because they showed lower variability due to improved sampling procedures. For Wells MW-1 and MW-5A, the data evaluated were collected during the 3-year period from 2009 to 2012. For the off-site Wells OW-1, OW-2, OW-4 and OW-9, data consisted of the most recent data value collected during the RI/FS (samples collected in December 2010 and January 2011).

The manganese and iron data for MW-1 and MW-5A were converted from mg/L to $\mu\text{g/L}$ and tested for normality using the probability plot method described by the Washington Department of Ecology (1992) for Model Toxic Control Act (MTCA) site cleanup work. Non-detects were evaluated as recorded in the data sets (equal to the method detection limit (MDL) or reporting limit (RL)) to be consistent with data from the off-site wells. Subsequently, the 50th and 90th percentiles of data from the 2009–2012 period were calculated, and background concentrations were determined in preparation for establishing cleanup levels for manganese and iron at the landfill. Non-detects were evaluated as one-half MDL.

MANGANESE

The manganese data for MW-1 (n=14) are neither log normally or normally distributed ($r^2 = 0.42$ and 0.42 respectively) with a median = $5 \mu\text{g/L}$. The manganese data for MW-5A (n=14) are neither log normally or normally distributed ($r^2 = 0.44$ and 0.44 respectively) with a median = $5 \mu\text{g/L}$.

These median data were combined as a data set with the manganese values for the off-site wells. The 90th and 50th percentiles were calculated using the log transformed data as described by Ecology (1992). The values of the percentiles are not applicable (N/A) and 2.5 respectively and the coefficient of variation (CV) is N/A. Non-parametric estimates of the 90th percentile and the CV cannot be calculated due to the small size of the data set. Although the 90th percentile cannot be estimated, that value must be greater than 32 (the highest value and rank of the manganese data set). The Ecology guidance (1992) indicates that the background value should be set at four times the 50th percentile when the 90th percentile is greater than four times the 50th percentile. Therefore, the background concentration of iron in groundwater at the Olalla Landfill was calculated to be $10 \mu\text{g/L}$. These

results were verified by hand calculation as shown in Example 5 of the Ecology guidance document (1992). A copy of the spreadsheets, including data and calculations, are attached to this memorandum.

IRON

The iron data for MW-1 (n=14) are neither log normally or normally distributed ($r^2 = 0.50$ and 0.48 respectively) with a median = $20 \mu\text{g/L}$. The iron data for MW-5A (n=14) could not be analyzed for log normal or normal distribution because all values but one were identical and all were non-detects. The MW-5A data have a median = $20 \mu\text{g/L}$.

These median data were combined as a data set with the iron values for the off-site wells. The 90th and 50th percentiles were calculated using the non-parametric method as described by Ecology (1992). The values of the percentiles are N/A and 10.00 respectively and the CV is N/A. Non-parametric estimates of the 90th percentile and the CV cannot be calculated due to the small size of the data set. Although the 90th percentile cannot be estimated, that value must be greater than 106 (the highest value and rank of the iron data set). The Ecology guidance (1992) indicates that the background value should be set at four times the 50th percentile when the 90th percentile is greater than four times the 50th percentile. Therefore the background concentration of iron in groundwater at the Olalla Landfill was calculated to be $40 \mu\text{g/L}$. These results were verified by hand calculation as shown in Example 5 of the Ecology guidance document (1992). Copies of the spreadsheets, including data and calculations, are attached to this memorandum.

REFERENCES

Blakely, N. 1992. Statistical Guidance for Ecology Site Managers. Washington Department of Ecology. Olympia, Washington.

Attachments

Attachment A
Distribution and Summary Statistics

Background calculations

20 3/23/2009 **MW-1 Iron_ug/L (2009-2012)**

20 6/2/2009
 20 9/30/2009
 20 12/8/2009
 20 12/28/2010
 20 3/23/2011
 20 6/1/2011
 20 9/26/2011
 20 12/5/2011
 20 3/6/2012
 20 6/21/2012
 25.1 10/28/2010
 47 3/3/2010
 50 6/9/2010

| MTCASat 3.0 | | | |
|----------------------------------|----|----------------------|--|
| Number of samples | | Uncensored values | |
| Uncensored | 14 | Mean | 24.44 |
| Censored | 0 | Lognormal mean | 24.28 |
| TOTAL | 14 | Std. devn. | 10.30 |
| | | Median | 20 |
| | | Min. | 20 |
| | | Max. | 50 |
| Lognormal distribution? | | Normal distribution? | |
| r-squared is: 0.50 | | r-squared is: 0.48 | |
| Recommendations: | | | |
| Use nonparametric method. | | | |
| Distribution selection | | Enter percentile | Value corresponding to that percentile is: |
| 1 | | 90 | |
| 1 = Lognormal | | | 50th |
| 2 = Normal | | | 4 X 50th |
| 3 = Nonparametric method | | | |

Background calculations

5 3/23/2009 **MW-1_ug/L Manganese (2009-2012)**

5 6/2/2009

5 9/30/2009

5 12/8/2009

5 3/3/2010

5 6/9/2010

5 6/1/2011

5 12/5/2011

5 3/6/2012

5 6/21/2012

5 10/28/2010

5 3/23/2011

10 12/28/2010

10 9/26/2011

| MTCASat 3.0 | | | | |
|-------------------------------------|----|----------------------|--|--|
| Number of samples | | Uncensored values | | |
| Uncensored | 14 | Mean | 5.71 | |
| Censored | 0 | Lognormal mean | 5.70 | |
| TOTAL | 14 | Std. devn. | 1.82 | |
| | | Median | 5 | |
| | | Min. | 5 | |
| | | Max. | 10 | |
| Lognormal distribution? | | Normal distribution? | | |
| r-squared is: 0.42 | | r-squared is: 0.42 | | |
| Recommendations: | | | | |
| Use nonparametric method. | | | | |
| Distribution selection | | Enter percentile | Value corresponding to that percentile is: | |
| 1 | | 90 | | |
| 1 = Lognormal | | | 50th | |
| 2 = Normal | | | 4 X 50th | |
| 3 = Nonparametric method | | | | |

Background calculations

10 10/28/2010 **MW-5A_ug/L Iron (2009-2012)**

20 3/23/2009

20 6/2/2009

20 9/30/2009

20 12/8/2009

20 3/3/2010

20 6/9/2010

20 12/28/2010

20 3/23/2011

20 6/2/2011

20 9/26/2011

20 12/5/2011

20 3/6/2012

| MTCASat 3.0 | | | | | |
|---|----|-------------------|--|--|--|
| Number of samples | | Uncensored values | | | |
| Uncensored | 13 | Mean | 19.23 | | |
| Censored | 0 | Lognormal mean | 19.32 | | |
| TOTAL | 13 | Std. devn. | 2.77 | | |
| | | Median | 20 | | |
| | | Min. | 10 | | |
| | | Max. | 20 | | |
| Lognormal distribution? | | | Normal distribution? | | |
| r-squared is: | | | r-squared is: | | |
| Recommendations: | | | | | |
| Unable to analyze probability plot for lognormal case. Unable to analyze probability plot for normal case. If data are OK, use nonparametric method. | | | | | |
| Distribution selection | | | | | |
| | | Enter percentile | Value corresponding to that percentile is: | | |
| 1 | | 90 | | | |
| 1 = Lognormal | | | 50th | | |
| 2 = Normal | | | 4 X 50th | | |
| 3 = Nonparametric method | | | | | |

Background calculations

5 3/23/2009 **MW-5A_ug/L Manganese (2009-2012)**

5 6/2/2009

5 9/30/2009

5 12/8/2009

5 3/3/2010

5 6/9/2010

5 12/28/2010

5 6/2/2011

5 9/26/2011

5 12/5/2011

5 3/6/2012

10 10/28/2010

10 3/23/2011

| MTCASat 3.0 | | | | |
|----------------------------------|----|----------------------|--|--|
| Number of samples | | Uncensored values | | |
| Uncensored | 13 | Mean | 5.77 | |
| Censored | 0 | Lognormal mean | 5.75 | |
| TOTAL | 13 | Std. devn. | 1.88 | |
| | | Median | 5 | |
| | | Min. | 5 | |
| | | Max. | 10 | |
| Lognormal distribution? | | Normal distribution? | | |
| r-squared is: 0.44 | | r-squared is: 0.44 | | |
| Recommendations: | | | | |
| Use nonparametric method. | | | | |
| Distribution selection | | Enter percentile | Value corresponding to that percentile is: | |
| 1 | | 90 | 50th | |
| 1 = Lognormal | | | 4 X 50th | |
| 2 = Normal | | | | |
| 3 = Nonparametric method | | | | |

Attachment B
Background Calculations Using *MTCASat*

Background calculations

10 MW-1
 10 MW-5A
 10 OW1
 10 OW2
 71 OW9
 106 OW4

Calculation of Iron Background for Olalla Landfill (NDs = [1/2 MDL or RL])

| MTCASat 3.0 | | | |
|----------------------------------|---|---------------------------------------|--|
| Number of samples | | Uncensored values | |
| Uncensored | 6 | Mean | 36.17 |
| Censored | 0 | Lognormal mean | 38.60 |
| TOTAL | 6 | Std. devn. | 42.02 |
| | | Median | 10 |
| | | Min. | 10 |
| | | Max. | 106 |
| Lognormal distribution? | | Normal distribution? | |
| r-squared is: 0.69 | | r-squared is: 0.71 | |
| Recommendations: | | | |
| Use nonparametric method. | | | |
| Distribution selection | | | |
| 3 | | Enter percentile | Value corresponding to that percentile is: |
| 1 = Lognormal | | 90 | N/A |
| 2 = Normal | | 50th | 10.00 |
| 3 = Nonparametric method | | 4 X 50th | 40.00 |
| | | Coefficient of Variation = N/A | |

Background calculations

- 2.5 MW-1
- 2.5 MW-5A
- 2.5 OW1
- 2.5 OW2
- 2.5 OW9
- 32 OW4

Calculation of Mn Background for Olalla Landfill (NDs = [1/2 MDL or RL])

| MTCASat 3.0 | | | |
|---|---|---------------------------------------|--|
| Number of samples | | Uncensored values | |
| Uncensored | 6 | Mean | 7.42 |
| Censored | 0 | Lognormal mean | 6.57 |
| TOTAL | 6 | Std. devn. | 12.04 |
| | | Median | 2.5 |
| | | Min. | 2.5 |
| | | Max. | 32 |
| Lognormal distribution? | | Normal distribution? | |
| r-squared is: | | r-squared is: | |
| Recommendations: | | | |
| Unable to analyze probability plot for lognormal case. Unable to analyze probability plot for normal case. If data are OK, use nonparametric method. | | | |
| Distribution selection | | Enter percentile | Value corresponding to that percentile is: |
| 3 | | 90 | N/A |
| 1 = Lognormal | | 50th | 2.50 |
| 2 = Normal | | 4 X 50th | 10.00 |
| 3 = Nonparametric method | | Coefficient of Variation = N/A | |

Appendix I

Groundwater and Surface Water
Laboratory Data Sheets



Aquatic Research Inc.
3927 Aurora Ave. N., Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-1 | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-1 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 0601006.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 79% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 105% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 92% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



Aquatic Research Inc.
3927 Aurora Ave. N., Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A3** Matrix: Water
Sample ID No.: **OL-MW-2** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 1001010.D**

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-2 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 1001010.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 77% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 108% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 86% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A4** Matrix: Water
Sample ID No.: **OL-MW-4** Sample Wt./Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 1101011.D**

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-4 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 1101011.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 72% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 104% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 84% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-5A | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



Aquatic Research Inc.
3927 Aurora Ave. N., Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-5A | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 0801008.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------------------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| Surrogate Recoveries | | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 80% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 106% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 92% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



Aquatic Research Inc.
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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A2 Dup** Matrix: Water
Sample ID No.: **OL-MW-5A Duplicate** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 0901009.D**

| CAS# | Name of Compound | Amount (ppb) | Flag |
|-----------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01- | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02- | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A2 Dup | Matrix: | Water |
| Sample ID No.: | OL-MW-5A Duplicate | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 0901009.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 70% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 106% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 84% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A7** Matrix: Water
Sample ID No.: **OL-ER** Sample Wt./Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 1201012.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A7 | Matrix: | Water |
| Sample ID No.: | OL-ER | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 1201012.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------------------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| Surrogate Recoveries | | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 72% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 109% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 88% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A1 MS | Matrix: | Water |
| Sample ID No.: | BKS00147A1 MS | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | | |
| Date Analyzed: | 03/31/11 | Analyst: | T. Meadows |
| Date of Report: | 3/31/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | 0701007.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 79% | 66% | 118% |
| Toluene-d8 | 110% | 51% | 143% |
| 4-Bromofluorobenzene | 86% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 102% | 75% | 125% |
| Benzene | 105% | 75% | 125% |
| Trichloroethene | 112% | 75% | 125% |
| Toluene | 112% | 75% | 125% |
| Chlorobenzene | 113% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | 3/31/11-LCS | Matrix: | Water |
| Sample ID No.: | 3/31/11-LCS | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed | 03/31/11 | Analyst: | T. Meadows |
| Date of Report: | 3/31/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 108% | 66% | 118% |
| Toluene-d8 | 107% | 51% | 143% |
| 4-Bromofluorobenzene | 107% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 103% | 75% | 125% |
| Benzene | 102% | 75% | 125% |
| Trichloroethene | 102% | 75% | 125% |
| Toluene | 102% | 75% | 125% |
| Chlorobenzene | 107% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | 3/31/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | 3/31/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 0401004.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 106% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 104% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 106% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A8** Matrix: Water
Sample ID No.: **OL-TB** Sample Wt./Vol. (gm/ml) **25.0**
Date Collected: **03/15/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 1301013.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A8 | Matrix: | Water |
| Sample ID No.: | OL-TB | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/15/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 1301013.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 74% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 106% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 83% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A1**
Sample ID No.: **OL-MW-1**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1050**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **0601006.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.76 | U | 86-73-7 | Fluorene | 0.76 | U |
| 76-01-7 | Pentachloroethane | 0.76 | U | 84-66-2 | Diethyl phthalate | 0.76 | U |
| 62-53-3 | Aniline | 0.76 | U | 100-01-6 | 4-Nitroaniline | 0.76 | U |
| 108-95-2 | Phenol | 0.76 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.76 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.76 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.76 | U |
| 95-57-8 | 2-Chlorophenol | 0.76 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.76 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 103-33-3 | Azobenzene | 0.76 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.76 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.76 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.76 | U | 319-84-6 | a-BHC | 0.76 | U |
| 100-51-6 | Benzyl Alcohol | 0.76 | U | 118-74-1 | Hexachlorobenzene | 0.76 | U |
| 95-48-7 | 2-Methyl phenol | 0.76 | U | 319-85-7 | b-BHC | 0.76 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.76 | U | 87-86-5 | Pentachlorophenol | 0.76 | U |
| 98-86-2 | Acetophenone | 0.76 | U | 58-89-9 | g-BHC (Lindane) | 0.76 | U |
| 67-72-1 | Hexachloroethane | 0.76 | U | 85-01-8 | Phenanthrene | 0.76 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.76 | U | 120-12-7 | Anthracene | 0.76 | U |
| 106-44-5 | 4-Methyl phenol | 0.76 | U | 319-86-8 | d-BHC | 0.76 | U |
| 98-95-3 | Nitrobenzene | 0.76 | U | 84-74-2 | Di-n-butyl phthalate | 0.76 | U |
| 100-75-4 | N-nitrosopiperidine | 0.76 | U | 76-44-8 | Heptachlor | 0.76 | U |
| 78-59-1 | Isophorone | 0.76 | U | 309-00-2 | Aldrin | 0.76 | U |
| 88-75-5 | 2-Nitrophenol | 0.76 | U | 1024-57-3 | Heptachlor epoxide | 0.76 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.76 | U | 206-44-0 | Fluoranthrene | 0.76 | U |
| 65-85-0 | Benzoic Acid | 0.76 | U | 129-00-0 | Pyrene | 0.76 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.76 | U | 959-98-8 | Endosulfan I | 0.76 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.76 | U | 92-87-5 | Benzidine | 0.76 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.76 | U | 72-55-9 | 4,4'-DDE | 0.76 | U |
| 91-20-3 | Naphthalene | 0.76 | U | 60-57-1 | Dieldrin | 0.76 | U |
| 106-47-8 | 4-Chloroaniline | 0.76 | U | 72-20-8 | Endrin | 0.76 | U |
| 87-68-3 | Hexachlorobutadiene | 0.76 | U | 33213-65-9 | Endosulfan II | 0.76 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.76 | U | 72-54-8 | 4,4'-DDD | 0.76 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.76 | U | 7421-92-4 | Endrin aldehyde | 0.76 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.76 | U | 85-68-7 | Butyl benzenyl phthalate | 0.76 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.76 | U | 1031-07-8 | Endosulfan sulfate | 0.76 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.76 | U | 50-29-3 | 4,4'-DDT | 0.76 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.76 | U | 56-55-3 | Benzo(a)anthracene | 0.76 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.76 | U | 218-01-9 | Chrysene | 0.76 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.76 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.76 | U |
| 88-74-4 | 2-Nitroaniline | 0.76 | U | 72-43-5 | Methoxychlor | 0.76 | U |
| 131-11-3 | Dimethyl phthalate | 0.76 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.76 | U |
| 208-96-8 | Acenaphthylene | 0.76 | U | 117-84-0 | Di-n-octyl phthalate | 0.76 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.76 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.76 | U |
| 99-09-2 | 3-Nitroaniline | 0.76 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.76 | U |
| 83-32-9 | Acenaphthene | 0.76 | U | 50-32-8 | Benzo(a)pyrene | 0.76 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.76 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.76 | U |
| 132-64-9 | Dibenzofuran | 0.76 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.76 | U |
| 100-02-7 | 4-Nitrophenol | 0.76 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.76 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.76 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.76 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-1 | Sample Vol. (ml) | 1050 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 50% | 24-132% | 65-135% |
| Phenol-d5 | 35% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 99% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 81% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 43% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 125% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.76 | U |
| 100-52-7 | Benzaldehyde | 0.76 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.76 | U |
| 95-53-4 | o-Toluidine | 0.76 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.76 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.76 | U |
| 1888-71-7 | Hexachloropropene | 0.76 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.76 | U |
| 87-65-0 | Caprolactam | 0.76 | U |
| 106-50-3 | p-Phenylenediamine | 0.76 | U |
| 120-58-1 | Isosafrole | 0.76 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.76 | U |
| 94-59-7 | Safrole | 0.76 | U |
| 92-52-4 | Biphenyl | 0.76 | U |
| 101-84-8 | Diphenyl ether | 0.76 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.76 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.76 | U |
| 608-93-5 | Pentachlorobenzene | 0.76 | U |
| 134-32-7 | 1-Naphthalenamine | 0.76 | U |
| 91-59-8 | 2-Naphthalenamine | 0.76 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.76 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.76 | U |
| 122-39-4 | Diphenylamine | 0.76 | U |
| 3689-24-5 | Sulfotep | 0.76 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.76 | U |
| 298-02-2 | Phorate | 0.76 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.76 | U |
| 62-44-2 | Phenacetin | 0.76 | U |
| 60-51-5 | Dimethoate | 0.76 | U |
| 1912-24-9 | Atrazine | 0.76 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.76 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.76 | U |
| 23950-58-! | Pronamide | 0.76 | U |
| 298-04-4 | Disulfoton | 0.76 | U |
| 298-00-0 | Methyl parathion | 0.76 | U |
| 56-38-2 | Parathion | 0.76 | U |
| 91-80-5 | Methapyrilene | 0.76 | U |
| 465-73-6 | Isodrin | 0.76 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.76 | U |
| 510-15-0 | Chlorobenzilate | 0.76 | U |
| 143-50-0 | Kepone | 0.76 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.76 | U |
| 52-85-7 | Famphur | 0.76 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.76 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.76 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.76 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.76 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A3**
Sample ID No.: **OL-MW-2**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1060**
Final Volume (ml) **2.0**
Dilution Factor: **1**
Analyst: **T. Meadows**
Supervisor's Initials: **0801008.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-2 | Sample Vol. (ml) | 1060 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 73% | 24-132% | 65-135% |
| Phenol-d5 | 51% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 108% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 82% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 46% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 137% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A4**
Sample ID No.: **OL-MW-4**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **0901009.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-4 | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 50% | 24-132% | 65-135% |
| Phenol-d5 | 43% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 110% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 96% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 38% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 117% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A2**
Sample ID No.: **OL-MW-5A**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials:
0701007.D

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-5A | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | 0701007.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 51% | 24-132% | 65-135% |
| Phenol-d5 | 43% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 98% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 78% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 37% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 102% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A7**
Sample ID No.: **OL-ER**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1040**
Final Volume (ml) **2.0**
Dilution Factor: **1**
Analyst: **T. Meadows**
Supervisor's Initials: **1001010.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.77 | U | 86-73-7 | Fluorene | 0.77 | U |
| 76-01-7 | Pentachloroethane | 0.77 | U | 84-66-2 | Diethyl phthalate | 0.77 | U |
| 62-53-3 | Aniline | 0.77 | U | 100-01-6 | 4-Nitroaniline | 0.77 | U |
| 108-95-2 | Phenol | 0.77 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.77 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.77 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.77 | U |
| 95-57-8 | 2-Chlorophenol | 0.77 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.77 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.77 | U | 103-33-3 | Azobenzene | 0.77 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.77 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.77 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.77 | U | 319-84-6 | a-BHC | 0.77 | U |
| 100-51-6 | Benzyl Alcohol | 0.77 | U | 118-74-1 | Hexachlorobenzene | 0.77 | U |
| 95-48-7 | 2-Methyl phenol | 0.77 | U | 319-85-7 | b-BHC | 0.77 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.77 | U | 87-86-5 | Pentachlorophenol | 0.77 | U |
| 98-86-2 | Acetophenone | 0.77 | U | 58-89-9 | g-BHC (Lindane) | 0.77 | U |
| 67-72-1 | Hexachloroethane | 0.77 | U | 85-01-8 | Phenanthrene | 0.77 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.77 | U | 120-12-7 | Anthracene | 0.77 | U |
| 106-44-5 | 4-Methyl phenol | 0.77 | U | 319-86-8 | d-BHC | 0.77 | U |
| 98-95-3 | Nitrobenzene | 0.77 | U | 84-74-2 | Di-n-butyl phthalate | 0.77 | U |
| 100-75-4 | N-nitrosopiperidine | 0.77 | U | 76-44-8 | Heptachlor | 0.77 | U |
| 78-59-1 | Isophorone | 0.77 | U | 309-00-2 | Aldrin | 0.77 | U |
| 88-75-5 | 2-Nitrophenol | 0.77 | U | 1024-57-3 | Heptachlor epoxide | 0.77 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.77 | U | 206-44-0 | Fluoranthrene | 0.77 | U |
| 65-85-0 | Benzoic Acid | 0.77 | U | 129-00-0 | Pyrene | 0.77 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.77 | U | 959-98-8 | Endosulfan I | 0.77 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.77 | U | 92-87-5 | Benzidine | 0.77 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.77 | U | 72-55-9 | 4,4'-DDE | 0.77 | U |
| 91-20-3 | Naphthalene | 0.77 | U | 60-57-1 | Dieldrin | 0.77 | U |
| 106-47-8 | 4-Chloroaniline | 0.77 | U | 72-20-8 | Endrin | 0.77 | U |
| 87-68-3 | Hexachlorobutadiene | 0.77 | U | 33213-65-9 | Endosulfan II | 0.77 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.77 | U | 72-54-8 | 4,4'-DDD | 0.77 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.77 | U | 7421-92-4 | Endrin aldehyde | 0.77 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.77 | U | 85-68-7 | Butyl benzenyl phthalate | 0.77 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.77 | U | 1031-07-8 | Endosulfan sulfate | 0.77 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.77 | U | 50-29-3 | 4,4'-DDT | 0.77 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.77 | U | 56-55-3 | Benzo(a)anthracene | 0.77 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.77 | U | 218-01-9 | Chrysene | 0.77 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.77 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.77 | U |
| 88-74-4 | 2-Nitroaniline | 0.77 | U | 72-43-5 | Methoxychlor | 0.77 | U |
| 131-11-3 | Dimethyl phthalate | 0.77 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.77 | U |
| 208-96-8 | Acenaphthylene | 0.77 | U | 117-84-0 | Di-n-octyl phthalate | 0.77 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.77 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.77 | U |
| 99-09-2 | 3-Nitroaniline | 0.77 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.77 | U |
| 83-32-9 | Acenaphthene | 0.77 | U | 50-32-8 | Benzo(a)pyrene | 0.77 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.77 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.77 | U |
| 132-64-9 | Dibenzofuran | 0.77 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.77 | U |
| 100-02-7 | 4-Nitrophenol | 0.77 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.77 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.77 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.77 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A7 | Matrix: | Water |
| Sample ID No.: | OL-ER | Sample Vol. (ml) | 1040 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | 1001010.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 53% | 24-132% | 65-135% |
| Phenol-d5 | 49% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 96% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 84% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 28% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 96% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.77 | U |
| 100-52-7 | Benzaldehyde | 0.77 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.77 | U |
| 95-53-4 | o-Toluidine | 0.77 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.77 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.77 | U |
| 1888-71-7 | Hexachloropropene | 0.77 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.77 | U |
| 87-65-0 | Caprolactam | 0.77 | U |
| 106-50-3 | p-Phenylenediamine | 0.77 | U |
| 120-58-1 | Isosafrole | 0.77 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.77 | U |
| 94-59-7 | Safrole | 0.77 | U |
| 92-52-4 | Biphenyl | 0.77 | U |
| 101-84-8 | Diphenyl ether | 0.77 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.77 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.77 | U |
| 608-93-5 | Pentachlorobenzene | 0.77 | U |
| 134-32-7 | 1-Naphthalenamine | 0.77 | U |
| 91-59-8 | 2-Naphthalenamine | 0.77 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.77 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.77 | U |
| 122-39-4 | Diphenylamine | 0.77 | U |
| 3689-24-5 | Sulfotep | 0.77 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.77 | U |
| 298-02-2 | Phorate | 0.77 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.77 | U |
| 62-44-2 | Phenacetin | 0.77 | U |
| 60-51-5 | Dimethoate | 0.77 | U |
| 1912-24-9 | Atrazine | 0.77 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.77 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.77 | U |
| 23950-58-! | Pronamide | 0.77 | U |
| 298-04-4 | Disulfoton | 0.77 | U |
| 298-00-0 | Methyl parathion | 0.77 | U |
| 56-38-2 | Parathion | 0.77 | U |
| 91-80-5 | Methapyrilene | 0.77 | U |
| 465-73-6 | Isodrin | 0.77 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.77 | U |
| 510-15-0 | Chlorobenzilate | 0.77 | U |
| 143-50-0 | Kepone | 0.77 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.77 | U |
| 52-85-7 | Famphur | 0.77 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.77 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.77 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.77 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.77 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
10/14/10chy/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | 3/25/11-LCS | Matrix: | Water |
| Sample ID No.: | Lab Control Spike | Sample Vol. (ml) | 1000 |
| Date Collected: | n/a | Final Volume (ml) | 2.0 |
| Date Received: | n/a | Dilution Factor: | 1 |
| Date Analyzed: | 04/06/11 | Analyst: | T. Meadows |
| Date of Report: | 04/08/11 | Supervisor's Initials: | |
| Data File Path: | Q:\5973\8270\110406\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 110% | 24-132% | 65-135% |
| Phenol-d5 | 68% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 87% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 76% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 37% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 139% | 80-141% | 65-135% |

| Matix Spike Recoveries | %Rec. | QC limits | |
|--------------------------|-------|-----------|---------|
| | | Water | Soil |
| Phenol | 66% | 50-150% | 50-150% |
| 2-Chlorophenol | 77% | 50-150% | 50-150% |
| 1,4-Dichlorobenzene | 60% | 50-150% | 50-150% |
| N-Nitroso-n-propyl amine | 57% | 50-150% | 50-150% |
| 1,2,4-Trichlorobenzene | 59% | 50-150% | 50-150% |
| 4-Chloro-3-methyl phenol | 77% | 50-150% | 50-150% |
| Acenaphthene | 82% | 50-150% | 50-150% |
| 4-Nitrophenol | 47% | 50-150% | 50-150% |
| 2,4-Dinitrotoluene | 50% | 50-150% | 50-150% |
| Pentachlorophenol | 0% | 50-150% | 50-150% |
| Pyrene | 115% | 50-150% | 50-150% |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **3/25/11-MB**
Sample ID No.: **Method Blank**
Date Collected: **n/a**
Date Received: **n/a**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1000**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **0401004.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.80 | U | 86-73-7 | Fluorene | 0.80 | U |
| 76-01-7 | Pentachloroethane | 0.80 | U | 84-66-2 | Diethyl phthalate | 0.80 | U |
| 62-53-3 | Aniline | 0.80 | U | 100-01-6 | 4-Nitroaniline | 0.80 | U |
| 108-95-2 | Phenol | 0.80 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.80 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.80 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.80 | U |
| 95-57-8 | 2-Chlorophenol | 0.80 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.80 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 103-33-3 | Azobenzene | 0.80 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.80 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.80 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.80 | U | 319-84-6 | a-BHC | 0.80 | U |
| 100-51-6 | Benzyl Alcohol | 0.80 | U | 118-74-1 | Hexachlorobenzene | 0.80 | U |
| 95-48-7 | 2-Methyl phenol | 0.80 | U | 319-85-7 | b-BHC | 0.80 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.80 | U | 87-86-5 | Pentachlorophenol | 0.80 | U |
| 98-86-2 | Acetophenone | 0.80 | U | 58-89-9 | g-BHC (Lindane) | 0.80 | U |
| 67-72-1 | Hexachloroethane | 0.80 | U | 85-01-8 | Phenanthrene | 0.80 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.80 | U | 120-12-7 | Anthracene | 0.80 | U |
| 106-44-5 | 4-Methyl phenol | 0.80 | U | 319-86-8 | d-BHC | 0.80 | U |
| 98-95-3 | Nitrobenzene | 0.80 | U | 84-74-2 | Di-n-butyl phthalate | 0.80 | U |
| 100-75-4 | N-nitrosopiperidine | 0.80 | U | 76-44-8 | Heptachlor | 0.80 | U |
| 78-59-1 | Isophorone | 0.80 | U | 309-00-2 | Aldrin | 0.80 | U |
| 88-75-5 | 2-Nitrophenol | 0.80 | U | 1024-57-3 | Heptachlor epoxide | 0.80 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.80 | U | 206-44-0 | Fluoranthrene | 0.80 | U |
| 65-85-0 | Benzoic Acid | 0.80 | U | 129-00-0 | Pyrene | 0.80 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.80 | U | 959-98-8 | Endosulfan I | 0.80 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.80 | U | 92-87-5 | Benzidine | 0.80 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.80 | U | 72-55-9 | 4,4'-DDE | 0.80 | U |
| 91-20-3 | Naphthalene | 0.80 | U | 60-57-1 | Dieldrin | 0.80 | U |
| 106-47-8 | 4-Chloroaniline | 0.80 | U | 72-20-8 | Endrin | 0.80 | U |
| 87-68-3 | Hexachlorobutadiene | 0.80 | U | 33213-65-9 | Endosulfan II | 0.80 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.80 | U | 72-54-8 | 4,4'-DDD | 0.80 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.80 | U | 7421-92-4 | Endrin aldehyde | 0.80 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.80 | U | 85-68-7 | Butyl benzenyl phthalate | 0.80 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.80 | U | 1031-07-8 | Endosulfan sulfate | 0.80 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.80 | U | 50-29-3 | 4,4'-DDT | 0.80 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.80 | U | 56-55-3 | Benzo(a)anthracene | 0.80 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.80 | U | 218-01-9 | Chrysene | 0.80 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.80 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.80 | U |
| 88-74-4 | 2-Nitroaniline | 0.80 | U | 72-43-5 | Methoxychlor | 0.80 | U |
| 131-11-3 | Dimethyl phthalate | 0.80 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.80 | U |
| 208-96-8 | Acenaphthylene | 0.80 | U | 117-84-0 | Di-n-octyl phthalate | 0.80 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.80 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.80 | U |
| 99-09-2 | 3-Nitroaniline | 0.80 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.80 | U |
| 83-32-9 | Acenaphthene | 0.80 | U | 50-32-8 | Benzo(a)pyrene | 0.80 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.80 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.80 | U |
| 132-64-9 | Dibenzofuran | 0.80 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.80 | U |
| 100-02-7 | 4-Nitrophenol | 0.80 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.80 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.80 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.80 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | 3/25/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Vol. (ml) | 1000 |
| Date Collected: | n/a | Final Volume (ml) | 2.0 |
| Date Received: | n/a | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purges Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 50% | 24-132% | 65-135% |
| Phenol-d5 | 47% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 92% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 86% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 41% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 118% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.80 | U |
| 100-52-7 | Benzaldehyde | 0.80 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.80 | U |
| 95-53-4 | o-Toluidine | 0.80 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.80 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.80 | U |
| 1888-71-7 | Hexachloropropene | 0.80 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.80 | U |
| 87-65-0 | Caprolactam | 0.80 | U |
| 106-50-3 | p-Phenylenediamine | 0.80 | U |
| 120-58-1 | Isosafrole | 0.80 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.80 | U |
| 94-59-7 | Safrole | 0.80 | U |
| 92-52-4 | Biphenyl | 0.80 | U |
| 101-84-8 | Diphenyl ether | 0.80 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.80 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.80 | U |
| 608-93-5 | Pentachlorobenzene | 0.80 | U |
| 134-32-7 | 1-Naphthalenamine | 0.80 | U |
| 91-59-8 | 2-Naphthalenamine | 0.80 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.80 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.80 | U |
| 122-39-4 | Diphenylamine | 0.80 | U |
| 3689-24-5 | Sulfotep | 0.80 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.80 | U |
| 298-02-2 | Phorate | 0.80 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.80 | U |
| 62-44-2 | Phenacetin | 0.80 | U |
| 60-51-5 | Dimethoate | 0.80 | U |
| 1912-24-9 | Atrazine | 0.80 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.80 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.80 | U |
| 23950-58-! | Pronamide | 0.80 | U |
| 298-04-4 | Disulfoton | 0.80 | U |
| 298-00-0 | Methyl parathion | 0.80 | U |
| 56-38-2 | Parathion | 0.80 | U |
| 91-80-5 | Methapyrilene | 0.80 | U |
| 465-73-6 | Isodrin | 0.80 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.80 | U |
| 510-15-0 | Chlorobenzilate | 0.80 | U |
| 143-50-0 | Kepone | 0.80 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.80 | U |
| 52-85-7 | Famphur | 0.80 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.80 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.80 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.80 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.80 | U |



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CASE FILE NUMBER: BKS001-47,48 PAGE 1
REPORT DATE: 04/06/11
DATE SAMPLED: 03/23,24/11 DATE RECEIVED: 03/24,25/11
FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER
SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL

CASE NARRATIVE

Fifteen water samples were received by the laboratory in good condition. Analysis was performed according to the chain of custody received with the samples. Dissolved arsenic analysis was performed by ICP/MS by concentrating the samples 10 fold. No difficulties were encountered in the preparation or analysis of these samples. Sample data follows while QA/QC data is contained on subsequent pages. Organic data are included as separate reports.

SAMPLE DATA

| SAMPLE ID | AMMONIA (mg/L) | NITRATE (mg/L) | NITRITE (mg/L) | TOC (mg/L) | COD (mg/L) | VINYL CHLORIDE (ug/L) |
|-----------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| OL-MW-1 | <0.010 | 0.254 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-MW-5A | <0.010 | 0.192 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-MW-2 | <0.010 | 0.389 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-MW-4 | <0.010 | 0.536 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-SW-3 | | 0.080 | 0.002 | | | |
| OL-SW-4 | | 0.147 | 0.002 | | | |
| OL-ER | <0.010 | <0.010 | <0.002 | <0.250 | <10.0 | 0.05 |
| OL-TB | | | | | | 0.09 |
| OL-MW-7 | 0.012 | 0.549 | <0.002 | 0.370 | <10.0 | <0.02 |
| OL-MW-8 | <0.010 | 0.131 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-MW-6 | <0.010 | 0.011 | <0.002 | 1.14 | <10.0 | <0.02 |
| OL-MW-10 | 0.017 | 0.292 | <0.002 | 3.04 | <10.0 | <0.02 |
| OL-MW-3 | <0.010 | 0.200 | 0.003 | 1.28 | <10.0 | <0.02 |
| OL-TB | | | | | | 0.05 |
| OL-MW-11 | <0.010 | 0.550 | <0.002 | 0.400 | <10.0 | <0.02 |



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CASE FILE NUMBER: BKS001-47,48 **PAGE 2**
REPORT DATE: 04/06/11
DATE SAMPLED: 03/23,24/11 03/24,25/11
FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER
SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL

SAMPLE DATA-CONTINUED

| SAMPLE ID | CHLORIDE (mg/L) | SULFATE (mg/L) | TDS (mg/L) | CYANIDE (mg/L) | TOTAL COLIFORM (#/100mls) | FECAL COLIFORM (#/100mls) |
|-----------|--------------------|-------------------|---------------|-------------------|------------------------------|------------------------------|
| OL-MW-1 | 2.64 | 4.03 | 85.0 | <0.010 | <2 | |
| OL-MW-5A | 2.05 | 3.82 | 148 | <0.010 | <2 | |
| OL-MW-2 | 2.15 | 2.62 | 100 | <0.010 | 2 | |
| OL-MW-4 | 2.54 | 2.46 | 101 | <0.010 | 4 | |
| OL-SW-3 | | | | | | 2 |
| OL-SW-4 | | | | | | <2 |
| OL-ER | <0.50 | <1.00 | 21.0 | <0.010 | 6 | |
| OL-MW-7 | 1.25 | 3.08 | 109 | <0.010 | <2 | |
| OL-MW-8 | 1.17 | 1.42 | 118 | <0.010 | <2 | |
| OL-MW-6 | 1.25 | 5.84 | 109 | <0.010 | <2 | |
| OL-MW-10 | 3.07 | 27.0 | 326 | <0.010 | <2 | |
| OL-MW-3 | 2.74 | 49.2 | 212 | <0.010 | <2 | |
| OL-MW-11 | 1.27 | 4.32 | 103 | <0.010 | <2 | |

| | pH | ALKALINITY (mgCaCO3/L) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|----------|------|---------------------------|---------------------|--------------------|
| OL-MW-1 | 6.58 | 60.8 | 38.3 | <1.00 |
| OL-MW-5A | 6.92 | 72.7 | 57.3 | <1.00 |
| OL-MW-2 | 7.13 | 55.2 | 47.3 | <1.00 |
| OL-MW-4 | 7.17 | 68.1 | 59.1 | <1.00 |
| OL-ER | 6.23 | <1.00 | <1.00 | <1.00 |
| OL-MW-7 | 6.98 | 67.0 | 54.3 | <1.00 |
| OL-MW-8 | 7.01 | 59.0 | 48.4 | <1.00 |
| OL-MW-6 | 6.92 | 67.5 | 53.2 | <1.00 |
| OL-MW-10 | 6.58 | 257 | 162 | <1.00 |
| OL-MW-3 | 6.36 | 134 | 67.8 | <1.00 |
| OL-MW-11 | 6.96 | 65.1 | 52.2 | <1.00 |



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| | | |
|--|---------------------|--------------------|
| CASE FILE NUMBER: | BKS001-47,48 | PAGE 3 |
| REPORT DATE: | 04/06/11 | |
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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

SAMPLE DATA-CONTINUED

| DISSOLVED METALS | | | | | | |
|------------------|-------------------|--------------------|------------------|---------------------|-------------------|--------------------|
| SAMPLE ID | ARSENIC (ug/L) | ANTIMONY (mg/L) | BARIUM (mg/L) | BERYLLIUM (mg/L) | CADMIUM (mg/L) | CHROMIUM (mg/L) |
| OL-MW-1 | 0.082 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0023 |
| OL-MW-5A | 0.113 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0036 |
| OL-MW-2 | 0.517 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0034 |
| OL-MW-4 | 0.188 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0024 |
| OL-MW-7 | 0.327 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0035 |
| OL-MW-8 | 1.49 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0013 |
| OL-MW-6 | 0.689 | <0.0020 | <0.005 | <0.0010 | <0.0010 | <0.0010 |
| OL-MW-10 | 1.03 | <0.0020 | 0.016 | <0.0010 | <0.0010 | 0.0020 |
| OL-MW-3 | 0.087 | <0.0020 | 0.012 | <0.0010 | <0.0010 | 0.0014 |
| OL-MW-11 | 0.349 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0034 |

| DISSOLVED METALS | | | | | | |
|------------------|------------------|------------------|----------------|--------------------|------------------|--------------------|
| SAMPLE ID | COBALT (mg/L) | COPPER (mg/L) | LEAD (mg/L) | SELENIUM (mg/L) | SILVER (mg/L) | THALLIUM (mg/L) |
| OL-MW-1 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-5A | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-2 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-4 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-7 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-8 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-6 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-10 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-3 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-11 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |

| DISSOLVED METALS | | | |
|------------------|---------------|--------------------|----------------|
| SAMPLE ID | TIN (mg/L) | VANADIUM (mg/L) | ZINC (mg/L) |
| OL-MW-1 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-5A | <0.040 | 0.0030 | <0.0050 |
| OL-MW-2 | <0.040 | 0.0040 | <0.0050 |
| OL-MW-4 | <0.040 | 0.0032 | <0.0050 |
| OL-MW-7 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-8 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-6 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-10 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-3 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-11 | <0.040 | <0.0030 | <0.0050 |



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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

SAMPLE DATA-CONTINUED

| SAMPLE ID | DISSOLVED METALS | | | | | |
|-----------|-------------------|---------------------|------------------|---------------------|----------------|---------------------|
| | CALCIUM (mg/L) | POTASSIUM (mg/L) | SODIUM (mg/L) | MAGNESIUM (mg/L) | IRON (mg/L) | MANGANESE (mg/L) |
| OL-MW-1 | 8.96 | 0.711 | 4.05 | 5.84 | <0.020 | <0.010 |
| OL-MW-5A | 8.73 | 0.811 | 4.49 | 8.45 | <0.020 | <0.010 |
| OL-MW-2 | 6.76 | 1.06 | 4.02 | 6.01 | <0.020 | <0.010 |
| OL-MW-4 | 8.54 | 0.942 | 4.54 | 7.64 | <0.020 | <0.010 |
| OL-MW-7 | 8.69 | 0.818 | 3.83 | 7.00 | <0.020 | <0.010 |
| OL-MW-8 | 9.28 | 0.680 | 4.11 | 4.69 | <0.020 | 0.143 |
| OL-MW-6 | 8.88 | 0.890 | 4.82 | 6.43 | 0.316 | 0.412 |
| OL-MW-10 | 39.3 | 1.44 | 10.8 | 24.8 | 0.084 | 4.85 |
| OL-MW-3 | 31.7 | 0.871 | 6.44 | 11.7 | <0.020 | 1.33 |
| OL-MW-11 | 8.86 | 0.821 | 3.83 | 7.21 | <0.020 | <0.010 |

| SAMPLE ID | TOTAL METALS | |
|-----------|-------------------|------------------|
| | MERCURY (mg/L) | NICKEL (mg/L) |
| OL-MW-1 | <0.0002 | <0.0050 |
| OL-MW-5A | <0.0002 | <0.0050 |
| OL-MW-2 | <0.0002 | <0.0050 |
| OL-MW-4 | <0.0002 | <0.0050 |
| OL-ER | <0.0002 | <0.0050 |
| OL-MW-7 | <0.0002 | <0.0050 |
| OL-MW-8 | <0.0002 | <0.0050 |
| OL-MW-6 | <0.0002 | <0.0050 |
| OL-MW-10 | <0.0002 | 0.0227 |
| OL-MW-3 | <0.0002 | <0.0050 |
| OL-MW-11 | <0.0002 | <0.0050 |



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| CASE FILE NUMBER: | BKS001-47,48 | PAGE 5 |
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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | AMMONIA (mg/L) | NITRATE (mg/L) | NITRITE (mg/L) | TOC (mg/L) | COD (mg/L) | VINYL CHLORIDE (ug/L) |
|-----------------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| METHOD | EPA 350.1 | EPA 353.2 | EPA 354.1 | EPA 415.1 | EPA 410.2 | EPA 8260-SIM |
| DATE ANALYZED | 03/24,25/11 | 03/24,25/11 | 03/25/11 | 04/01/11 | 04/01/11 | 04/04/11 |
| DETECTION LIMIT | 0.006 | 0.005 | 0.001 | 0.100 | 6.00 | 0.01 |
| REPORTING LIMIT | 0.010 | 0.010 | 0.002 | 0.250 | 10.0 | 0.02 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-10 | OL-ER | OL-MW-11 | OL-MW-11 | OL-MW-6 |
| ORIGINAL | 0.017 | 0.292 | <0.002 | 0.400 | <10.0 | <0.02 |
| DUPLICATE | 0.017 | 0.294 | <0.002 | 0.380 | <10.0 | <0.02 |
| RPD | 2.48% | 0.63% | NC | 5.07% | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-10 | OL-ER | OL-MW-11 | OL-MW-11 | |
| ORIGINAL | 0.017 | 0.292 | <0.002 | 0.400 | <10.0 | |
| SPIKED SAMPLE | 0.218 | 0.498 | 0.040 | 4.79 | 50.0 | |
| SPIKE ADDED | 0.200 | 0.200 | 0.040 | 4.50 | 50.0 | |
| % RECOVERY | 100.27% | 102.75% | 100.00% | 97.46% | 100.05% | NA |
| QC CHECK | | | | | | |
| FOUND | 0.313 | 0.400 | 0.041 | 4.12 | 94.4 | 0.10 |
| TRUE | 0.324 | 0.408 | 0.040 | 4.00 | 100 | 0.10 |
| % RECOVERY | 96.51% | 98.05% | 102.50% | 102.93% | 94.40% | 100.00% |
| BLANK | | | | | | |
| | <0.010 | <0.010 | <0.002 | <0.250 | <10.0 | <0.02 |

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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | CHLORIDE (mg/l) | SULFATE (mg/l) | TDS (mg/L) | CYANIDE (mg/L) | TOTAL COLIFORM (#/100mls) | FECAL COLIFORM (#/100mls) |
|-----------------|--------------------|-------------------|---------------|-------------------|------------------------------|------------------------------|
| METHOD | EPA 325.3 | EPA 375.4 | SM18 2540C | SM 4500CNE | SM18 9222B | SM18 9222D |
| DATE ANALYZED | 04/04/11 | 03/30/11 | 03/29/11 | 03/30/11 | 03/24,25/11 | 03/24/11 |
| DETECTION LIMIT | 0.16 | 0.76 | 5.0 | 0.010 | 2 | 2 |
| REPORTING LIMIT | 0.50 | 1.00 | 5.0 | 0.010 | 2 | 2 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-11 | OL-MW-10 | OL-MW-10 | OL-MW-10 | OL-SW-4 |
| ORIGINAL | 3.07 | 4.32 | 326 | <0.010 | <2 | <2 |
| DUPLICATE | 3.03 | 4.46 | 316 | <0.010 | <2 | <2 |
| RPD | 1.19% | 3.00% | 3.12% | NC | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-11 | | OL-MW-10 | | |
| ORIGINAL | 3.07 | 4.32 | | <0.010 | | |
| SPIKED SAMPLE | 12.7 | 14.5 | | 0.179 | | |
| SPIKE ADDED | 10.0 | 10.0 | | 0.200 | | |
| % RECOVERY | 96.41% | 101.53% | NA | 89.50% | NA | NA |
| QC CHECK | | | | | | |
| FOUND | 28.3 | 9.95 | | 0.184 | | |
| TRUE | 30.0 | 10.0 | | 0.200 | | |
| % RECOVERY | 94.49% | 99.51% | NA | 92.00% | NA | NA |
| PREP BLANK | | | | | | |
| | <0.50 | <1.00 | <5.0 | <0.010 | < 1 | < 1 |

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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | pH | ALKALINITY (mgCaCO3/l) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|-----------------|-------------|---------------------------|---------------------|--------------------|
| METHOD | EPA 150.1 | EPA 310.1 | EPA 310.1 | EPA 310.1 |
| DATE ANALYZED | 03/24,25/11 | 03/28/11 | 03/28/11 | 03/28/11 |
| DETECTION LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| REPORTING LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| DUPLICATE | | | | |
| SAMPLE ID | | OL-MW-10 | | |
| ORIGINAL | | 257 | | |
| DUPLICATE | | 265 | | |
| RPD | NA | 3.01% | NA | NA |
| SPIKE SAMPLE | | | | |
| SAMPLE ID | | | | |
| ORIGINAL | | | | |
| SPIKED SAMPLE | | | | |
| SPIKE ADDED | | | | |
| % RECOVERY | NA | NA | NA | NA |
| QC CHECK | | | | |
| FOUND | | 102 | | |
| TRUE | | 100 | | |
| % RECOVERY | NA | 101.50% | NA | NA |
| PREP BLANK | NA | NA | NA | NA |

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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | ARSENIC (ug/l) | ANTIMONY (mg/L) | BARIUM (mg/L) | BERYLLIUM (mg/L) | CADMIUM (mg/L) | CHROMIUM (mg/L) |
|-----------------|-------------------|--------------------|------------------|---------------------|-------------------|--------------------|
| METHOD | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 03/28,30/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 |
| DETECTION LIMIT | 0.010 | 0.0010 | 0.001 | 0.0003 | 0.0002 | 0.0005 |
| REPORTING LIMIT | 0.050 | 0.0020 | 0.005 | 0.0010 | 0.0010 | 0.0010 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | 1.03 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0024 |
| DUPLICATE | 1.07 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0025 |
| RPD | 3.81% | NC | NC | NC | NC | 1.30% |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | 1.03 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0024 |
| SPIKED SAMPLE | 1.21 | 0.0534 | 0.052 | 0.0555 | 0.0564 | 0.0491 |
| SPIKE ADDED | 0.200 | 0.0500 | 0.050 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 90.00% | 106.82% | 103.94% | 111.00% | 112.76% | 93.35% |
| QC CHECK | | | | | | |
| FOUND | 0.702 | 0.0489 | 0.048 | 0.0512 | 0.0470 | 0.0497 |
| TRUE | 0.668 | 0.0500 | 0.050 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 105.09% | 97.74% | 95.00% | 102.32% | 93.92% | 99.32% |
| PREP BLANK | | | | | | |
| | <0.050 | <0.0020 | <0.010 | <0.0010 | <0.0010 | <0.0010 |

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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | COBALT (mg/L) | COPPER (mg/L) | LEAD (mg/L) | SELENIUM (mg/L) | SILVER (mg/L) | THALLIUM (mg/L) |
|-----------------|------------------|------------------|----------------|--------------------|------------------|--------------------|
| METHOD | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 |
| DETECTION LIMIT | 0.001 | 0.0004 | 0.0003 | 0.0010 | 0.0004 | 0.0010 |
| REPORTING LIMIT | 0.010 | 0.0010 | 0.0010 | 0.0050 | 0.0010 | 0.0020 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| DUPLICATE | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| RPD | NC | NC | NC | NC | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| SPIKED SAMPLE | 0.048 | 0.0487 | 0.0467 | 0.0598 | 0.0513 | 0.0475 |
| SPIKE ADDED | 0.050 | 0.0500 | 0.0500 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 96.12% | 97.34% | 93.34% | 119.50% | 102.68% | 94.90% |
| QC CHECK | | | | | | |
| FOUND | 0.051 | 0.0515 | 0.0505 | 0.0487 | 0.0520 | 0.0480 |
| TRUE | 0.050 | 0.0500 | 0.0500 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 101.04% | 102.92% | 100.90% | 97.30% | 103.92% | 96.08% |
| PREP BLANK | | | | | | |
| | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |

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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | TIN (mg/L) | VANADIUM (mg/L) | ZINC (mg/L) |
|-----------------|---------------|--------------------|----------------|
| METHOD | EPA 6010 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 03/25,29/11 | 03/25,28/11 | 03/25,28/11 |
| DETECTION LIMIT | 0.010 | 0.0010 | 0.0010 |
| REPORTING LIMIT | 0.040 | 0.0030 | 0.0050 |
| DUPLICATE | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | <0.040 | 0.0032 | <0.0050 |
| DUPLICATE | <0.040 | 0.0032 | <0.0050 |
| RPD | NC | 0.57% | NC |
| SPIKE SAMPLE | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | <0.040 | 0.0032 | <0.0050 |
| SPIKED SAMPLE | 1.06 | 0.0510 | 0.0580 |
| SPIKE ADDED | 1.00 | 0.0500 | 0.0500 |
| % RECOVERY | 105.73% | 95.56% | 115.92% |
| QC CHECK | | | |
| FOUND | 1.08 | 0.0496 | 0.0474 |
| TRUE | 1.00 | 0.0500 | 0.0500 |
| % RECOVERY | 108.20% | 99.14% | 94.86% |
| PREP BLANK | | | |
| | <0.040 | <0.0030 | <0.0050 |

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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | CALCIUM (mg/L) | POTASSIUM (mg/L) | SODIUM (mg/L) | MAGNESIUM (mg/L) | IRON (mg/L) | MANGANESE (mg/L) |
|-----------------|-------------------|---------------------|------------------|---------------------|----------------|---------------------|
| METHOD | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 |
| DATE ANALYZED | 03/25,29/11 | 03/25,29/11 | 03/25,29/11 | 03/25,29/11 | 03/25,29/11 | 03/25,29/11 |
| DETECTION LIMIT | 0.050 | 0.300 | 0.300 | 0.050 | 0.010 | 0.003 |
| REPORTING LIMIT | 0.100 | 0.500 | 0.500 | 0.100 | 0.020 | 0.010 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | 8.54 | 0.942 | 4.54 | 7.64 | <0.020 | <0.010 |
| DUPLICATE | 8.53 | 0.948 | 4.51 | 7.62 | <0.020 | <0.010 |
| RPD | 0.14% | 0.58% | 0.56% | 0.24% | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | 8.54 | 0.942 | 4.54 | 7.64 | <0.020 | <0.010 |
| SPIKED SAMPLE | 16.6 | 12.7 | 14.0 | 16.6 | 9.87 | 1.02 |
| SPIKE ADDED | 10.0 | 10.0 | 10.0 | 10.0 | 10.0 | 1.00 |
| % RECOVERY | 80.91% | 117.46% | 94.18% | 89.60% | 98.70% | 102.00% |
| QC CHECK | | | | | | |
| FOUND | 0.977 | 9.81 | 10.3 | 10.0 | 0.981 | 1.02 |
| TRUE | 1.00 | 10.0 | 10.0 | 10.0 | 1.00 | 1.00 |
| % RECOVERY | 97.66% | 98.08% | 103.37% | 100.04% | 98.11% | 101.65% |
| PREP BLANK | <0.100 | <0.500 | <0.500 | <0.100 | <0.020 | <0.010 |

RPD = RELATIVE PERCENT DIFFERENCE.
 NA = NOT APPLICABLE OR NOT AVAILABLE.
 NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.
 OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.



AQUATIC RESEARCH INCORPORATED

LABORATORY & CONSULTING SERVICES
3927 AURORA AVENUE NORTH, SEATTLE, WA 98103
PHONE: (206) 632-2715 FAX: (206) 632-2417

| | | |
|---|--------------|----------------------------|
| CASE FILE NUMBER: | BKS001-47,48 | PAGE 12 |
| REPORT DATE: | 04/06/11 | |
| DATE SAMPLED: | 03/23,24/11 | DATE RECEIVED: 03/24,25/11 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - TOTAL METALS

| QC PARAMETER | MERCURY (mg/l) | NICKEL (mg/l) |
|-----------------|-------------------|------------------|
| METHOD | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 03/25,28/11 | 03/25,28/11 |
| DETECTION LIMIT | 0.0001 | 0.0010 |
| REPORTING LIMIT | 0.0002 | 0.0050 |
| DUPLICATE | | |
| SAMPLE ID | OL-MW-1 | OL-MW-1 |
| ORIGINAL | <0.0002 | <0.0050 |
| DUPLICATE | <0.0002 | <0.0050 |
| RPD | NC | NC |
| SPIKE SAMPLE | | |
| SAMPLE ID | OL-MW-1 | OL-MW-1 |
| ORIGINAL | <0.0002 | <0.0050 |
| SPIKED SAMPLE | 0.0022 | 0.0517 |
| SPIKE ADDED | 0.0020 | 0.0500 |
| % RECOVERY | 110.00% | 103.46% |
| QC CHECK | | |
| FOUND | 0.0021 | 0.0503 |
| TRUE | 0.0020 | 0.0500 |
| % RECOVERY | 104.00% | 100.62% |
| PREP BLANK | <0.0002 | <0.0050 |

RPD = RELATIVE PERCENT DIFFERENCE.

NA = NOT APPLICABLE OR NOT AVAILABLE.

NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.

OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.

Submitted By:

Steven Lazoff
Laboratory Director



Aquatic Research, Inc.

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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-1 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 85% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



Aquatic Research, Inc.

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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-2 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 109% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-4 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 71% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-5A | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 70% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A7 | Matrix: | Water |
| Sample ID No.: | OL-ER | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 93% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs LABORATORY FORTIFIED BLANK REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|---------------|-------------------------------|-------|
| Case File Number: | LFB (3/25/11) | Matrix: | Water |
| Sample ID No.: | LFB | Sample Vol. (ml) | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Arochlor | RECOVERY | LCL | UCL |
|------------|----------|-----|------|
| Total PCBs | 75% | 70% | 130% |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 82% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs LABORATORY REAGENT BLANK REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|--------------|-------------------------------|-------|
| Case File Number: | MB (3/25/11) | Matrix: | Water |
| Sample ID No.: | LRB | Sample Vol. (ml): | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Aroclor | Result (ug/L) | UCL (ug/L) | RL | (ug/L) |
|------------|------------------|---------------|----|--------|
| Total PCBs | < 0.1 | < 0.1 | | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 73% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit

Vinyl Chloride by SIM

T. Meadows

4/4/11

| Sample ID | Client ID | Sample Amount (ml) | Dilution Factor | Vinyl Chloride (µg/L) | 1,2-Dichloroethane-d4 (µg/L) | % R Surr | % R VC |
|----------------|--------------|--------------------|-----------------|-----------------------|------------------------------|----------|--------|
| CCV | CCV | 25 | 1 | 0.10 | 0.11 | 110.0% | 100.0% |
| CCV | CCV | 25 | 1 | 0.09 | 0.08 | 80.0% | 90.0% |
| 4/4/11-MB | Method Blank | 25 | 1 | <0.02 | 0.11 | 110.0% | |
| BKS00147A1 | OL-MW-1 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00147A2 | OL-MW-5A | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00147A3 | OL-MW-2 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00147A4 | OL-MW-4 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00147A7 | OL-ER | 25 | 1 | 0.05 | 0.07 | 70.0% | |
| BKS00147A8 | OL-TB | 25 | 1 | 0.09 | 0.08 | 80.0% | |
| BKS00148A1 | OL-MW-7 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A2 | OL-MW-8 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A3 | OL-MW-6 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A3 Dup | OL-MW-6 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A4 | OL-MW-10 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A5 | OL-MW-3 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A6 | OL-TB | 25 | 1 | 0.05 | 0.07 | 70.0% | |
| BKS00148A7 | OL-MW-11 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| | | | RPD | 0% | | 0% | |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A5** Matrix: Water
Sample ID No.: **OL-MW-3** Sample Wt./Vol. (gm/ml) **25.0**
Date Collected: **03/24/11** Dilution Factor: **1**
Date Received: **03/25/11**
Date Analyzed: **04/ 1/11** Analyst: **T. Meadows**
Date of Report: **04/01/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110401\ 1201012.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-3 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 1201012.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 84% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 110% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 84% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0801008.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 78% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 107% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 82% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-7 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-7 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0601006.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 82% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 107% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 88% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-8 | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-8 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0701007.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 81% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 109% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 79% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A4** Matrix: Water
Sample ID No.: **OL-MW-10** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **03/24/11** Dilution Factor: **1**
Date Received: **03/25/11**
Date Analyzed: **04/ 1/11** Analyst: **T. Meadows**
Date of Report: **04/01/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110401\ 0901009.D**

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0901009.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 73% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 101% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 80% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A7 | Matrix: | Water |
| Sample ID No.: | OL-MW-11 | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A7 | Matrix: | Water |
| Sample ID No.: | OL-MW-11 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 1401014.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 81% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 107% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 83% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00148A4 MS / MSD | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 4/1/11 | | |
| Data File Paths: | C:\HPCHEM\1\DATA\VOA\110401\ 1001010.D | | |
| | C:\HPCHEM\1\DATA\VOA\110401\ 1101011.D | | |

| Surrogate Recoveries | % Rec. | | |
|----------------------|--------|-----|-----|
| | MS | MSD | RPD |
| Dibromofluoromethane | 72% | 70% | 3% |
| Toluene-d8 | 106% | 97% | 10% |
| 4-Bromofluorobenzene | 84% | 71% | 15% |

| Spike Recoveries | % Rec. | | |
|--------------------|--------|------|-----|
| | MS | MSD | RPD |
| 1,1-Dichloroethene | 116% | 114% | 2% |
| Benzene | 94% | 97% | 3% |
| Trichloroethene | 119% | 108% | 10% |
| Toluene | 106% | 105% | 1% |
| Chlorobenzene | 106% | 112% | 5% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|------------------------------|-------------------------|------------|
| Case File Number: | 4/1/11-LCS | Matrix: | Water |
| Sample ID No.: | 4/1/11-LCS | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed | 04/ 1/11 | Analyst: | T. Meadows |
| Date of Report: | 4/1/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 94% | 66% | 118% |
| Toluene-d8 | 97% | 51% | 143% |
| 4-Bromofluorobenzene | 83% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 91% | 75% | 125% |
| Benzene | 96% | 75% | 125% |
| Trichloroethene | 100% | 75% | 125% |
| Toluene | 99% | 75% | 125% |
| Chlorobenzene | 100% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00148A4 MS | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | | |
| Date Analyzed: | 04/ 1/11 | Analyst: | T. Meadows |
| Date of Report: | 4/1/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ 1001010.D | | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 72% | 66% | 118% |
| Toluene-d8 | 106% | 51% | 143% |
| 4-Bromofluorobenzene | 84% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 116% | 75% | 125% |
| Benzene | 94% | 75% | 125% |
| Trichloroethene | 119% | 75% | 125% |
| Toluene | 106% | 75% | 125% |
| Chlorobenzene | 106% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00148A4 MSD | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | | |
| Date Analyzed: | 04/ 1/11 | Analyst: | T. Meadows |
| Date of Report: | 4/1/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\1101011.D | | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 70% | 66% | 118% |
| Toluene-d8 | 97% | 51% | 143% |
| 4-Bromofluorobenzene | 71% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 114% | 75% | 125% |
| Benzene | 97% | 75% | 125% |
| Trichloroethene | 108% | 75% | 125% |
| Toluene | 105% | 75% | 125% |
| Chlorobenzene | 112% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | 4/1/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | 4/1/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0401004.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 108% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 105% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 91% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A6** Matrix: Water
Sample ID No.: **OL-TB** Sample Wt./Vol. (gm/ml) **25.0**
Date Collected: **03/24/11** Dilution Factor: **1**
Date Received: **03/25/11**
Date Analyzed: **04/ 1/11** Analyst: **T. Meadows**
Date of Report: **04/01/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110401\ 1301013.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A6 | Matrix: | Water |
| Sample ID No.: | OL-TB | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 1301013.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 74% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 108% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 82% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A5**
Sample ID No.: **OL-MW-3**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1060**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3701019.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-3 | Sample Vol. (ml) | 1060 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purges Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 39% | 24-132% | 65-135% |
| Phenol-d5 | 39% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 99% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 82% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 28% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 110% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A3**
Sample ID No.: **OL-MW-6**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3301015.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6 | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | 3301015.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 61% | 24-132% | 65-135% |
| Phenol-d5 | 57% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 119% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 113% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 32% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 127% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A1**
Sample ID No.: **OL-MW-7**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/07/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1060**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials:
3101013.D

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-7 | Sample Vol. (ml) | 1060 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/07/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 77% | 24-132% | 65-135% |
| Phenol-d5 | 62% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 129% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 95% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 42% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 118% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A2**
Sample ID No.: **OL-MW-8**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3201014.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-8 | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | 3201014.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 41% | 24-132% | 65-135% |
| Phenol-d5 | 36% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 133% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 101% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 36% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 114% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A4**
Sample ID No.: **OL-MW-10**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1060**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3401016.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Vol. (ml) | 1060 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | 3401016.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 10% | 24-132% | 65-135% |
| Phenol-d5 | 22% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 127% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 114% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 19% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 125% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-1 | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A7**
Sample ID No.: **OL-MW-11**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3801020.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A7 | Matrix: | Water |
| Sample ID No.: | OL-MW-11 | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 44% | 24-132% | 65-135% |
| Phenol-d5 | 37% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 125% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 102% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 47% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 118% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
10/14/10chy/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | 3/28/11-LCS | Matrix: | Water |
| Sample ID No.: | Lab Control Spike | Sample Vol. (ml) | 1000 |
| Date Collected: | n/a | Final Volume (ml) | 2.0 |
| Date Received: | n/a | Dilution Factor: | 1 |
| Date Analyzed: | 04/07/11 | Analyst: | T. Meadows |
| Date of Report: | 04/11/11 | Supervisor's Initials: | |
| Data File Path: | Q:\5973\PAM\040711\ | 3001012.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 87% | 24-132% | 65-135% |
| Phenol-d5 | 59% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 116% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 106% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 26% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 119% | 80-141% | 65-135% |

| Matix Spike Recoveries | %Rec. | QC limits | |
|--------------------------|-------|-----------|---------|
| | | Water | Soil |
| Phenol | 52% | 50-150% | 50-150% |
| 2-Chlorophenol | 58% | 50-150% | 50-150% |
| 1,4-Dichlorobenzene | 76% | 50-150% | 50-150% |
| N-Nitroso-n-propyl amine | 84% | 50-150% | 50-150% |
| 1,2,4-Trichlorobenzene | 68% | 50-150% | 50-150% |
| 4-Chloro-3-methyl phenol | 48% | 50-150% | 50-150% |
| Acenaphthene | 96% | 50-150% | 50-150% |
| 4-Nitrophenol | 44% | 50-150% | 50-150% |
| 2,4-Dinitrotoluene | 48% | 50-150% | 50-150% |
| Pentachlorophenol | 0% | 50-150% | 50-150% |
| Pyrene | 90% | 50-150% | 50-150% |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
10/14/10ghy/Mass Spectrometry

| | | | |
|-------------------|------------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A4 MS | Matrix: | Water |
| Sample ID No.: | OL-MW-10 Matrix Spike | Sample Vol. (ml) | 1070 |
| Date Collected: | 3/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 3/25/11 | Dilution Factor: | 1 |
| Date Analyzed: | 04/08/11 | Analyst: | T. Meadows |
| Date of Report: | 04/11/11 | Supervisor's Initials: | |
| Data File Path: | Q:\5973\PAM\040711\ | 3501017.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 27% | 24-132% | 65-135% |
| Phenol-d5 | 29% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 108% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 95% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 12% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 114% | 80-141% | 65-135% |

| Matix Spike Recoveries | %Rec. | QC limits | |
|--------------------------|-------|-----------|---------|
| | | Water | Soil |
| Phenol | 35% | 50-150% | 50-150% |
| 2-Chlorophenol | 37% | 50-150% | 50-150% |
| 1,4-Dichlorobenzene | 76% | 50-150% | 50-150% |
| N-Nitroso-n-propyl amine | 100% | 50-150% | 50-150% |
| 1,2,4-Trichlorobenzene | 63% | 50-150% | 50-150% |
| 4-Chloro-3-methyl phenol | 33% | 50-150% | 50-150% |
| Acenaphthene | 104% | 50-150% | 50-150% |
| 4-Nitrophenol | 51% | 50-150% | 50-150% |
| 2,4-Dinitrotoluene | 62% | 50-150% | 50-150% |
| Pentachlorophenol | 4% | 50-150% | 50-150% |
| Pyrene | 91% | 50-150% | 50-150% |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
10/14/10chy/Mass Spectrometry

| | | | |
|-------------------|--|------------------------|-------------------|
| Case File Number: | BKS00148A4 MSD | Matrix: | Water |
| Sample ID No.: | OL-MW-10 Matrix Spike Duplicate | Sample Vol. (ml) | 1070 |
| Date Collected: | 3/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 3/25/11 | Dilution Factor: | 1 |
| Date Analyzed: | 04/08/11 | Analyst: | T. Meadows |
| Date of Report: | 04/11/11 | Supervisor's Initials: | |
| Data File Path: | Q:\5973\PAM\040711\ | 3601018.D | |

| Surrogate Recoveries | %Rec. | QC limits | | RPD |
|----------------------|-------|-----------|---------|-------|
| | | Water | Soil | |
| 2-Fluorophenol | 25% | 24-132% | 65-135% | 5.0% |
| Phenol-d5 | 30% | 0-105% | 65-135% | 5.0% |
| Nitrobenzene-d5 | 125% | 24-136% | 65-135% | 14.2% |
| 2-Fluorobiphenyl | 109% | 41-134% | 65-135% | 13.7% |
| 2,4,6-Tribromophenol | 10% | 0-183% | 65-135% | 19.8% |
| p-Terphenyl-d14 | 115% | 80-141% | 65-135% | 0.8% |

| Matix Spike Recoveries | %Rec. | QC limits | | RPD |
|--------------------------|-------|-----------|---------|-------|
| | | Water | Soil | |
| Phenol | 40% | 50-150% | 50-150% | 12.3% |
| 2-Chlorophenol | 40% | 50-150% | 50-150% | 7.0% |
| 1,4-Dichlorobenzene | 90% | 50-150% | 50-150% | 17.0% |
| N-Nitroso-n-propyl amine | 104% | 50-150% | 50-150% | 4.5% |
| 1,2,4-Trichlorobenzene | 75% | 50-150% | 50-150% | 17.2% |
| 4-Chloro-3-methyl phenol | 35% | 50-150% | 50-150% | 6.9% |
| Acenaphthene | 109% | 50-150% | 50-150% | 4.3% |
| 4-Nitrophenol | 50% | 50-150% | 50-150% | 1.3% |
| 2,4-Dinitrotoluene | 64% | 50-150% | 50-150% | 3.0% |
| Pentachlorophenol | 4% | 50-150% | 50-150% | 8.4% |
| Pyrene | 90% | 50-150% | 50-150% | 0.4% |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **3/28/11-MB**
Sample ID No.: **Method Blank**
Date Collected: **n/a**
Date Received: **n/a**
Date Extracted: **03/28/11**
Date Analyzed: **04/07/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1000**
Final Volume (ml) **2.0**
Dilution Factor: **1**
Analyst: **T. Meadows**
Supervisor's Initials: **2901011.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.80 | U | 86-73-7 | Fluorene | 0.80 | U |
| 76-01-7 | Pentachloroethane | 0.80 | U | 84-66-2 | Diethyl phthalate | 0.80 | U |
| 62-53-3 | Aniline | 0.80 | U | 100-01-6 | 4-Nitroaniline | 0.80 | U |
| 108-95-2 | Phenol | 0.80 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.80 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.80 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.80 | U |
| 95-57-8 | 2-Chlorophenol | 0.80 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.80 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 103-33-3 | Azobenzene | 0.80 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.80 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.80 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.80 | U | 319-84-6 | a-BHC | 0.80 | U |
| 100-51-6 | Benzyl Alcohol | 0.80 | U | 118-74-1 | Hexachlorobenzene | 0.80 | U |
| 95-48-7 | 2-Methyl phenol | 0.80 | U | 319-85-7 | b-BHC | 0.80 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.80 | U | 87-86-5 | Pentachlorophenol | 0.80 | U |
| 98-86-2 | Acetophenone | 0.80 | U | 58-89-9 | g-BHC (Lindane) | 0.80 | U |
| 67-72-1 | Hexachloroethane | 0.80 | U | 85-01-8 | Phenanthrene | 0.80 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.80 | U | 120-12-7 | Anthracene | 0.80 | U |
| 106-44-5 | 4-Methyl phenol | 0.80 | U | 319-86-8 | d-BHC | 0.80 | U |
| 98-95-3 | Nitrobenzene | 0.80 | U | 84-74-2 | Di-n-butyl phthalate | 0.80 | U |
| 100-75-4 | N-nitrosopiperidine | 0.80 | U | 76-44-8 | Heptachlor | 0.80 | U |
| 78-59-1 | Isophorone | 0.80 | U | 309-00-2 | Aldrin | 0.80 | U |
| 88-75-5 | 2-Nitrophenol | 0.80 | U | 1024-57-3 | Heptachlor epoxide | 0.80 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.80 | U | 206-44-0 | Fluoranthrene | 0.80 | U |
| 65-85-0 | Benzoic Acid | 0.80 | U | 129-00-0 | Pyrene | 0.80 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.80 | U | 959-98-8 | Endosulfan I | 0.80 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.80 | U | 92-87-5 | Benzidine | 0.80 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.80 | U | 72-55-9 | 4,4'-DDE | 0.80 | U |
| 91-20-3 | Naphthalene | 0.80 | U | 60-57-1 | Dieldrin | 0.80 | U |
| 106-47-8 | 4-Chloroaniline | 0.80 | U | 72-20-8 | Endrin | 0.80 | U |
| 87-68-3 | Hexachlorobutadiene | 0.80 | U | 33213-65-9 | Endosulfan II | 0.80 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.80 | U | 72-54-8 | 4,4'-DDD | 0.80 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.80 | U | 7421-92-4 | Endrin aldehyde | 0.80 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.80 | U | 85-68-7 | Butyl benzenyl phthalate | 0.80 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.80 | U | 1031-07-8 | Endosulfan sulfate | 0.80 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.80 | U | 50-29-3 | 4,4'-DDT | 0.80 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.80 | U | 56-55-3 | Benzo(a)anthracene | 0.80 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.80 | U | 218-01-9 | Chrysene | 0.80 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.80 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.80 | U |
| 88-74-4 | 2-Nitroaniline | 0.80 | U | 72-43-5 | Methoxychlor | 0.80 | U |
| 131-11-3 | Dimethyl phthalate | 0.80 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.80 | U |
| 208-96-8 | Acenaphthylene | 0.80 | U | 117-84-0 | Di-n-octyl phthalate | 0.80 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.80 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.80 | U |
| 99-09-2 | 3-Nitroaniline | 0.80 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.80 | U |
| 83-32-9 | Acenaphthene | 0.80 | U | 50-32-8 | Benzo(a)pyrene | 0.80 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.80 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.80 | U |
| 132-64-9 | Dibenzofuran | 0.80 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.80 | U |
| 100-02-7 | 4-Nitrophenol | 0.80 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.80 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.80 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.80 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | 3/28/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Vol. (ml) | 1000 |
| Date Collected: | n/a | Final Volume (ml) | 2.0 |
| Date Received: | n/a | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/07/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 79% | 24-132% | 65-135% |
| Phenol-d5 | 78% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 128% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 102% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 36% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 110% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.80 | U |
| 100-52-7 | Benzaldehyde | 0.80 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.80 | U |
| 95-53-4 | o-Toluidine | 0.80 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.80 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.80 | U |
| 1888-71-7 | Hexachloropropene | 0.80 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.80 | U |
| 87-65-0 | Caprolactam | 0.80 | U |
| 106-50-3 | p-Phenylenediamine | 0.80 | U |
| 120-58-1 | Isosafrole | 0.80 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.80 | U |
| 94-59-7 | Safrole | 0.80 | U |
| 92-52-4 | Biphenyl | 0.80 | U |
| 101-84-8 | Diphenyl ether | 0.80 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.80 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.80 | U |
| 608-93-5 | Pentachlorobenzene | 0.80 | U |
| 134-32-7 | 1-Naphthalenamine | 0.80 | U |
| 91-59-8 | 2-Naphthalenamine | 0.80 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.80 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.80 | U |
| 122-39-4 | Diphenylamine | 0.80 | U |
| 3689-24-5 | Sulfotep | 0.80 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.80 | U |
| 298-02-2 | Phorate | 0.80 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.80 | U |
| 62-44-2 | Phenacetin | 0.80 | U |
| 60-51-5 | Dimethoate | 0.80 | U |
| 1912-24-9 | Atrazine | 0.80 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.80 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.80 | U |
| 23950-58-! | Pronamide | 0.80 | U |
| 298-04-4 | Disulfoton | 0.80 | U |
| 298-00-0 | Methyl parathion | 0.80 | U |
| 56-38-2 | Parathion | 0.80 | U |
| 91-80-5 | Methapyrilene | 0.80 | U |
| 465-73-6 | Isodrin | 0.80 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.80 | U |
| 510-15-0 | Chlorobenzilate | 0.80 | U |
| 143-50-0 | Kepone | 0.80 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.80 | U |
| 52-85-7 | Famphur | 0.80 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.80 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.80 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.80 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.80 | U |



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-3 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/8/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 107% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 74% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-7 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 103% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-8 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 119% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 115% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A7 | Matrix: | Water |
| Sample ID No.: | OL-MW-11 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/8/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 95% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs MATRIX SPIKE REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|---------------|-------------------------------|-------|
| Case File Number: | BKS00148A4 MS | Matrix: | Water |
| Sample ID No.: | OL-MW-10 MS | Sample Vol. (ml) | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Aroclor | RECOVERY | LCL | UCL |
|------------|----------|-----|------|
| Total PCBs | 97% | 70% | 130% |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 99% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



Aquatic Research, Inc.

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TOTAL PCBs MATRIX SPIKE REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|----------------|-------------------------------|-------|
| Case File Number: | BKS00148A4 MSD | Matrix: | Water |
| Sample ID No.: | OL-MW-10 MSD | Sample Vol. (ml) | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/8/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Aroclor | RECOVERY | LCL | UCL |
|----------------|-----------------|------------|------------|
| Total PCBs | 104% | 70% | 130% |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 99% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



Aquatic Research, Inc.

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TOTAL PCBs LABORATORY FORTIFIED BLANK REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|---------------|-------------------------------|-------|
| Case File Number: | LFB (3/28/11) | Matrix: | Water |
| Sample ID No.: | LFB | Sample Vol. (ml) | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Arochlor | RECOVERY | LCL | UCL |
|------------|----------|-----|------|
| Total PCBs | 121% | 70% | 130% |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 95% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



Aquatic Research, Inc.

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TOTAL PCBs LABORATORY REAGENT BLANK REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|--------------|-------------------------------|-------|
| Case File Number: | MB (3/28/11) | Matrix: | Water |
| Sample ID No.: | LRB | Sample Vol. (ml): | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Aroclor | Result (ug/L) | UCL (ug/L) | RL | (ug/L) |
|------------|------------------|---------------|----|--------|
| Total PCBs | < 0.1 | < 0.1 | | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 95% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A1** Matrix: Water
Sample ID No.: **OL-MW-1** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\0601006.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-1 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 0601006.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 79% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 105% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 92% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A3** Matrix: Water
Sample ID No.: **OL-MW-2** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 1001010.D**

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-2 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 1001010.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 77% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 108% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 86% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A4** Matrix: Water
Sample ID No.: **OL-MW-4** Sample Wt./Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 1101011.D**

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



Aquatic Research Inc.
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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-4 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 1101011.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 72% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 104% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 84% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



Aquatic Research Inc.
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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A2** Matrix: Water
Sample ID No.: **OL-MW-5A** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 0801008.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-5A | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 0801008.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 80% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 106% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 92% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number:
Sample ID No.:
Date Collected:
Date Received:
Date Analyzed:
Date of Report:
Data File Path:

BKS00147A2 Dup
OL-MW-5A Duplicate
03/23/11
03/24/11
03/31/11
03/31/11
C:\HPCHEM\1\DATA\VOA\110331

Matrix:
Sample Wt/Vol. (gm/ml)
Dilution Factor:

Water
25.0
1

Analyst:
Supervisor's Initials:
0901009.D

T. Meadows

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A2 Dup | Matrix: | Water |
| Sample ID No.: | OL-MW-5A Duplicate | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 0901009.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 70% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 106% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 84% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A7** Matrix: Water
Sample ID No.: **OL-ER** Sample Wt./Vol. (gm/ml) **25.0**
Date Collected: **03/23/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 1201012.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A7 | Matrix: | Water |
| Sample ID No.: | OL-ER | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/23/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 1201012.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 72% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 109% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 88% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A1 MS | Matrix: | Water |
| Sample ID No.: | BKS00147A1 MS | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | | |
| Date Analyzed: | 03/31/11 | Analyst: | T. Meadows |
| Date of Report: | 3/31/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | 0701007.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 79% | 66% | 118% |
| Toluene-d8 | 110% | 51% | 143% |
| 4-Bromofluorobenzene | 86% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 102% | 75% | 125% |
| Benzene | 105% | 75% | 125% |
| Trichloroethene | 112% | 75% | 125% |
| Toluene | 112% | 75% | 125% |
| Chlorobenzene | 113% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | 3/31/11-LCS | Matrix: | Water |
| Sample ID No.: | 3/31/11-LCS | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed | 03/31/11 | Analyst: | T. Meadows |
| Date of Report: | 3/31/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 108% | 66% | 118% |
| Toluene-d8 | 107% | 51% | 143% |
| 4-Bromofluorobenzene | 107% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 103% | 75% | 125% |
| Benzene | 102% | 75% | 125% |
| Trichloroethene | 102% | 75% | 125% |
| Toluene | 102% | 75% | 125% |
| Chlorobenzene | 107% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | 3/31/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | 3/31/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 0401004.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 106% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 104% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 106% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A8** Matrix: Water
Sample ID No.: **OL-TB** Sample Wt./Vol. (gm/ml) **25.0**
Date Collected: **03/15/11** Dilution Factor: **1**
Date Received: **03/24/11**
Date Analyzed: **03/31/11** Analyst: **T. Meadows**
Date of Report: **03/31/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110331\ 1301013.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A8 | Matrix: | Water |
| Sample ID No.: | OL-TB | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/15/11 | Dilution Factor: | 1 |
| Date Received: | 03/24/11 | Analyst: | T. Meadows |
| Date Analyzed: | 03/31/11 | Supervisor's Initials: | |
| Date of Report: | 03/31/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110331\ | | 1301013.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 74% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 106% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 83% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



Aquatic Research Inc.
3927 Aurora Ave. N., Seattle, WA 98103 | (206) 632-2715

SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A1**
Sample ID No.: **OL-MW-1**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1050**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **0601006.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.76 | U | 86-73-7 | Fluorene | 0.76 | U |
| 76-01-7 | Pentachloroethane | 0.76 | U | 84-66-2 | Diethyl phthalate | 0.76 | U |
| 62-53-3 | Aniline | 0.76 | U | 100-01-6 | 4-Nitroaniline | 0.76 | U |
| 108-95-2 | Phenol | 0.76 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.76 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.76 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.76 | U |
| 95-57-8 | 2-Chlorophenol | 0.76 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.76 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.76 | U | 103-33-3 | Azobenzene | 0.76 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.76 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.76 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.76 | U | 319-84-6 | a-BHC | 0.76 | U |
| 100-51-6 | Benzyl Alcohol | 0.76 | U | 118-74-1 | Hexachlorobenzene | 0.76 | U |
| 95-48-7 | 2-Methyl phenol | 0.76 | U | 319-85-7 | b-BHC | 0.76 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.76 | U | 87-86-5 | Pentachlorophenol | 0.76 | U |
| 98-86-2 | Acetophenone | 0.76 | U | 58-89-9 | g-BHC (Lindane) | 0.76 | U |
| 67-72-1 | Hexachloroethane | 0.76 | U | 85-01-8 | Phenanthrene | 0.76 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.76 | U | 120-12-7 | Anthracene | 0.76 | U |
| 106-44-5 | 4-Methyl phenol | 0.76 | U | 319-86-8 | d-BHC | 0.76 | U |
| 98-95-3 | Nitrobenzene | 0.76 | U | 84-74-2 | Di-n-butyl phthalate | 0.76 | U |
| 100-75-4 | N-nitrosopiperidine | 0.76 | U | 76-44-8 | Heptachlor | 0.76 | U |
| 78-59-1 | Isophorone | 0.76 | U | 309-00-2 | Aldrin | 0.76 | U |
| 88-75-5 | 2-Nitrophenol | 0.76 | U | 1024-57-3 | Heptachlor epoxide | 0.76 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.76 | U | 206-44-0 | Fluoranthrene | 0.76 | U |
| 65-85-0 | Benzoic Acid | 0.76 | U | 129-00-0 | Pyrene | 0.76 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.76 | U | 959-98-8 | Endosulfan I | 0.76 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.76 | U | 92-87-5 | Benzidine | 0.76 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.76 | U | 72-55-9 | 4,4'-DDE | 0.76 | U |
| 91-20-3 | Naphthalene | 0.76 | U | 60-57-1 | Dieldrin | 0.76 | U |
| 106-47-8 | 4-Chloroaniline | 0.76 | U | 72-20-8 | Endrin | 0.76 | U |
| 87-68-3 | Hexachlorobutadiene | 0.76 | U | 33213-65-9 | Endosulfan II | 0.76 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.76 | U | 72-54-8 | 4,4'-DDD | 0.76 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.76 | U | 7421-92-4 | Endrin aldehyde | 0.76 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.76 | U | 85-68-7 | Butyl benzenyl phthalate | 0.76 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.76 | U | 1031-07-8 | Endosulfan sulfate | 0.76 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.76 | U | 50-29-3 | 4,4'-DDT | 0.76 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.76 | U | 56-55-3 | Benzo(a)anthracene | 0.76 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.76 | U | 218-01-9 | Chrysene | 0.76 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.76 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.76 | U |
| 88-74-4 | 2-Nitroaniline | 0.76 | U | 72-43-5 | Methoxychlor | 0.76 | U |
| 131-11-3 | Dimethyl phthalate | 0.76 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.76 | U |
| 208-96-8 | Acenaphthylene | 0.76 | U | 117-84-0 | Di-n-octyl phthalate | 0.76 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.76 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.76 | U |
| 99-09-2 | 3-Nitroaniline | 0.76 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.76 | U |
| 83-32-9 | Acenaphthene | 0.76 | U | 50-32-8 | Benzo(a)pyrene | 0.76 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.76 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.76 | U |
| 132-64-9 | Dibenzofuran | 0.76 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.76 | U |
| 100-02-7 | 4-Nitrophenol | 0.76 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.76 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.76 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.76 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-1 | Sample Vol. (ml) | 1050 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 50% | 24-132% | 65-135% |
| Phenol-d5 | 35% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 99% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 81% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 43% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 125% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.76 | U |
| 100-52-7 | Benzaldehyde | 0.76 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.76 | U |
| 95-53-4 | o-Toluidine | 0.76 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.76 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.76 | U |
| 1888-71-7 | Hexachloropropene | 0.76 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.76 | U |
| 87-65-0 | Caprolactam | 0.76 | U |
| 106-50-3 | p-Phenylenediamine | 0.76 | U |
| 120-58-1 | Isosafrole | 0.76 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.76 | U |
| 94-59-7 | Safrole | 0.76 | U |
| 92-52-4 | Biphenyl | 0.76 | U |
| 101-84-8 | Diphenyl ether | 0.76 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.76 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.76 | U |
| 608-93-5 | Pentachlorobenzene | 0.76 | U |
| 134-32-7 | 1-Naphthalenamine | 0.76 | U |
| 91-59-8 | 2-Naphthalenamine | 0.76 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.76 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.76 | U |
| 122-39-4 | Diphenylamine | 0.76 | U |
| 3689-24-5 | Sulfotep | 0.76 | U |

| | | | |
|------------|----------------------------------|------|---|
| 2303-16-4 | Diallate | 0.76 | U |
| 298-02-2 | Phorate | 0.76 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.76 | U |
| 62-44-2 | Phenacetin | 0.76 | U |
| 60-51-5 | Dimethoate | 0.76 | U |
| 1912-24-9 | Atrazine | 0.76 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.76 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.76 | U |
| 23950-58-1 | Pronamide | 0.76 | U |
| 298-04-4 | Disulfoton | 0.76 | U |
| 298-00-0 | Methyl parathion | 0.76 | U |
| 56-38-2 | Parathion | 0.76 | U |
| 91-80-5 | Methapyrilene | 0.76 | U |
| 465-73-6 | Isodrin | 0.76 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.76 | U |
| 510-15-0 | Chlorobenzilate | 0.76 | U |
| 143-50-0 | Kepone | 0.76 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.76 | U |
| 52-85-7 | Famphur | 0.76 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.76 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracene | 0.76 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.76 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.76 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A3**
Sample ID No.: **OL-MW-2**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1060**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **0801008.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-2 | Sample Vol. (ml) | 1060 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 73% | 24-132% | 65-135% |
| Phenol-d5 | 51% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 108% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 82% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 46% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 137% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A4**
Sample ID No.: **OL-MW-4**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **0901009.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-4 | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 50% | 24-132% | 65-135% |
| Phenol-d5 | 43% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 110% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 96% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 38% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 117% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A2**
Sample ID No.: **OL-MW-5A**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials:
0701007.D

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-5A | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | 0701007.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 51% | 24-132% | 65-135% |
| Phenol-d5 | 43% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 98% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 78% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 37% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 102% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00147A7**
Sample ID No.: **OL-ER**
Date Collected: **03/23/11**
Date Received: **03/24/11**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1040**
Final Volume (ml) **2.0**
Dilution Factor: **1**
Analyst: **T. Meadows**
Supervisor's Initials: **1001010.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.77 | U | 86-73-7 | Fluorene | 0.77 | U |
| 76-01-7 | Pentachloroethane | 0.77 | U | 84-66-2 | Diethyl phthalate | 0.77 | U |
| 62-53-3 | Aniline | 0.77 | U | 100-01-6 | 4-Nitroaniline | 0.77 | U |
| 108-95-2 | Phenol | 0.77 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.77 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.77 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.77 | U |
| 95-57-8 | 2-Chlorophenol | 0.77 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.77 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.77 | U | 103-33-3 | Azobenzene | 0.77 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.77 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.77 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.77 | U | 319-84-6 | a-BHC | 0.77 | U |
| 100-51-6 | Benzyl Alcohol | 0.77 | U | 118-74-1 | Hexachlorobenzene | 0.77 | U |
| 95-48-7 | 2-Methyl phenol | 0.77 | U | 319-85-7 | b-BHC | 0.77 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.77 | U | 87-86-5 | Pentachlorophenol | 0.77 | U |
| 98-86-2 | Acetophenone | 0.77 | U | 58-89-9 | g-BHC (Lindane) | 0.77 | U |
| 67-72-1 | Hexachloroethane | 0.77 | U | 85-01-8 | Phenanthrene | 0.77 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.77 | U | 120-12-7 | Anthracene | 0.77 | U |
| 106-44-5 | 4-Methyl phenol | 0.77 | U | 319-86-8 | d-BHC | 0.77 | U |
| 98-95-3 | Nitrobenzene | 0.77 | U | 84-74-2 | Di-n-butyl phthalate | 0.77 | U |
| 100-75-4 | N-nitrosopiperidine | 0.77 | U | 76-44-8 | Heptachlor | 0.77 | U |
| 78-59-1 | Isophorone | 0.77 | U | 309-00-2 | Aldrin | 0.77 | U |
| 88-75-5 | 2-Nitrophenol | 0.77 | U | 1024-57-3 | Heptachlor epoxide | 0.77 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.77 | U | 206-44-0 | Fluoranthrene | 0.77 | U |
| 65-85-0 | Benzoic Acid | 0.77 | U | 129-00-0 | Pyrene | 0.77 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.77 | U | 959-98-8 | Endosulfan I | 0.77 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.77 | U | 92-87-5 | Benzidine | 0.77 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.77 | U | 72-55-9 | 4,4'-DDE | 0.77 | U |
| 91-20-3 | Naphthalene | 0.77 | U | 60-57-1 | Dieldrin | 0.77 | U |
| 106-47-8 | 4-Chloroaniline | 0.77 | U | 72-20-8 | Endrin | 0.77 | U |
| 87-68-3 | Hexachlorobutadiene | 0.77 | U | 33213-65-9 | Endosulfan II | 0.77 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.77 | U | 72-54-8 | 4,4'-DDD | 0.77 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.77 | U | 7421-92-4 | Endrin aldehyde | 0.77 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.77 | U | 85-68-7 | Butyl benzenyl phthalate | 0.77 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.77 | U | 1031-07-8 | Endosulfan sulfate | 0.77 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.77 | U | 50-29-3 | 4,4'-DDT | 0.77 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.77 | U | 56-55-3 | Benzo(a)anthracene | 0.77 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.77 | U | 218-01-9 | Chrysene | 0.77 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.77 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.77 | U |
| 88-74-4 | 2-Nitroaniline | 0.77 | U | 72-43-5 | Methoxychlor | 0.77 | U |
| 131-11-3 | Dimethyl phthalate | 0.77 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.77 | U |
| 208-96-8 | Acenaphthylene | 0.77 | U | 117-84-0 | Di-n-octyl phthalate | 0.77 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.77 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.77 | U |
| 99-09-2 | 3-Nitroaniline | 0.77 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.77 | U |
| 83-32-9 | Acenaphthene | 0.77 | U | 50-32-8 | Benzo(a)pyrene | 0.77 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.77 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.77 | U |
| 132-64-9 | Dibenzofuran | 0.77 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.77 | U |
| 100-02-7 | 4-Nitrophenol | 0.77 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.77 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.77 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.77 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | BKS00147A7 | Matrix: | Water |
| Sample ID No.: | OL-ER | Sample Vol. (ml) | 1040 |
| Date Collected: | 03/23/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/24/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | 1001010.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 53% | 24-132% | 65-135% |
| Phenol-d5 | 49% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 96% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 84% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 28% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 96% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.77 | U |
| 100-52-7 | Benzaldehyde | 0.77 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.77 | U |
| 95-53-4 | o-Toluidine | 0.77 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.77 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.77 | U |
| 1888-71-7 | Hexachloropropene | 0.77 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.77 | U |
| 87-65-0 | Caprolactam | 0.77 | U |
| 106-50-3 | p-Phenylenediamine | 0.77 | U |
| 120-58-1 | Isosafrole | 0.77 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.77 | U |
| 94-59-7 | Safrole | 0.77 | U |
| 92-52-4 | Biphenyl | 0.77 | U |
| 101-84-8 | Diphenyl ether | 0.77 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.77 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.77 | U |
| 608-93-5 | Pentachlorobenzene | 0.77 | U |
| 134-32-7 | 1-Naphthalenamine | 0.77 | U |
| 91-59-8 | 2-Naphthalenamine | 0.77 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.77 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.77 | U |
| 122-39-4 | Diphenylamine | 0.77 | U |
| 3689-24-5 | Sulfotep | 0.77 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.77 | U |
| 298-02-2 | Phorate | 0.77 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.77 | U |
| 62-44-2 | Phenacetin | 0.77 | U |
| 60-51-5 | Dimethoate | 0.77 | U |
| 1912-24-9 | Atrazine | 0.77 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.77 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.77 | U |
| 23950-58-! | Pronamide | 0.77 | U |
| 298-04-4 | Disulfoton | 0.77 | U |
| 298-00-0 | Methyl parathion | 0.77 | U |
| 56-38-2 | Parathion | 0.77 | U |
| 91-80-5 | Methapyrilene | 0.77 | U |
| 465-73-6 | Isodrin | 0.77 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.77 | U |
| 510-15-0 | Chlorobenzilate | 0.77 | U |
| 143-50-0 | Kepone | 0.77 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.77 | U |
| 52-85-7 | Famphur | 0.77 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.77 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.77 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.77 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.77 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
10/14/10chy/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | 3/25/11-LCS | Matrix: | Water |
| Sample ID No.: | Lab Control Spike | Sample Vol. (ml) | 1000 |
| Date Collected: | n/a | Final Volume (ml) | 2.0 |
| Date Received: | n/a | Dilution Factor: | 1 |
| Date Analyzed: | 04/06/11 | Analyst: | T. Meadows |
| Date of Report: | 04/08/11 | Supervisor's Initials: | |
| Data File Path: | Q:\5973\8270\110406\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 110% | 24-132% | 65-135% |
| Phenol-d5 | 68% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 87% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 76% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 37% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 139% | 80-141% | 65-135% |

| Matix Spike Recoveries | %Rec. | QC limits | |
|--------------------------|-------|-----------|---------|
| | | Water | Soil |
| Phenol | 66% | 50-150% | 50-150% |
| 2-Chlorophenol | 77% | 50-150% | 50-150% |
| 1,4-Dichlorobenzene | 60% | 50-150% | 50-150% |
| N-Nitroso-n-propyl amine | 57% | 50-150% | 50-150% |
| 1,2,4-Trichlorobenzene | 59% | 50-150% | 50-150% |
| 4-Chloro-3-methyl phenol | 77% | 50-150% | 50-150% |
| Acenaphthene | 82% | 50-150% | 50-150% |
| 4-Nitrophenol | 47% | 50-150% | 50-150% |
| 2,4-Dinitrotoluene | 50% | 50-150% | 50-150% |
| Pentachlorophenol | 0% | 50-150% | 50-150% |
| Pyrene | 115% | 50-150% | 50-150% |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **3/25/11-MB**
Sample ID No.: **Method Blank**
Date Collected: **n/a**
Date Received: **n/a**
Date Extracted: **03/25/11**
Date Analyzed: **04/06/11**
Date of Report: **04/08/11**
Data File Path: **Q:\5973\8270\110406**

Matrix: **Water**
Sample Vol. (ml) **1000**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **0401004.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.80 | U | 86-73-7 | Fluorene | 0.80 | U |
| 76-01-7 | Pentachloroethane | 0.80 | U | 84-66-2 | Diethyl phthalate | 0.80 | U |
| 62-53-3 | Aniline | 0.80 | U | 100-01-6 | 4-Nitroaniline | 0.80 | U |
| 108-95-2 | Phenol | 0.80 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.80 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.80 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.80 | U |
| 95-57-8 | 2-Chlorophenol | 0.80 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.80 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 103-33-3 | Azobenzene | 0.80 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.80 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.80 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.80 | U | 319-84-6 | a-BHC | 0.80 | U |
| 100-51-6 | Benzyl Alcohol | 0.80 | U | 118-74-1 | Hexachlorobenzene | 0.80 | U |
| 95-48-7 | 2-Methyl phenol | 0.80 | U | 319-85-7 | b-BHC | 0.80 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.80 | U | 87-86-5 | Pentachlorophenol | 0.80 | U |
| 98-86-2 | Acetophenone | 0.80 | U | 58-89-9 | g-BHC (Lindane) | 0.80 | U |
| 67-72-1 | Hexachloroethane | 0.80 | U | 85-01-8 | Phenanthrene | 0.80 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.80 | U | 120-12-7 | Anthracene | 0.80 | U |
| 106-44-5 | 4-Methyl phenol | 0.80 | U | 319-86-8 | d-BHC | 0.80 | U |
| 98-95-3 | Nitrobenzene | 0.80 | U | 84-74-2 | Di-n-butyl phthalate | 0.80 | U |
| 100-75-4 | N-nitrosopiperidine | 0.80 | U | 76-44-8 | Heptachlor | 0.80 | U |
| 78-59-1 | Isophorone | 0.80 | U | 309-00-2 | Aldrin | 0.80 | U |
| 88-75-5 | 2-Nitrophenol | 0.80 | U | 1024-57-3 | Heptachlor epoxide | 0.80 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.80 | U | 206-44-0 | Fluoranthrene | 0.80 | U |
| 65-85-0 | Benzoic Acid | 0.80 | U | 129-00-0 | Pyrene | 0.80 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.80 | U | 959-98-8 | Endosulfan I | 0.80 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.80 | U | 92-87-5 | Benzidine | 0.80 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.80 | U | 72-55-9 | 4,4'-DDE | 0.80 | U |
| 91-20-3 | Naphthalene | 0.80 | U | 60-57-1 | Dieldrin | 0.80 | U |
| 106-47-8 | 4-Chloroaniline | 0.80 | U | 72-20-8 | Endrin | 0.80 | U |
| 87-68-3 | Hexachlorobutadiene | 0.80 | U | 33213-65-9 | Endosulfan II | 0.80 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.80 | U | 72-54-8 | 4,4'-DDD | 0.80 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.80 | U | 7421-92-4 | Endrin aldehyde | 0.80 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.80 | U | 85-68-7 | Butyl benzenyl phthalate | 0.80 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.80 | U | 1031-07-8 | Endosulfan sulfate | 0.80 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.80 | U | 50-29-3 | 4,4'-DDT | 0.80 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.80 | U | 56-55-3 | Benzo(a)anthracene | 0.80 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.80 | U | 218-01-9 | Chrysene | 0.80 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.80 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.80 | U |
| 88-74-4 | 2-Nitroaniline | 0.80 | U | 72-43-5 | Methoxychlor | 0.80 | U |
| 131-11-3 | Dimethyl phthalate | 0.80 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.80 | U |
| 208-96-8 | Acenaphthylene | 0.80 | U | 117-84-0 | Di-n-octyl phthalate | 0.80 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.80 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.80 | U |
| 99-09-2 | 3-Nitroaniline | 0.80 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.80 | U |
| 83-32-9 | Acenaphthene | 0.80 | U | 50-32-8 | Benzo(a)pyrene | 0.80 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.80 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.80 | U |
| 132-64-9 | Dibenzofuran | 0.80 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.80 | U |
| 100-02-7 | 4-Nitrophenol | 0.80 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.80 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.80 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.80 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-----------------------------|------------------------|-------------------|
| Case File Number: | 3/25/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Vol. (ml) | 1000 |
| Date Collected: | n/a | Final Volume (ml) | 2.0 |
| Date Received: | n/a | Dilution Factor: | 1 |
| Date Extracted: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/06/11 | Supervisor's Initials: | |
| Date of Report: | 04/08/11 | | |
| Data File Path: | Q:\5973\8270\110406\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 50% | 24-132% | 65-135% |
| Phenol-d5 | 47% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 92% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 86% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 41% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 118% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.80 | U |
| 100-52-7 | Benzaldehyde | 0.80 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.80 | U |
| 95-53-4 | o-Toluidine | 0.80 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.80 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.80 | U |
| 1888-71-7 | Hexachloropropene | 0.80 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.80 | U |
| 87-65-0 | Caprolactam | 0.80 | U |
| 106-50-3 | p-Phenylenediamine | 0.80 | U |
| 120-58-1 | Isosafrole | 0.80 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.80 | U |
| 94-59-7 | Safrole | 0.80 | U |
| 92-52-4 | Biphenyl | 0.80 | U |
| 101-84-8 | Diphenyl ether | 0.80 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.80 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.80 | U |
| 608-93-5 | Pentachlorobenzene | 0.80 | U |
| 134-32-7 | 1-Naphthalenamine | 0.80 | U |
| 91-59-8 | 2-Naphthalenamine | 0.80 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.80 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.80 | U |
| 122-39-4 | Diphenylamine | 0.80 | U |
| 3689-24-5 | Sulfotep | 0.80 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.80 | U |
| 298-02-2 | Phorate | 0.80 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.80 | U |
| 62-44-2 | Phenacetin | 0.80 | U |
| 60-51-5 | Dimethoate | 0.80 | U |
| 1912-24-9 | Atrazine | 0.80 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.80 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.80 | U |
| 23950-58-! | Pronamide | 0.80 | U |
| 298-04-4 | Disulfoton | 0.80 | U |
| 298-00-0 | Methyl parathion | 0.80 | U |
| 56-38-2 | Parathion | 0.80 | U |
| 91-80-5 | Methapyrilene | 0.80 | U |
| 465-73-6 | Isodrin | 0.80 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.80 | U |
| 510-15-0 | Chlorobenzilate | 0.80 | U |
| 143-50-0 | Kepone | 0.80 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.80 | U |
| 52-85-7 | Famphur | 0.80 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.80 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.80 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.80 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.80 | U |



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| | | |
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| CASE FILE NUMBER: | BKS001-47,48 | PAGE 1 |
| REPORT DATE: | 04/06/11 | |
| DATE SAMPLED: | 03/23,24/11 | DATE RECEIVED: 03/24,25/11 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

CASE NARRATIVE

Fifteen water samples were received by the laboratory in good condition. Analysis was performed according to the chain of custody received with the samples. Dissolved arsenic analysis was performed by ICP/MS by concentrating the samples 10 fold. No difficulties were encountered in the preparation or analysis of these samples. Sample data follows while QA/QC data is contained on subsequent pages. Organic data are included as separate reports.

SAMPLE DATA

| SAMPLE ID | AMMONIA (mg/L) | NITRATE (mg/L) | NITRITE (mg/L) | TOC (mg/L) | COD (mg/L) | VINYL CHLORIDE (ug/L) |
|-----------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| OL-MW-1 | <0.010 | 0.254 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-MW-5A | <0.010 | 0.192 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-MW-2 | <0.010 | 0.389 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-MW-4 | <0.010 | 0.536 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-SW-3 | | 0.080 | 0.002 | | | |
| OL-SW-4 | | 0.147 | 0.002 | | | |
| OL-ER | <0.010 | <0.010 | <0.002 | <0.250 | <10.0 | 0.05 |
| OL-TB | | | | | | 0.09 |
| OL-MW-7 | 0.012 | 0.549 | <0.002 | 0.370 | <10.0 | <0.02 |
| OL-MW-8 | <0.010 | 0.131 | <0.002 | <0.250 | <10.0 | <0.02 |
| OL-MW-6 | <0.010 | 0.011 | <0.002 | 1.14 | <10.0 | <0.02 |
| OL-MW-10 | 0.017 | 0.292 | <0.002 | 3.04 | <10.0 | <0.02 |
| OL-MW-3 | <0.010 | 0.200 | 0.003 | 1.28 | <10.0 | <0.02 |
| OL-TB | | | | | | 0.05 |
| OL-MW-11 | <0.010 | 0.550 | <0.002 | 0.400 | <10.0 | <0.02 |



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SAMPLE DATA-CONTINUED

| SAMPLE ID | CHLORIDE (mg/L) | SULFATE (mg/L) | TDS (mg/L) | CYANIDE (mg/L) | TOTAL COLIFORM (#/100mls) | FECAL COLIFORM (#/100mls) |
|-----------|--------------------|-------------------|---------------|-------------------|------------------------------|------------------------------|
| OL-MW-1 | 2.64 | 4.03 | 85.0 | <0.010 | <2 | |
| OL-MW-5A | 2.05 | 3.82 | 148 | <0.010 | <2 | |
| OL-MW-2 | 2.15 | 2.62 | 100 | <0.010 | 2 | |
| OL-MW-4 | 2.54 | 2.46 | 101 | <0.010 | 4 | |
| OL-SW-3 | | | | | | 2 |
| OL-SW-4 | | | | | | <2 |
| OL-ER | <0.50 | <1.00 | 21.0 | <0.010 | 6 | |
| OL-MW-7 | 1.25 | 3.08 | 109 | <0.010 | <2 | |
| OL-MW-8 | 1.17 | 1.42 | 118 | <0.010 | <2 | |
| OL-MW-6 | 1.25 | 5.84 | 109 | <0.010 | <2 | |
| OL-MW-10 | 3.07 | 27.0 | 326 | <0.010 | <2 | |
| OL-MW-3 | 2.74 | 49.2 | 212 | <0.010 | <2 | |
| OL-MW-11 | 1.27 | 4.32 | 103 | <0.010 | <2 | |

| | pH | ALKALINITY (mgCaCO3/L) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|----------|------|---------------------------|---------------------|--------------------|
| OL-MW-1 | 6.58 | 60.8 | 38.3 | <1.00 |
| OL-MW-5A | 6.92 | 72.7 | 57.3 | <1.00 |
| OL-MW-2 | 7.13 | 55.2 | 47.3 | <1.00 |
| OL-MW-4 | 7.17 | 68.1 | 59.1 | <1.00 |
| OL-ER | 6.23 | <1.00 | <1.00 | <1.00 |
| OL-MW-7 | 6.98 | 67.0 | 54.3 | <1.00 |
| OL-MW-8 | 7.01 | 59.0 | 48.4 | <1.00 |
| OL-MW-6 | 6.92 | 67.5 | 53.2 | <1.00 |
| OL-MW-10 | 6.58 | 257 | 162 | <1.00 |
| OL-MW-3 | 6.36 | 134 | 67.8 | <1.00 |
| OL-MW-11 | 6.96 | 65.1 | 52.2 | <1.00 |



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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

SAMPLE DATA-CONTINUED

| DISSOLVED METALS | | | | | | |
|------------------|-------------------|--------------------|------------------|---------------------|-------------------|--------------------|
| SAMPLE ID | ARSENIC (ug/L) | ANTIMONY (mg/L) | BARIUM (mg/L) | BERYLLIUM (mg/L) | CADMIUM (mg/L) | CHROMIUM (mg/L) |
| OL-MW-1 | 0.082 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0023 |
| OL-MW-5A | 0.113 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0036 |
| OL-MW-2 | 0.517 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0034 |
| OL-MW-4 | 0.188 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0024 |
| OL-MW-7 | 0.327 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0035 |
| OL-MW-8 | 1.49 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0013 |
| OL-MW-6 | 0.689 | <0.0020 | <0.005 | <0.0010 | <0.0010 | <0.0010 |
| OL-MW-10 | 1.03 | <0.0020 | 0.016 | <0.0010 | <0.0010 | 0.0020 |
| OL-MW-3 | 0.087 | <0.0020 | 0.012 | <0.0010 | <0.0010 | 0.0014 |
| OL-MW-11 | 0.349 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0034 |

| DISSOLVED METALS | | | | | | |
|------------------|------------------|------------------|----------------|--------------------|------------------|--------------------|
| SAMPLE ID | COBALT (mg/L) | COPPER (mg/L) | LEAD (mg/L) | SELENIUM (mg/L) | SILVER (mg/L) | THALLIUM (mg/L) |
| OL-MW-1 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-5A | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-2 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-4 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-7 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-8 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-6 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-10 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-3 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| OL-MW-11 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |

| DISSOLVED METALS | | | |
|------------------|---------------|--------------------|----------------|
| SAMPLE ID | TIN (mg/L) | VANADIUM (mg/L) | ZINC (mg/L) |
| OL-MW-1 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-5A | <0.040 | 0.0030 | <0.0050 |
| OL-MW-2 | <0.040 | 0.0040 | <0.0050 |
| OL-MW-4 | <0.040 | 0.0032 | <0.0050 |
| OL-MW-7 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-8 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-6 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-10 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-3 | <0.040 | <0.0030 | <0.0050 |
| OL-MW-11 | <0.040 | <0.0030 | <0.0050 |



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SAMPLE DATA-CONTINUED

| SAMPLE ID | DISSOLVED METALS | | | | | |
|-----------|-------------------|---------------------|------------------|---------------------|----------------|---------------------|
| | CALCIUM (mg/L) | POTASSIUM (mg/L) | SODIUM (mg/L) | MAGNESIUM (mg/L) | IRON (mg/L) | MANGANESE (mg/L) |
| OL-MW-1 | 8.96 | 0.711 | 4.05 | 5.84 | <0.020 | <0.010 |
| OL-MW-5A | 8.73 | 0.811 | 4.49 | 8.45 | <0.020 | <0.010 |
| OL-MW-2 | 6.76 | 1.06 | 4.02 | 6.01 | <0.020 | <0.010 |
| OL-MW-4 | 8.54 | 0.942 | 4.54 | 7.64 | <0.020 | <0.010 |
| OL-MW-7 | 8.69 | 0.818 | 3.83 | 7.00 | <0.020 | <0.010 |
| OL-MW-8 | 9.28 | 0.680 | 4.11 | 4.69 | <0.020 | 0.143 |
| OL-MW-6 | 8.88 | 0.890 | 4.82 | 6.43 | 0.316 | 0.412 |
| OL-MW-10 | 39.3 | 1.44 | 10.8 | 24.8 | 0.084 | 4.85 |
| OL-MW-3 | 31.7 | 0.871 | 6.44 | 11.7 | <0.020 | 1.33 |
| OL-MW-11 | 8.86 | 0.821 | 3.83 | 7.21 | <0.020 | <0.010 |

| SAMPLE ID | TOTAL METALS | |
|-----------|-------------------|------------------|
| | MERCURY (mg/L) | NICKEL (mg/L) |
| OL-MW-1 | <0.0002 | <0.0050 |
| OL-MW-5A | <0.0002 | <0.0050 |
| OL-MW-2 | <0.0002 | <0.0050 |
| OL-MW-4 | <0.0002 | <0.0050 |
| OL-ER | <0.0002 | <0.0050 |
| OL-MW-7 | <0.0002 | <0.0050 |
| OL-MW-8 | <0.0002 | <0.0050 |
| OL-MW-6 | <0.0002 | <0.0050 |
| OL-MW-10 | <0.0002 | 0.0227 |
| OL-MW-3 | <0.0002 | <0.0050 |
| OL-MW-11 | <0.0002 | <0.0050 |



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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | AMMONIA (mg/L) | NITRATE (mg/L) | NITRITE (mg/L) | TOC (mg/L) | COD (mg/L) | VINYL CHLORIDE (ug/L) |
|-----------------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| METHOD | EPA 350.1 | EPA 353.2 | EPA 354.1 | EPA 415.1 | EPA 410.2 | EPA 8260-SIM |
| DATE ANALYZED | 03/24,25/11 | 03/24,25/11 | 03/25/11 | 04/01/11 | 04/01/11 | 04/04/11 |
| DETECTION LIMIT | 0.006 | 0.005 | 0.001 | 0.100 | 6.00 | 0.01 |
| REPORTING LIMIT | 0.010 | 0.010 | 0.002 | 0.250 | 10.0 | 0.02 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-10 | OL-ER | OL-MW-11 | OL-MW-11 | OL-MW-6 |
| ORIGINAL | 0.017 | 0.292 | <0.002 | 0.400 | <10.0 | <0.02 |
| DUPLICATE | 0.017 | 0.294 | <0.002 | 0.380 | <10.0 | <0.02 |
| RPD | 2.48% | 0.63% | NC | 5.07% | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-10 | OL-ER | OL-MW-11 | OL-MW-11 | |
| ORIGINAL | 0.017 | 0.292 | <0.002 | 0.400 | <10.0 | |
| SPIKED SAMPLE | 0.218 | 0.498 | 0.040 | 4.79 | 50.0 | |
| SPIKE ADDED | 0.200 | 0.200 | 0.040 | 4.50 | 50.0 | |
| % RECOVERY | 100.27% | 102.75% | 100.00% | 97.46% | 100.05% | NA |
| QC CHECK | | | | | | |
| FOUND | 0.313 | 0.400 | 0.041 | 4.12 | 94.4 | 0.10 |
| TRUE | 0.324 | 0.408 | 0.040 | 4.00 | 100 | 0.10 |
| % RECOVERY | 96.51% | 98.05% | 102.50% | 102.93% | 94.40% | 100.00% |
| BLANK | | | | | | |
| | <0.010 | <0.010 | <0.002 | <0.250 | <10.0 | <0.02 |

RPD = RELATIVE PERCENT DIFFERENCE.
 NA = NOT APPLICABLE OR NOT AVAILABLE.
 NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.
 OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.



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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | CHLORIDE (mg/l) | SULFATE (mg/l) | TDS (mg/L) | CYANIDE (mg/L) | TOTAL COLIFORM (#/100mls) | FECAL COLIFORM (#/100mls) |
|-----------------|--------------------|-------------------|---------------|-------------------|------------------------------|------------------------------|
| METHOD | EPA 325.3 | EPA 375.4 | SM18 2540C | SM 4500CNE | SM18 9222B | SM18 9222D |
| DATE ANALYZED | 04/04/11 | 03/30/11 | 03/29/11 | 03/30/11 | 03/24,25/11 | 03/24/11 |
| DETECTION LIMIT | 0.16 | 0.76 | 5.0 | 0.010 | 2 | 2 |
| REPORTING LIMIT | 0.50 | 1.00 | 5.0 | 0.010 | 2 | 2 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-11 | OL-MW-10 | OL-MW-10 | OL-MW-10 | OL-SW-4 |
| ORIGINAL | 3.07 | 4.32 | 326 | <0.010 | <2 | <2 |
| DUPLICATE | 3.03 | 4.46 | 316 | <0.010 | <2 | <2 |
| RPD | 1.19% | 3.00% | 3.12% | NC | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-11 | | OL-MW-10 | | |
| ORIGINAL | 3.07 | 4.32 | | <0.010 | | |
| SPIKED SAMPLE | 12.7 | 14.5 | | 0.179 | | |
| SPIKE ADDED | 10.0 | 10.0 | | 0.200 | | |
| % RECOVERY | 96.41% | 101.53% | NA | 89.50% | NA | NA |
| QC CHECK | | | | | | |
| FOUND | 28.3 | 9.95 | | 0.184 | | |
| TRUE | 30.0 | 10.0 | | 0.200 | | |
| % RECOVERY | 94.49% | 99.51% | NA | 92.00% | NA | NA |
| PREP BLANK | | | | | | |
| | <0.50 | <1.00 | <5.0 | <0.010 | < 1 | < 1 |

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QA/QC DATA

| QC PARAMETER | pH | ALKALINITY (mgCaCO3/l) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|-----------------|-------------|---------------------------|---------------------|--------------------|
| METHOD | EPA 150.1 | EPA 310.1 | EPA 310.1 | EPA 310.1 |
| DATE ANALYZED | 03/24,25/11 | 03/28/11 | 03/28/11 | 03/28/11 |
| DETECTION LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| REPORTING LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| DUPLICATE | | | | |
| SAMPLE ID | | OL-MW-10 | | |
| ORIGINAL | | 257 | | |
| DUPLICATE | | 265 | | |
| RPD | NA | 3.01% | NA | NA |
| SPIKE SAMPLE | | | | |
| SAMPLE ID | | | | |
| ORIGINAL | | | | |
| SPIKED SAMPLE | | | | |
| SPIKE ADDED | | | | |
| % RECOVERY | NA | NA | NA | NA |
| QC CHECK | | | | |
| FOUND | | 102 | | |
| TRUE | | 100 | | |
| % RECOVERY | NA | 101.50% | NA | NA |
| PREP BLANK | NA | NA | NA | NA |

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| REPORT DATE: | 04/06/11 | |
| DATE SAMPLED: | 03/23,24/11 | 03/24,25/11 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | ARSENIC (ug/l) | ANTIMONY (mg/L) | BARIUM (mg/L) | BERYLLIUM (mg/L) | CADMIUM (mg/L) | CHROMIUM (mg/L) |
|-----------------|-------------------|--------------------|------------------|---------------------|-------------------|--------------------|
| METHOD | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 03/28,30/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 |
| DETECTION LIMIT | 0.010 | 0.0010 | 0.001 | 0.0003 | 0.0002 | 0.0005 |
| REPORTING LIMIT | 0.050 | 0.0020 | 0.005 | 0.0010 | 0.0010 | 0.0010 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | 1.03 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0024 |
| DUPLICATE | 1.07 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0025 |
| RPD | 3.81% | NC | NC | NC | NC | 1.30% |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-10 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | 1.03 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0024 |
| SPIKED SAMPLE | 1.21 | 0.0534 | 0.052 | 0.0555 | 0.0564 | 0.0491 |
| SPIKE ADDED | 0.200 | 0.0500 | 0.050 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 90.00% | 106.82% | 103.94% | 111.00% | 112.76% | 93.35% |
| QC CHECK | | | | | | |
| FOUND | 0.702 | 0.0489 | 0.048 | 0.0512 | 0.0470 | 0.0497 |
| TRUE | 0.668 | 0.0500 | 0.050 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 105.09% | 97.74% | 95.00% | 102.32% | 93.92% | 99.32% |
| PREP BLANK | | | | | | |
| | <0.050 | <0.0020 | <0.010 | <0.0010 | <0.0010 | <0.0010 |

RPD = RELATIVE PERCENT DIFFERENCE.
 NA = NOT APPLICABLE OR NOT AVAILABLE.
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AQUATIC RESEARCH INCORPORATED
LABORATORY & CONSULTING SERVICES
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| | | |
|--|--------------|---------------|
| CASE FILE NUMBER: | BKS001-47,48 | PAGE 9 |
| REPORT DATE: | 04/06/11 | |
| DATE SAMPLED: | 03/23,24/11 | 03/24,25/11 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | COBALT (mg/L) | COPPER (mg/L) | LEAD (mg/L) | SELENIUM (mg/L) | SILVER (mg/L) | THALLIUM (mg/L) |
|-----------------|------------------|------------------|----------------|--------------------|------------------|--------------------|
| METHOD | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 | 03/25,28/11 |
| DETECTION LIMIT | 0.001 | 0.0004 | 0.0003 | 0.0010 | 0.0004 | 0.0010 |
| REPORTING LIMIT | 0.010 | 0.0010 | 0.0010 | 0.0050 | 0.0010 | 0.0020 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| DUPLICATE | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| RPD | NC | NC | NC | NC | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| SPIKED SAMPLE | 0.048 | 0.0487 | 0.0467 | 0.0598 | 0.0513 | 0.0475 |
| SPIKE ADDED | 0.050 | 0.0500 | 0.0500 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 96.12% | 97.34% | 93.34% | 119.50% | 102.68% | 94.90% |
| QC CHECK | | | | | | |
| FOUND | 0.051 | 0.0515 | 0.0505 | 0.0487 | 0.0520 | 0.0480 |
| TRUE | 0.050 | 0.0500 | 0.0500 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 101.04% | 102.92% | 100.90% | 97.30% | 103.92% | 96.08% |
| PREP BLANK | | | | | | |
| | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |

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| | | |
|--|--------------|----------------|
| CASE FILE NUMBER: | BKS001-47,48 | PAGE 10 |
| REPORT DATE: | 04/06/11 | |
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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | TIN (mg/L) | VANADIUM (mg/L) | ZINC (mg/L) |
|-----------------|---------------|--------------------|----------------|
| METHOD | EPA 6010 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 03/25,29/11 | 03/25,28/11 | 03/25,28/11 |
| DETECTION LIMIT | 0.010 | 0.0010 | 0.0010 |
| REPORTING LIMIT | 0.040 | 0.0030 | 0.0050 |
| DUPLICATE | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | <0.040 | 0.0032 | <0.0050 |
| DUPLICATE | <0.040 | 0.0032 | <0.0050 |
| RPD | NC | 0.57% | NC |
| SPIKE SAMPLE | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | <0.040 | 0.0032 | <0.0050 |
| SPIKED SAMPLE | 1.06 | 0.0510 | 0.0580 |
| SPIKE ADDED | 1.00 | 0.0500 | 0.0500 |
| % RECOVERY | 105.73% | 95.56% | 115.92% |
| QC CHECK | | | |
| FOUND | 1.08 | 0.0496 | 0.0474 |
| TRUE | 1.00 | 0.0500 | 0.0500 |
| % RECOVERY | 108.20% | 99.14% | 94.86% |
| PREP BLANK | | | |
| | <0.040 | <0.0030 | <0.0050 |

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| | | |
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| CASE FILE NUMBER: | BKS001-47,48 | PAGE 11 |
| REPORT DATE: | 04/06/11 | |
| DATE SAMPLED: | 03/23,24/11 | DATE RECEIVED: 03/24,25/11 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | CALCIUM (mg/L) | POTASSIUM (mg/L) | SODIUM (mg/L) | MAGNESIUM (mg/L) | IRON (mg/L) | MANGANESE (mg/L) |
|-----------------|-------------------|---------------------|------------------|---------------------|----------------|---------------------|
| METHOD | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 |
| DATE ANALYZED | 03/25,29/11 | 03/25,29/11 | 03/25,29/11 | 03/25,29/11 | 03/25,29/11 | 03/25,29/11 |
| DETECTION LIMIT | 0.050 | 0.300 | 0.300 | 0.050 | 0.010 | 0.003 |
| REPORTING LIMIT | 0.100 | 0.500 | 0.500 | 0.100 | 0.020 | 0.010 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | 8.54 | 0.942 | 4.54 | 7.64 | <0.020 | <0.010 |
| DUPLICATE | 8.53 | 0.948 | 4.51 | 7.62 | <0.020 | <0.010 |
| RPD | 0.14% | 0.58% | 0.56% | 0.24% | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 | OL-MW-4 |
| ORIGINAL | 8.54 | 0.942 | 4.54 | 7.64 | <0.020 | <0.010 |
| SPIKED SAMPLE | 16.6 | 12.7 | 14.0 | 16.6 | 9.87 | 1.02 |
| SPIKE ADDED | 10.0 | 10.0 | 10.0 | 10.0 | 10.0 | 1.00 |
| % RECOVERY | 80.91% | 117.46% | 94.18% | 89.60% | 98.70% | 102.00% |
| QC CHECK | | | | | | |
| FOUND | 0.977 | 9.81 | 10.3 | 10.0 | 0.981 | 1.02 |
| TRUE | 1.00 | 10.0 | 10.0 | 10.0 | 1.00 | 1.00 |
| % RECOVERY | 97.66% | 98.08% | 103.37% | 100.04% | 98.11% | 101.65% |
| PREP BLANK | <0.100 | <0.500 | <0.500 | <0.100 | <0.020 | <0.010 |

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|---|--------------|----------------------------|
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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - TOTAL METALS

| QC PARAMETER | MERCURY (mg/l) | NICKEL (mg/l) |
|-----------------|-------------------|------------------|
| METHOD | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 03/25,28/11 | 03/25,28/11 |
| DETECTION LIMIT | 0.0001 | 0.0010 |
| REPORTING LIMIT | 0.0002 | 0.0050 |
| DUPLICATE | | |
| SAMPLE ID | OL-MW-1 | OL-MW-1 |
| ORIGINAL | <0.0002 | <0.0050 |
| DUPLICATE | <0.0002 | <0.0050 |
| RPD | NC | NC |
| SPIKE SAMPLE | | |
| SAMPLE ID | OL-MW-1 | OL-MW-1 |
| ORIGINAL | <0.0002 | <0.0050 |
| SPIKED SAMPLE | 0.0022 | 0.0517 |
| SPIKE ADDED | 0.0020 | 0.0500 |
| % RECOVERY | 110.00% | 103.46% |
| QC CHECK | | |
| FOUND | 0.0021 | 0.0503 |
| TRUE | 0.0020 | 0.0500 |
| % RECOVERY | 104.00% | 100.62% |
| PREP BLANK | <0.0002 | <0.0050 |

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OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.

Submitted By:

Steven Lazoff
Laboratory Director



Aquatic Research, Inc.

3927 Aurora Ave. N. , Seattle, WA 98103 | (206) 632-2715

TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-1 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 85% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-2 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 109% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-4 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 71% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-5A | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 70% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00147A7 | Matrix: | Water |
| Sample ID No.: | OL-ER | Sample Vol. (ml) | 990 |
| Date Collected: | 3/23/11 | Dilution Factor: | 1 |
| Date Received: | 3/24/11 | Prepped: | TM |
| Date Prepped: | 3/25/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 93% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs LABORATORY FORTIFIED BLANK REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|---------------|-------------------------------|-------|
| Case File Number: | LFB (3/25/11) | Matrix: | Water |
| Sample ID No.: | LFB | Sample Vol. (ml) | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Arochlor | RECOVERY | LCL | UCL |
|------------|----------|-----|------|
| Total PCBs | 75% | 70% | 130% |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 82% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs LABORATORY REAGENT BLANK REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|--------------|-------------------------------|-------|
| Case File Number: | MB (3/25/11) | Matrix: | Water |
| Sample ID No.: | LRB | Sample Vol. (ml): | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Aroclor | Result (ug/L) | UCL (ug/L) | RL | (ug/L) |
|------------|------------------|---------------|----|--------|
| Total PCBs | < 0.1 | < 0.1 | | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 73% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit

Vinyl Chloride by SIM

T. Meadows

4/4/11

| Sample ID | Client ID | Sample Amount (ml) | Dilution Factor | Vinyl Chloride (µg/L) | 1,2-Dichloroethane-d4 (µg/L) | % R Surr | % R VC |
|----------------|--------------|--------------------|-----------------|-----------------------|------------------------------|----------|--------|
| CCV | CCV | 25 | 1 | 0.10 | 0.11 | 110.0% | 100.0% |
| CCV | CCV | 25 | 1 | 0.09 | 0.08 | 80.0% | 90.0% |
| 4/4/11-MB | Method Blank | 25 | 1 | <0.02 | 0.11 | 110.0% | |
| BKS00147A1 | OL-MW-1 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00147A2 | OL-MW-5A | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00147A3 | OL-MW-2 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00147A4 | OL-MW-4 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00147A7 | OL-ER | 25 | 1 | 0.05 | 0.07 | 70.0% | |
| BKS00147A8 | OL-TB | 25 | 1 | 0.09 | 0.08 | 80.0% | |
| BKS00148A1 | OL-MW-7 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A2 | OL-MW-8 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A3 | OL-MW-6 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A3 Dup | OL-MW-6 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A4 | OL-MW-10 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A5 | OL-MW-3 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00148A6 | OL-TB | 25 | 1 | 0.05 | 0.07 | 70.0% | |
| BKS00148A7 | OL-MW-11 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| | | | RPD | 0% | | 0% | |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-3 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-3 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 1201012.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 84% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 110% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 84% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00147A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0801008.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 78% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 107% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 82% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-7 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-7 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0601006.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 82% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 107% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 88% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



Aquatic Research Inc.
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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A2** Matrix: Water
Sample ID No.: **OL-MW-8** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **03/24/11** Dilution Factor: **1**
Date Received: **03/25/11**
Date Analyzed: **04/ 1/11** Analyst: **T. Meadows**
Date of Report: **04/01/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110401\ 0701007.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|-----------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01- | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02- | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-8 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0701007.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 81% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 109% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 79% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A4** Matrix: Water
Sample ID No.: **OL-MW-10** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **03/24/11** Dilution Factor: **1**
Date Received: **03/25/11**
Date Analyzed: **04/ 1/11** Analyst: **T. Meadows**
Date of Report: **04/01/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110401\ 0901009.D**

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0901009.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 73% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 101% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 80% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A7 | Matrix: | Water |
| Sample ID No.: | OL-MW-11 | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A7 | Matrix: | Water |
| Sample ID No.: | OL-MW-11 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 1401014.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 81% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 107% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 83% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



Aquatic Research Inc.
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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00148A4 MS / MSD | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | | |
| Date Analyzed: | 04/ 1/11 | Analyst: | T. Meadows |
| Date of Report: | 4/1/11 | Supervisor's Initials: | |
| Data File Paths: | C:\HPCHEM\1\DATA\VOA\110401\ 1001010.D | | |
| | C:\HPCHEM\1\DATA\VOA\110401\ 1101011.D | | |

| Surrogate Recoveries | % Rec. | | |
|----------------------|--------|-----|-----|
| | MS | MSD | RPD |
| Dibromofluoromethane | 72% | 70% | 3% |
| Toluene-d8 | 106% | 97% | 10% |
| 4-Bromofluorobenzene | 84% | 71% | 15% |

| Spike Recoveries | % Rec. | | |
|--------------------|--------|------|-----|
| | MS | MSD | RPD |
| 1,1-Dichloroethene | 116% | 114% | 2% |
| Benzene | 94% | 97% | 3% |
| Trichloroethene | 119% | 108% | 10% |
| Toluene | 106% | 105% | 1% |
| Chlorobenzene | 106% | 112% | 5% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|------------------------------|-------------------------|------------|
| Case File Number: | 4/1/11-LCS | Matrix: | Water |
| Sample ID No.: | 4/1/11-LCS | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed | 04/ 1/11 | Analyst: | T. Meadows |
| Date of Report: | 4/1/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 94% | 66% | 118% |
| Toluene-d8 | 97% | 51% | 143% |
| 4-Bromofluorobenzene | 83% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 91% | 75% | 125% |
| Benzene | 96% | 75% | 125% |
| Trichloroethene | 100% | 75% | 125% |
| Toluene | 99% | 75% | 125% |
| Chlorobenzene | 100% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00148A4 MS | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | | |
| Date Analyzed: | 04/ 1/11 | Analyst: | T. Meadows |
| Date of Report: | 4/1/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ 1001010.D | | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 72% | 66% | 118% |
| Toluene-d8 | 106% | 51% | 143% |
| 4-Bromofluorobenzene | 84% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 116% | 75% | 125% |
| Benzene | 94% | 75% | 125% |
| Trichloroethene | 119% | 75% | 125% |
| Toluene | 106% | 75% | 125% |
| Chlorobenzene | 106% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00148A4 MSD | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | | |
| Date Analyzed: | 04/ 1/11 | Analyst: | T. Meadows |
| Date of Report: | 4/1/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\1101011.D | | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 70% | 66% | 118% |
| Toluene-d8 | 97% | 51% | 143% |
| 4-Bromofluorobenzene | 71% | 63% | 119% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 114% | 75% | 125% |
| Benzene | 97% | 75% | 125% |
| Trichloroethene | 108% | 75% | 125% |
| Toluene | 105% | 75% | 125% |
| Chlorobenzene | 112% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | 4/1/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|-----------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01- | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02- | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | 4/1/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 0401004.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | | | | QC limits | | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | | |
| | Dibromofluoromethane | 108% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 105% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 91% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A6 | Matrix: | Water |
| Sample ID No.: | OL-TB | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U |
| 74-95-3 | Dibromomethane | 0.4 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U |
| 103-65-1 | n-Propylbenzene | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U |

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|---------------------------|--------------|------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 127-18-4 | Tetrachloroethene | 0.4 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 79-01-6 | Trichloroethene | 0.4 | U |
| 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 75-01-4 | Vinyl chloride | 0.4 | U |

BTEX

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------|--------------|------|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

TRIHALOMETHANES (THM)

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|----------------------|--------------|------|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

KETONES, CS2, 2-CEVE

| CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-------------------------|--------------|------|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00148A6 | Matrix: | Water |
| Sample ID No.: | OL-TB | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 03/24/11 | Dilution Factor: | 1 |
| Date Received: | 03/25/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/ 1/11 | Supervisor's Initials: | |
| Date of Report: | 04/01/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110401\ | | 1301013.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|-----------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | Water | Soil | QC limits | | | |
| | Dibromofluoromethane | 74% | 66-118% | 66-118% | | | | |
| | Toluene-d8 | 108% | 51-143% | 51-143% | | | | |
| | 4-Bromofluorobenzene | 82% | 63 - 119% | 63 - 119% | | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.



Aquatic Research Inc.
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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A5**
Sample ID No.: **OL-MW-3**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1060**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3701019.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-3 | Sample Vol. (ml) | 1060 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purges Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 39% | 24-132% | 65-135% |
| Phenol-d5 | 39% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 99% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 82% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 28% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 110% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A3**
Sample ID No.: **OL-MW-6**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3301015.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6 | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 61% | 24-132% | 65-135% |
| Phenol-d5 | 57% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 119% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 113% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 32% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 127% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A1**
Sample ID No.: **OL-MW-7**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/07/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1060**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials:
3101013.D

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-7 | Sample Vol. (ml) | 1060 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/07/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 77% | 24-132% | 65-135% |
| Phenol-d5 | 62% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 129% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 95% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 42% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 118% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A2**
Sample ID No.: **OL-MW-8**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials:
3201014.D

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-8 | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | 3201014.D |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 41% | 24-132% | 65-135% |
| Phenol-d5 | 36% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 133% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 101% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 36% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 114% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|----------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-1 | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracene | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A4**
Sample ID No.: **OL-MW-10**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1060**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3401016.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Vol. (ml) | 1060 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 10% | 24-132% | 65-135% |
| Phenol-d5 | 22% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 127% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 114% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 19% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 125% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00148A7**
Sample ID No.: **OL-MW-11**
Date Collected: **03/24/11**
Date Received: **03/25/11**
Date Extracted: **03/28/11**
Date Analyzed: **04/08/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1070**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **3801020.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.75 | U | 86-73-7 | Fluorene | 0.75 | U |
| 76-01-7 | Pentachloroethane | 0.75 | U | 84-66-2 | Diethyl phthalate | 0.75 | U |
| 62-53-3 | Aniline | 0.75 | U | 100-01-6 | 4-Nitroaniline | 0.75 | U |
| 108-95-2 | Phenol | 0.75 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.75 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.75 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.75 | U |
| 95-57-8 | 2-Chlorophenol | 0.75 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.75 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.75 | U | 103-33-3 | Azobenzene | 0.75 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.75 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.75 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.75 | U | 319-84-6 | a-BHC | 0.75 | U |
| 100-51-6 | Benzyl Alcohol | 0.75 | U | 118-74-1 | Hexachlorobenzene | 0.75 | U |
| 95-48-7 | 2-Methyl phenol | 0.75 | U | 319-85-7 | b-BHC | 0.75 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.75 | U | 87-86-5 | Pentachlorophenol | 0.75 | U |
| 98-86-2 | Acetophenone | 0.75 | U | 58-89-9 | g-BHC (Lindane) | 0.75 | U |
| 67-72-1 | Hexachloroethane | 0.75 | U | 85-01-8 | Phenanthrene | 0.75 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.75 | U | 120-12-7 | Anthracene | 0.75 | U |
| 106-44-5 | 4-Methyl phenol | 0.75 | U | 319-86-8 | d-BHC | 0.75 | U |
| 98-95-3 | Nitrobenzene | 0.75 | U | 84-74-2 | Di-n-butyl phthalate | 0.75 | U |
| 100-75-4 | N-nitrosopiperidine | 0.75 | U | 76-44-8 | Heptachlor | 0.75 | U |
| 78-59-1 | Isophorone | 0.75 | U | 309-00-2 | Aldrin | 0.75 | U |
| 88-75-5 | 2-Nitrophenol | 0.75 | U | 1024-57-3 | Heptachlor epoxide | 0.75 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.75 | U | 206-44-0 | Fluoranthrene | 0.75 | U |
| 65-85-0 | Benzoic Acid | 0.75 | U | 129-00-0 | Pyrene | 0.75 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.75 | U | 959-98-8 | Endosulfan I | 0.75 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.75 | U | 92-87-5 | Benzidine | 0.75 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.75 | U | 72-55-9 | 4,4'-DDE | 0.75 | U |
| 91-20-3 | Naphthalene | 0.75 | U | 60-57-1 | Dieldrin | 0.75 | U |
| 106-47-8 | 4-Chloroaniline | 0.75 | U | 72-20-8 | Endrin | 0.75 | U |
| 87-68-3 | Hexachlorobutadiene | 0.75 | U | 33213-65-9 | Endosulfan II | 0.75 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.75 | U | 72-54-8 | 4,4'-DDD | 0.75 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.75 | U | 7421-92-4 | Endrin aldehyde | 0.75 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.75 | U | 85-68-7 | Butyl benzenyl phthalate | 0.75 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.75 | U | 1031-07-8 | Endosulfan sulfate | 0.75 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.75 | U | 50-29-3 | 4,4'-DDT | 0.75 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.75 | U | 56-55-3 | Benzo(a)anthracene | 0.75 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.75 | U | 218-01-9 | Chrysene | 0.75 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.75 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.75 | U |
| 88-74-4 | 2-Nitroaniline | 0.75 | U | 72-43-5 | Methoxychlor | 0.75 | U |
| 131-11-3 | Dimethyl phthalate | 0.75 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.75 | U |
| 208-96-8 | Acenaphthylene | 0.75 | U | 117-84-0 | Di-n-octyl phthalate | 0.75 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.75 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.75 | U |
| 99-09-2 | 3-Nitroaniline | 0.75 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.75 | U |
| 83-32-9 | Acenaphthene | 0.75 | U | 50-32-8 | Benzo(a)pyrene | 0.75 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.75 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.75 | U |
| 132-64-9 | Dibenzofuran | 0.75 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.75 | U |
| 100-02-7 | 4-Nitrophenol | 0.75 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.75 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.75 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.75 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A7 | Matrix: | Water |
| Sample ID No.: | OL-MW-11 | Sample Vol. (ml) | 1070 |
| Date Collected: | 03/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 03/25/11 | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/08/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 44% | 24-132% | 65-135% |
| Phenol-d5 | 37% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 125% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 102% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 47% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 118% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.75 | U |
| 100-52-7 | Benzaldehyde | 0.75 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.75 | U |
| 95-53-4 | o-Toluidine | 0.75 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.75 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.75 | U |
| 1888-71-7 | Hexachloropropene | 0.75 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.75 | U |
| 87-65-0 | Caprolactam | 0.75 | U |
| 106-50-3 | p-Phenylenediamine | 0.75 | U |
| 120-58-1 | Isosafrole | 0.75 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.75 | U |
| 94-59-7 | Safrole | 0.75 | U |
| 92-52-4 | Biphenyl | 0.75 | U |
| 101-84-8 | Diphenyl ether | 0.75 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.75 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.75 | U |
| 608-93-5 | Pentachlorobenzene | 0.75 | U |
| 134-32-7 | 1-Naphthalenamine | 0.75 | U |
| 91-59-8 | 2-Naphthalenamine | 0.75 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.75 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.75 | U |
| 122-39-4 | Diphenylamine | 0.75 | U |
| 3689-24-5 | Sulfotep | 0.75 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.75 | U |
| 298-02-2 | Phorate | 0.75 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.75 | U |
| 62-44-2 | Phenacetin | 0.75 | U |
| 60-51-5 | Dimethoate | 0.75 | U |
| 1912-24-9 | Atrazine | 0.75 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.75 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.75 | U |
| 23950-58-! | Pronamide | 0.75 | U |
| 298-04-4 | Disulfoton | 0.75 | U |
| 298-00-0 | Methyl parathion | 0.75 | U |
| 56-38-2 | Parathion | 0.75 | U |
| 91-80-5 | Methapyrilene | 0.75 | U |
| 465-73-6 | Isodrin | 0.75 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.75 | U |
| 510-15-0 | Chlorobenzilate | 0.75 | U |
| 143-50-0 | Kepone | 0.75 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.75 | U |
| 52-85-7 | Famphur | 0.75 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.75 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.75 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.75 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.75 | U |



Aquatic Research Inc.
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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
10/14/10chy/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | 3/28/11-LCS | Matrix: | Water |
| Sample ID No.: | Lab Control Spike | Sample Vol. (ml) | 1000 |
| Date Collected: | n/a | Final Volume (ml) | 2.0 |
| Date Received: | n/a | Dilution Factor: | 1 |
| Date Analyzed: | 04/07/11 | Analyst: | T. Meadows |
| Date of Report: | 04/11/11 | Supervisor's Initials: | |
| Data File Path: | Q:\5973\PAM\040711\ | 3001012.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 87% | 24-132% | 65-135% |
| Phenol-d5 | 59% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 116% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 106% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 26% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 119% | 80-141% | 65-135% |

| Matix Spike Recoveries | %Rec. | QC limits | |
|--------------------------|-------|-----------|---------|
| | | Water | Soil |
| Phenol | 52% | 50-150% | 50-150% |
| 2-Chlorophenol | 58% | 50-150% | 50-150% |
| 1,4-Dichlorobenzene | 76% | 50-150% | 50-150% |
| N-Nitroso-n-propyl amine | 84% | 50-150% | 50-150% |
| 1,2,4-Trichlorobenzene | 68% | 50-150% | 50-150% |
| 4-Chloro-3-methyl phenol | 48% | 50-150% | 50-150% |
| Acenaphthene | 96% | 50-150% | 50-150% |
| 4-Nitrophenol | 44% | 50-150% | 50-150% |
| 2,4-Dinitrotoluene | 48% | 50-150% | 50-150% |
| Pentachlorophenol | 0% | 50-150% | 50-150% |
| Pyrene | 90% | 50-150% | 50-150% |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
10/14/10ghy/Mass Spectrometry

| | | | |
|-------------------|------------------------------|------------------------|-------------------|
| Case File Number: | BKS00148A4 MS | Matrix: | Water |
| Sample ID No.: | OL-MW-10 Matrix Spike | Sample Vol. (ml) | 1070 |
| Date Collected: | 3/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 3/25/11 | Dilution Factor: | 1 |
| Date Analyzed: | 04/08/11 | Analyst: | T. Meadows |
| Date of Report: | 04/11/11 | Supervisor's Initials: | |
| Data File Path: | Q:\5973\PAM\040711\ | 3501017.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 27% | 24-132% | 65-135% |
| Phenol-d5 | 29% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 108% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 95% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 12% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 114% | 80-141% | 65-135% |

| Matix Spike Recoveries | %Rec. | QC limits | |
|--------------------------|-------|-----------|---------|
| | | Water | Soil |
| Phenol | 35% | 50-150% | 50-150% |
| 2-Chlorophenol | 37% | 50-150% | 50-150% |
| 1,4-Dichlorobenzene | 76% | 50-150% | 50-150% |
| N-Nitroso-n-propyl amine | 100% | 50-150% | 50-150% |
| 1,2,4-Trichlorobenzene | 63% | 50-150% | 50-150% |
| 4-Chloro-3-methyl phenol | 33% | 50-150% | 50-150% |
| Acenaphthene | 104% | 50-150% | 50-150% |
| 4-Nitrophenol | 51% | 50-150% | 50-150% |
| 2,4-Dinitrotoluene | 62% | 50-150% | 50-150% |
| Pentachlorophenol | 4% | 50-150% | 50-150% |
| Pyrene | 91% | 50-150% | 50-150% |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
10/14/10chy/Mass Spectrometry

| | | | |
|-------------------|--|------------------------|-------------------|
| Case File Number: | BKS00148A4 MSD | Matrix: | Water |
| Sample ID No.: | OL-MW-10 Matrix Spike Duplicate | Sample Vol. (ml) | 1070 |
| Date Collected: | 3/24/11 | Final Volume (ml) | 2.0 |
| Date Received: | 3/25/11 | Dilution Factor: | 1 |
| Date Analyzed: | 04/08/11 | Analyst: | T. Meadows |
| Date of Report: | 04/11/11 | Supervisor's Initials: | |
| Data File Path: | Q:\5973\PAM\040711\ | 3601018.D | |

| Surrogate Recoveries | %Rec. | QC limits | | RPD |
|----------------------|-------|-----------|---------|-------|
| | | Water | Soil | |
| 2-Fluorophenol | 25% | 24-132% | 65-135% | 5.0% |
| Phenol-d5 | 30% | 0-105% | 65-135% | 5.0% |
| Nitrobenzene-d5 | 125% | 24-136% | 65-135% | 14.2% |
| 2-Fluorobiphenyl | 109% | 41-134% | 65-135% | 13.7% |
| 2,4,6-Tribromophenol | 10% | 0-183% | 65-135% | 19.8% |
| p-Terphenyl-d14 | 115% | 80-141% | 65-135% | 0.8% |

| Matix Spike Recoveries | %Rec. | QC limits | | RPD |
|--------------------------|-------|-----------|---------|-------|
| | | Water | Soil | |
| Phenol | 40% | 50-150% | 50-150% | 12.3% |
| 2-Chlorophenol | 40% | 50-150% | 50-150% | 7.0% |
| 1,4-Dichlorobenzene | 90% | 50-150% | 50-150% | 17.0% |
| N-Nitroso-n-propyl amine | 104% | 50-150% | 50-150% | 4.5% |
| 1,2,4-Trichlorobenzene | 75% | 50-150% | 50-150% | 17.2% |
| 4-Chloro-3-methyl phenol | 35% | 50-150% | 50-150% | 6.9% |
| Acenaphthene | 109% | 50-150% | 50-150% | 4.3% |
| 4-Nitrophenol | 50% | 50-150% | 50-150% | 1.3% |
| 2,4-Dinitrotoluene | 64% | 50-150% | 50-150% | 3.0% |
| Pentachlorophenol | 4% | 50-150% | 50-150% | 8.4% |
| Pyrene | 90% | 50-150% | 50-150% | 0.4% |



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SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **3/28/11-MB**
Sample ID No.: **Method Blank**
Date Collected: **n/a**
Date Received: **n/a**
Date Extracted: **03/28/11**
Date Analyzed: **04/07/11**
Date of Report: **04/11/11**
Data File Path: **Q:\5973\PAM\040711**

Matrix: **Water**
Sample Vol. (ml) **1000**
Final Volume (ml) **2.0**
Dilution Factor: **1**

Analyst: **T. Meadows**
Supervisor's Initials: **2901011.D**

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|------------------------------|--------------|------|------------|-----------------------------|--------------|------|
| 55-18-5 | N-nitrosodiethylamine | 0.80 | U | 86-73-7 | Fluorene | 0.80 | U |
| 76-01-7 | Pentachloroethane | 0.80 | U | 84-66-2 | Diethyl phthalate | 0.80 | U |
| 62-53-3 | Aniline | 0.80 | U | 100-01-6 | 4-Nitroaniline | 0.80 | U |
| 108-95-2 | Phenol | 0.80 | U | 7005-72-3 | 4-Chlorophenyl phenyl ether | 0.80 | U |
| 111-44-4 | Bis(2-chloroethyl) ether | 0.80 | U | 534-52-1 | 2-Methyl-4,6-dinitrophenol | 0.80 | U |
| 95-57-8 | 2-Chlorophenol | 0.80 | U | 86-30-6 | N-Nitrosodiphenyl amine | 0.80 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 0.80 | U | 103-33-3 | Azobenzene | 0.80 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 0.80 | U | 101-55-3 | 4-Bromophenyl phenyl ether | 0.80 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 0.80 | U | 319-84-6 | a-BHC | 0.80 | U |
| 100-51-6 | Benzyl Alcohol | 0.80 | U | 118-74-1 | Hexachlorobenzene | 0.80 | U |
| 95-48-7 | 2-Methyl phenol | 0.80 | U | 319-85-7 | b-BHC | 0.80 | U |
| 108-60-1 | Bis(2-chloroisopropyl) ether | 0.80 | U | 87-86-5 | Pentachlorophenol | 0.80 | U |
| 98-86-2 | Acetophenone | 0.80 | U | 58-89-9 | g-BHC (Lindane) | 0.80 | U |
| 67-72-1 | Hexachloroethane | 0.80 | U | 85-01-8 | Phenanthrene | 0.80 | U |
| 621-64-7 | N-Nitroso-n-propyl amine | 0.80 | U | 120-12-7 | Anthracene | 0.80 | U |
| 106-44-5 | 4-Methyl phenol | 0.80 | U | 319-86-8 | d-BHC | 0.80 | U |
| 98-95-3 | Nitrobenzene | 0.80 | U | 84-74-2 | Di-n-butyl phthalate | 0.80 | U |
| 100-75-4 | N-nitrosopiperidine | 0.80 | U | 76-44-8 | Heptachlor | 0.80 | U |
| 78-59-1 | Isophorone | 0.80 | U | 309-00-2 | Aldrin | 0.80 | U |
| 88-75-5 | 2-Nitrophenol | 0.80 | U | 1024-57-3 | Heptachlor epoxide | 0.80 | U |
| 105-67-9 | 2,4-Dimethylphenol | 0.80 | U | 206-44-0 | Fluoranthrene | 0.80 | U |
| 65-85-0 | Benzoic Acid | 0.80 | U | 129-00-0 | Pyrene | 0.80 | U |
| 111-91-1 | Bis(2-chloroethoxy) methane | 0.80 | U | 959-98-8 | Endosulfan I | 0.80 | U |
| 120-83-2 | 2,4-Dichlorophenol | 0.80 | U | 92-87-5 | Benzidine | 0.80 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.80 | U | 72-55-9 | 4,4'-DDE | 0.80 | U |
| 91-20-3 | Naphthalene | 0.80 | U | 60-57-1 | Dieldrin | 0.80 | U |
| 106-47-8 | 4-Chloroaniline | 0.80 | U | 72-20-8 | Endrin | 0.80 | U |
| 87-68-3 | Hexachlorobutadiene | 0.80 | U | 33213-65-9 | Endosulfan II | 0.80 | U |
| 924-16-3 | N-nitrosodi-n-butylamine | 0.80 | U | 72-54-8 | 4,4'-DDD | 0.80 | U |
| 59-50-7 | 4-Chloro-3-methyl phenol | 0.80 | U | 7421-92-4 | Endrin aldehyde | 0.80 | U |
| 91-57-6 | 2-Methyl naphthalene | 0.80 | U | 85-68-7 | Butyl benzenyl phthalate | 0.80 | U |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 0.80 | U | 1031-07-8 | Endosulfan sulfate | 0.80 | U |
| 77-47-4 | Hexachlorocyclopentadiene | 0.80 | U | 50-29-3 | 4,4'-DDT | 0.80 | U |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.80 | U | 56-55-3 | Benzo(a)anthracene | 0.80 | U |
| 95-95-4 | 2,4,5-Trichlorophenol | 0.80 | U | 218-01-9 | Chrysene | 0.80 | U |
| 91-58-7 | 2-Chloronaphthalene | 0.80 | U | 91-94-1 | 3,3'-Dichlorobenzidine | 0.80 | U |
| 88-74-4 | 2-Nitroaniline | 0.80 | U | 72-43-5 | Methoxychlor | 0.80 | U |
| 131-11-3 | Dimethyl phthalate | 0.80 | U | 117-81-7 | Bis(2-ethylhexyl) phthalate | 0.80 | U |
| 208-96-8 | Acenaphthylene | 0.80 | U | 117-84-0 | Di-n-octyl phthalate | 0.80 | U |
| 606-20-2 | 2,6-Dinitrotoluene | 0.80 | U | 205-97-2 | Benzo(b)fluoranthrene | 0.80 | U |
| 99-09-2 | 3-Nitroaniline | 0.80 | U | 207-08-9 | Benzo(k)fluoranthrene | 0.80 | U |
| 83-32-9 | Acenaphthene | 0.80 | U | 50-32-8 | Benzo(a)pyrene | 0.80 | U |
| 51-28-5 | 2,4-Dinitrophenol | 0.80 | U | 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.80 | U |
| 132-64-9 | Dibenzofuran | 0.80 | U | 53-70-3 | Dibenzo(a,h)anthracene | 0.80 | U |
| 100-02-7 | 4-Nitrophenol | 0.80 | U | 91-24-2 | Benzo(g,h,i)perylene | 0.80 | U |
| 121-14+2 | 2,4-Dinitrotoluene | 0.80 | U | | | | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 0.80 | U | | | | |



SEMI-VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8270
Measurement of Extractable Organic Compounds in Water by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|----------------------------|------------------------|-------------------|
| Case File Number: | 3/28/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Vol. (ml) | 1000 |
| Date Collected: | n/a | Final Volume (ml) | 2.0 |
| Date Received: | n/a | Dilution Factor: | 1 |
| Date Extracted: | 03/28/11 | Analyst: | T. Meadows |
| Date Analyzed: | 04/07/11 | Supervisor's Initials: | |
| Date of Report: | 04/11/11 | | |
| Data File Path: | Q:\5973\PAM\040711\ | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|---|---|--------------|------|------|------------------|--------------|------|
| FLAGS: | | | | | | | |
| U | Indicates compound was analyzed for, but not detected at the specified detection limit. | | | | | | |
| B | Blank contaminated with this analyte. | | | | | | |
| J | Estimated value - compound positively identified, but below specified detection limit. | | | | | | |
| E | Estimated value - compound exceeded calibration range. | | | | | | |
| D | Compound analyzed at a secondary dilution factor of _____ from data file: _____ | | | | | | |
| PP | Compound Purged Poorly, requiring elevated detection limit. | | | | | | |
| NOTE: ppm Amounts are in mg/L or mg/KG dry weight. | | | | | | | |

Tentatively Identified Compounds

R.T. CAS# Compound Rel. Conc.

The above compounds have been tentatively identified as present. In general, specific identities of isomers of alkanes/alkenes is not possible to resolve.

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|---------|
| | | Water | Soil |
| 2-Fluorophenol | 79% | 24-132% | 65-135% |
| Phenol-d5 | 78% | 0-105% | 65-135% |
| Nitrobenzene-d5 | 128% | 24-136% | 65-135% |
| 2-Fluorobiphenyl | 102% | 41-134% | 65-135% |
| 2,4,6-Tribromophenol | 36% | 0-183% | 65-135% |
| p-Terphenyl-d14 | 110% | 80-141% | 65-135% |

| | | | |
|-----------|----------------------------------|------|---|
| 62-50-0 | Ethyl methanesulfonic acid | 0.80 | U |
| 100-52-7 | Benzaldehyde | 0.80 | U |
| 930-55-2 | N-nitrosopyrrolidine | 0.80 | U |
| 95-53-4 | o-Toluidine | 0.80 | U |
| 106-44-5 | 3 + 4-Methyl phenol (unresolved) | 0.80 | U |
| 126-68-1 | O,O,O-Triethyl Phosphorothioate | 0.80 | U |
| 1888-71-7 | Hexachloropropene | 0.80 | U |
| 87-65-0 | 2,6-Dichlorophenol | 0.80 | U |
| 87-65-0 | Caprolactam | 0.80 | U |
| 106-50-3 | p-Phenylenediamine | 0.80 | U |
| 120-58-1 | Isosafrole | 0.80 | U |
| 90-12-0 | 1-Methyl naphthalene | 0.80 | U |
| 94-59-7 | Safrole | 0.80 | U |
| 92-52-4 | Biphenyl | 0.80 | U |
| 101-84-8 | Diphenyl ether | 0.80 | U |
| 130-15-4 | 1,4-Naphthoquinone | 0.80 | U |
| 99-65-0 | 1,3-Dinitrobenzene | 0.80 | U |
| 608-93-5 | Pentachlorobenzene | 0.80 | U |
| 134-32-7 | 1-Naphthalenamine | 0.80 | U |
| 91-59-8 | 2-Naphthalenamine | 0.80 | U |
| 99-55-8 | 5-Nitro-o-Toluidine | 0.80 | U |
| 297-97-2 | Thionazin (Zinophos) | 0.80 | U |
| 122-39-4 | Diphenylamine | 0.80 | U |
| 3689-24-5 | Sulfotep | 0.80 | U |

| | | | |
|------------|---------------------------------|------|---|
| 2303-16-4 | Diallate | 0.80 | U |
| 298-02-2 | Phorate | 0.80 | U |
| 99-35-4 | 1,3,5-Trinitrobenzene | 0.80 | U |
| 62-44-2 | Phenacetin | 0.80 | U |
| 60-51-5 | Dimethoate | 0.80 | U |
| 1912-24-9 | Atrazine | 0.80 | U |
| 92-67-1 | 4-Aminobiphenyl | 0.80 | U |
| 82-68-8 | Pentachloronitrobenzene | 0.80 | U |
| 23950-58-! | Pronamide | 0.80 | U |
| 298-04-4 | Disulfoton | 0.80 | U |
| 298-00-0 | Methyl parathion | 0.80 | U |
| 56-38-2 | Parathion | 0.80 | U |
| 91-80-5 | Methapyrilene | 0.80 | U |
| 465-73-6 | Isodrin | 0.80 | U |
| 60-11-7 | p-(dimethyl amino) azobenzene | 0.80 | U |
| 510-15-0 | Chlorobenzilate | 0.80 | U |
| 143-50-0 | Kepone | 0.80 | U |
| 119-93-7 | 3,3'-Dimethylbenzidine | 0.80 | U |
| 52-85-7 | Famphur | 0.80 | U |
| 53-96-3 | 2-Acetylaminofluorene | 0.80 | U |
| 57-97-6 | 7,12-Dimethyl benz(a) anthracen | 0.80 | U |
| 56-49-5 | 3-Methylcholanthrene | 0.80 | U |
| 226-36-8 | Dibenzo(a,h)acridine | 0.80 | U |



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-3 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/8/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 107% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 74% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



Aquatic Research, Inc.

3927 Aurora Ave. N. , Seattle, WA 98103 | (206) 632-2715

TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-7 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 103% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-8 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 119% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



Aquatic Research, Inc.

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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-10 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/7/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 115% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs SAMPLE REPORT

Results of Analysis by EPA Method 8082A

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|------------|-------------------------------|-------|
| Case File Number: | BKS00148A7 | Matrix: | Water |
| Sample ID No.: | OL-MW-11 | Sample Vol. (ml) | 990 |
| Date Collected: | 3/24/11 | Dilution Factor: | 1 |
| Date Received: | 3/25/11 | Prepped: | TM |
| Date Prepped: | 3/28/11 | Analyst: | JDS |
| Date Analyzed | 4/8/11 | Supervisor's Initials: | |
| Date of Report: | 4/8/11 | | |

Total PCBs

| Aroclor | Result (ug/L) | RL (ug/L) |
|----------------|----------------------|------------------|
| 1016 | < 0.1 | 0.1 |
| 1221 | < 0.1 | 0.1 |
| 1232 | < 0.1 | 0.1 |
| 1242 | < 0.1 | 0.1 |
| 1248 | < 0.1 | 0.1 |
| 1254 | < 0.1 | 0.1 |
| 1260 | < 0.1 | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 95% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs MATRIX SPIKE REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|---------------|-------------------------------|-------|
| Case File Number: | BKS00148A4 MS | Matrix: | Water |
| Sample ID No.: | OL-MW-10 MS | Sample Vol. (ml) | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Aroclor | RECOVERY | LCL | UCL |
|------------|----------|-----|------|
| Total PCBs | 97% | 70% | 130% |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 99% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs MATRIX SPIKE REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|----------------|-------------------------------|-------|
| Case File Number: | BKS00148A4 MSD | Matrix: | Water |
| Sample ID No.: | OL-MW-10 MSD | Sample Vol. (ml) | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/8/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Aroclor | RECOVERY | LCL | UCL |
|----------------|-----------------|------------|------------|
| Total PCBs | 104% | 70% | 130% |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|---------------|------------|------------|
| Decachlorobiphenyl (DCB) | 99% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs LABORATORY FORTIFIED BLANK REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|---------------|-------------------------------|-------|
| Case File Number: | LFB (3/28/11) | Matrix: | Water |
| Sample ID No.: | LFB | Sample Vol. (ml) | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Arochlor | RECOVERY | LCL | UCL |
|------------|----------|-----|------|
| Total PCBs | 121% | 70% | 130% |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 95% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



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TOTAL PCBs LABORATORY REAGENT BLANK REPORT

Results of Analysis by EPA Method 508.1

Measurement of Total PCBs in Water by GC/ECD

| | | | |
|--------------------------|--------------|-------------------------------|-------|
| Case File Number: | MB (3/28/11) | Matrix: | Water |
| Sample ID No.: | LRB | Sample Vol. (ml): | 1000 |
| Date Collected: | NA | Dilution Factor: | 1 |
| Date Received: | NA | Prepped: | TM |
| Date Analyzed | 4/7/11 | Analyst: | JDS |
| Date of Report: | 4/8/11 | Supervisor's Initials: | |

Total PCBs

| Aroclor | Result (ug/L) | UCL (ug/L) | RL | (ug/L) |
|------------|------------------|---------------|----|--------|
| Total PCBs | < 0.1 | < 0.1 | | 0.1 |

Surrogate Compounds

| Parameter | % Rec. | LCL | UCL |
|--------------------------|--------|-----|------|
| Decachlorobiphenyl (DCB) | 95% | 70% | 130% |

MCL: Maximum Contaminant Level

RL: Reporting Limit

LCL: Lower Control Limit

UCL: Upper Control Limit



Aquatic Research Inc.
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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00145A2 | Matrix: | Water |
| Sample ID No.: | MW-1 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\11010_0801008.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-3 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-4 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|-------------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRihalOMETHANES (THM) | | | |
|------------------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|-----------------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00145A2 | Matrix: | Water |
| Sample ID No.: | MW-1 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\11010_0801008.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|------|--------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | | Water | Soil | | |
| | Dibromofluoromethane | 71% | | 66-1189 | 66-118% | | |
| | Toluene-d8 | 93% | | 51-1439 | 51-143% | | |
| | 4-Bromofluorobenzene | 99% | | 46-1039 | 46-103% | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00145A2 MS | Matrix: | Water |
| Sample ID No.: | BKS00145A2 MS | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | | |
| Date Analyzed: | 1/4/11 | Analyst: | T. Meadows |
| Date of Report: | 1/5/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\ 0901009.D | | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 73% | 66% | 118% |
| Toluene-d8 | 92% | 51% | 143% |
| 4-Bromofluorobenzene | 86% | 46% | 103% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 84% | 75% | 125% |
| Benzene | 94% | 75% | 125% |
| Trichloroethene | 113% | 75% | 125% |
| Toluene | 116% | 75% | 125% |
| Chlorobenzene | 111% | 75% | 125% |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00145A4 | Matrix: | Water |
| Sample ID No.: | MW-2 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\11010 1101011.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------------------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| Surrogate Recoveries | | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 67% | 66-118 | 66-118% | | | |
| | Toluene-d8 | 93% | 51-143 | 51-143% | | | |
| | 4-Bromofluorobenzene | 95% | 46-103 | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00145A5 | Matrix: | Water |
| Sample ID No.: | MW-4 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\1101041201012.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | | Water | Soil | | |
| | Dibromofluoromethane | 76% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 96% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 94% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00145A3 | Matrix: | Water |
| Sample ID No.: | MW-5A | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\ 1001010.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 70% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 91% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 93% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00145A1 | Matrix: | Water |
| Sample ID No.: | MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104 0601006.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 70% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 91% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 85% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



Aquatic Research Inc.
3927 Aurora Ave. N., Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00145A1 Dup | Matrix: | Water |
| Sample ID No.: | MW-10 Duplicate | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\0701007.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRihalOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00145A1 Dup | Matrix: | Water |
| Sample ID No.: | MW-10 Duplicate | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\0701007.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | | Water | Soil | | |
| | Dibromofluoromethane | 68% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 89% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 86% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|------------------------|-------------------|
| Case File Number: | BKS00145A8 | Matrix: | Water |
| Sample ID No.: | SW-3 | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\ 1501015.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRihalomethanes (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00145A8 | Matrix: | Water |
| Sample ID No.: | SW-3 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\ 1501015.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | | QC limits | | | |
| | | | | Water | Soil | | |
| | Dibromofluoromethane | 72% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 93% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 88% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00145A6 | Matrix: | Water |
| Sample ID No.: | SW-2 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\ 1301013.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 74% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 92% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 99% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---------------------------------------|-------------------------|------------|
| Case File Number: | 1/4/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed: | 01/04/11 | Analyst: | T. Meadows |
| Date of Report: | 01/05/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\0401004.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIHALOMETHANES (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | 1/4/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed: | 01/04/11 | Analyst: | T. Meadows |
| Date of Report: | 01/05/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\0401004.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | | QC limits | | | |
| | | | Water | Soil | | | |
| | Dibromofluoromethane | 106% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 95% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 99% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|------------------------------|------------------------|------------|
| Case File Number: | 1/4/11-LCS | Matrix: | Water |
| Sample ID No.: | 1/4/11-LCS | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed | 1/4/11 | Analyst: | T. Meadows |
| Date of Report: | 1/5/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 103% | 66% | 118% |
| Toluene-d8 | 95% | 51% | 143% |
| 4-Bromofluorobenzene | 103% | 46% | 103% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 78% | 75% | 125% |
| Benzene | 90% | 75% | 125% |
| Trichloroethene | 89% | 75% | 125% |
| Toluene | 100% | 75% | 125% |
| Chlorobenzene | 97% | 75% | 125% |

Vinyl Chloride by SIM

T. Meadows

1/7/11

| Sample ID | Client ID | Sample Amount (ml) | Dilution Factor | Vinyl Chloride (µg/L) | 1,2-Dichloroethane-d4 (µg/L) | % R Surr | % R VC |
|------------|--------------|--------------------|-----------------|-----------------------|------------------------------|----------|--------|
| CCV | CCV | 25 | 1 | 0.10 | 0.10 | 100.0% | 100.0% |
| CCV | CCV | 25 | 1 | 0.09 | 0.09 | 90.0% | 90.0% |
| 1/6/11-MB | Method Blank | 25 | 1 | <0.02 | 0.10 | 100.0% | |
| BKS00145A1 | MW-10 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00145A2 | MW-1 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00145A3 | MW-5A | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00145A4 | MW-2 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00145A5 | MW-4 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00145A6 | SW-2 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00145A7 | SW-4 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00145A8 | SW-3 | 25 | 1 | <0.02 | 0.07 | 70.0% | |

RPD 0%



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| | | |
|---|-----------|-------------------------|
| CASE FILE NUMBER: | BKS001-45 | PAGE 1 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/28/10 | DATE RECEIVED: 12/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

CASE NARRATIVE

Eight water samples were received by the laboratory in good condition. Analysis was performed according to the chain of custody received with the samples. Dissolved arsenic analysis was performed by ICP/MS by concentrating the samples 10 fold. No difficulties were encountered in the preparation or analysis of these samples. Sample data follows while QA/QC data is contained on subsequent pages. VOC data and Vinyl Chloride QC data are included as separate reports.

SAMPLE DATA

| SAMPLE ID | AMMONIA (mg/l) | NITRATE (mg/l) | NITRITE (mg/l) | TOC (mg/l) | COD (mg/l) | VINYL CHLORIDE (ug/l) |
|-----------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| MW-10 | 0.022 | <0.010 | <0.002 | 2.86 | <10.0 | <0.02 |
| MW-1 | <0.010 | 0.122 | <0.002 | <0.250 | <10.0 | <0.02 |
| MW-5A | <0.010 | 0.441 | <0.002 | 0.328 | <10.0 | <0.02 |
| MW-2 | <0.010 | 0.369 | 0.002 | 0.429 | 86.5 | <0.02 |
| MW-4 | <0.010 | 0.550 | <0.002 | <0.250 | <10.0 | <0.02 |
| SW-2 | <0.010 | <0.010 | 0.003 | 7.95 | 24.2 | <0.02 |
| SW-4 | <0.010 | 0.227 | 0.003 | 8.55 | 25.5 | <0.02 |
| SW-3 | <0.010 | 0.171 | 0.002 | 5.83 | 17.8 | <0.02 |



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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-45 | PAGE 2 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/28/10 | 12/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

SAMPLE DATA-CONTINUED

| SAMPLE ID | TOTAL COLIFORM (#/100mls) | FECAL COLIFORM (#/100mls) | CHLORIDE (mg/l) | SULFATE (mg/l) |
|-----------|------------------------------|------------------------------|--------------------|-------------------|
| MW-10 | <1 | | 0.88 | 14.2 |
| MW-1 | <1 | | 2.40 | 3.69 |
| MW-5A | <1 | | 1.86 | 3.18 |
| MW-2 | <1 | | 1.76 | 3.34 |
| MW-4 | 10 | | 2.54 | 3.02 |
| SW-2 | | 4 | 1.56 | <1.00 |
| SW-4 | | 1 | 1.76 | <1.00 |
| SW-3 | | 46 | 2.15 | <1.00 |

| SAMPLE ID | pH | ALKALINITY (mgCaCO3/l) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|-----------|------|---------------------------|---------------------|--------------------|
| MW-10 | 6.57 | 196 | 122 | <1.00 |
| MW-1 | 6.41 | 66.3 | 35.5 | <1.00 |
| MW-5A | 6.53 | 103 | 62.2 | <1.00 |
| MW-2 | 7.03 | 58.1 | 48.1 | <1.00 |
| MW-4 | 7.13 | 60.6 | 51.9 | <1.00 |
| SW-2 | 7.17 | 18.3 | 15.9 | <1.00 |
| SW-4 | 6.78 | 15.8 | 11.5 | <1.00 |
| SW-3 | 6.30 | 6.18 | 2.91 | <1.00 |



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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-45 | PAGE 3 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/28/10 | 12/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

SAMPLE DATA-CONTINUED

| SAMPLE ID | DISSOLVED METALS | | | | |
|-----------|-------------------|------------------|----------------|---------------------|----------------|
| | ARSENIC (ug/l) | BARIUM (mg/l) | IRON (mg/l) | MANGANESE (mg/l) | ZINC (mg/l) |
| MW-10 | 1.05 | 0.007 | <0.020 | 3.34 | <0.005 |
| MW-1 | 0.098 | <0.005 | <0.020 | <0.005 | <0.005 |
| MW-5A | 0.160 | <0.005 | <0.020 | <0.005 | <0.005 |
| MW-2 | 0.652 | <0.005 | <0.020 | <0.005 | <0.005 |
| MW-4 | 0.216 | <0.005 | <0.020 | <0.005 | <0.005 |
| SW-2 | 0.737 | 0.007 | 0.087 | <0.005 | 0.018 |
| SW-4 | 0.566 | 0.005 | 0.105 | 0.005 | 0.019 |
| SW-3 | 0.230 | 0.006 | 0.091 | 0.007 | 0.006 |

| SAMPLE ID | METALS | | |
|-----------|-------------------|---------------------|------------------|
| | CALCIUM (mg/l) | POTASSIUM (mg/l) | SODIUM (mg/l) |
| MW-10 | 30.1 | 1.12 | 9.51 |
| MW-1 | 9.93 | 0.708 | 4.27 |
| MW-5A | 12.1 | 0.915 | 5.36 |
| MW-2 | 7.24 | 1.06 | 4.32 |
| MW-4 | 8.19 | 0.882 | 4.54 |
| SW-2 | 6.81 | <0.500 | 2.28 |
| SW-4 | 6.55 | <0.500 | 2.11 |
| SW-3 | 2.55 | <0.500 | 1.71 |



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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-45 | PAGE 4 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/28/10 | 12/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | AMMONIA (mg/l) | NITRATE (mg/l) | NITRITE (mg/l) | TOC (mg/l) | COD (mg/l) | VINYL CHLORIDE (ug/l) |
|-----------------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| METHOD | EPA 350.1 | EPA 353.2 | EPA 354.1 | EPA 415.1 | EPA 410.2 | EPA 8260-SIM |
| DATE ANALYZED | 12/30/10 | 12/30/10 | 12/30/10 | 01/11/11 | 01/11/11 | 01/06/11 |
| DETECTION LIMIT | 0.006 | 0.005 | 0.001 | 0.100 | 6.00 | 0.01 |
| REPORTING LIMIT | 0.010 | 0.010 | 0.002 | 0.250 | 10.0 | 0.02 |
| DUPLICATE | | | | | | |
| SAMPLE ID | BATCH | BATCH | SW-3 | MW-10 | BATCH | BATCH |
| ORIGINAL | <0.010 | 0.594 | 0.002 | 2.86 | <10.0 | <0.02 |
| DUPLICATE | <0.010 | 0.593 | 0.002 | 2.87 | <10.0 | <0.02 |
| RPD | NC | 0.21% | 0.00% | 0.14% | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | BATCH | BATCH | SW-3 | MW-10 | BATCH | |
| ORIGINAL | <0.010 | 0.594 | 0.002 | 2.86 | <10.0 | |
| SPIKED SAMPLE | 0.194 | 0.796 | 0.042 | 6.87 | 54.3 | |
| SPIKE ADDED | 0.200 | 0.200 | 0.040 | 4.50 | 50.0 | |
| % RECOVERY | 96.87% | 100.64% | 100.00% | 89.11% | 108.66% | NA |
| QC CHECK | | | | | | |
| FOUND | 0.332 | 0.409 | 0.040 | 3.88 | 97.7 | 0.10 |
| TRUE | 0.324 | 0.408 | 0.040 | 4.00 | 100 | 0.10 |
| % RECOVERY | 102.42% | 100.17% | 100.00% | 96.93% | 97.69% | 100.00% |
| BLANK | | | | | | |
| | <0.010 | <0.010 | <0.002 | <0.250 | <10.0 | <0.02 |

RPD = RELATIVE PERCENT DIFFERENCE.
 NA = NOT APPLICABLE OR NOT AVAILABLE.
 NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.
 OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.



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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-45 | PAGE 5 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/28/10 | 12/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | TOTAL COLIFORM (#/100mls) | FECAL COLIFORM (#/100mls) | CHLORIDE (mg/l) | SULFATE (mg/l) |
|-----------------|------------------------------|------------------------------|--------------------|-------------------|
| METHOD | SM18 9222B | SM18 9222D | EPA 325.3 | EPA 375.4 |
| DATE ANALYZED | 12/29/10 | 12/29/10 | 01/04/11 | 01/05/11 |
| DETECTION LIMIT | 1 | 1 | 0.16 | 0.76 |
| REPORTING LIMIT | 1 | 1 | 0.50 | 1.00 |
| DUPLICATE | | | | |
| SAMPLE ID | MW-4 | SW-3 | BATCH | BATCH |
| ORIGINAL | 10 | 46 | 1.37 | 3.50 |
| DUPLICATE | 7 | 51 | 1.37 | 3.42 |
| RPD | 35.29% | 10.31% | 0.00% | 2.31% |
| SPIKE SAMPLE | | | | |
| SAMPLE ID | | | BATCH | BATCH |
| ORIGINAL | | | 1.37 | 3.50 |
| SPIKED SAMPLE | | | 11.0 | 13.2 |
| SPIKE ADDED | | | 10.0 | 10.0 |
| % RECOVERY | NA | NA | 96.78% | 97.43% |
| QC CHECK | | | | |
| FOUND | | | 28.4 | 9.76 |
| TRUE | | | 30.0 | 10.0 |
| % RECOVERY | NA | NA | 94.82% | 97.56% |
| PREP BLANK | < 1 | < 1 | <0.50 | <1.00 |

RPD = RELATIVE PERCENT DIFFERENCE.
 NA = NOT APPLICABLE OR NOT AVAILABLE.
 NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.
 OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.



AQUATIC RESEARCH INCORPORATED
LABORATORY & CONSULTING SERVICES
 3927 AURORA AVENUE NORTH, SEATTLE, WA 98103
 PHONE: (206) 632-2715 FAX: (206) 632-2417

| | | |
|--|-----------|---------------|
| CASE FILE NUMBER: | BKS001-45 | PAGE 6 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/28/10 | 12/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | pH | ALKALINITY (mgCaCO3/l) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|-----------------|-----------|---------------------------|---------------------|--------------------|
| METHOD | EPA 150.1 | EPA 310.1 | EPA 310.1 | EPA 310.1 |
| DATE ANALYZED | 12/29/10 | 01/03/11 | 01/03/11 | 01/03/11 |
| DETECTION LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| REPORTING LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| DUPLICATE | | | | |
| SAMPLE ID | | BATCH | | |
| ORIGINAL | | 380 | | |
| DUPLICATE | | 392 | | |
| RPD | NA | 3.08% | NA | NA |
| SPIKE SAMPLE | | | | |
| SAMPLE ID | | | | |
| ORIGINAL | | | | |
| SPIKED SAMPLE | | | | |
| SPIKE ADDED | | | | |
| % RECOVERY | NA | NA | NA | NA |
| QC CHECK | | | | |
| FOUND | | 102 | | |
| TRUE | | 100 | | |
| % RECOVERY | NA | 102.20% | NA | NA |
| PREP BLANK | | | | |
| | NA | NA | NA | NA |

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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-45 | PAGE 7 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/28/10 | 12/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | ARSENIC (ug/l) | BARIUM (mg/l) | IRON (mg/l) | MANGANESE (mg/l) | ZINC (mg/l) |
|-----------------|-------------------|------------------|----------------|---------------------|----------------|
| METHOD | EPA 6020 | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 |
| DATE ANALYZED | 01/05/11 | 01/04/11 | 01/04/11 | 01/04/11 | 01/04/11 |
| DETECTION LIMIT | 0.010 | 0.002 | 0.008 | 0.003 | 0.003 |
| REPORTING LIMIT | 0.050 | 0.005 | 0.020 | 0.005 | 0.005 |
| DUPLICATE | | | | | |
| SAMPLE ID | BATCH | SW-3 | SW-3 | SW-3 | SW-3 |
| ORIGINAL | 0.318 | 0.006 | 0.091 | 0.007 | 0.006 |
| DUPLICATE | 0.365 | 0.006 | 0.092 | 0.007 | 0.006 |
| RPD | 13.76% | 1.77% | 1.42% | 2.86% | 3.17% |
| SPIKE SAMPLE | | | | | |
| SAMPLE ID | BATCH | SW-3 | SW-3 | SW-3 | SW-3 |
| ORIGINAL | 0.318 | 0.006 | 0.091 | 0.007 | 0.006 |
| SPIKED SAMPLE | 0.522 | 1.01 | 9.86 | 0.972 | 1.03 |
| SPIKE ADDED | 0.200 | 1.00 | 10.0 | 1.00 | 1.00 |
| % RECOVERY | 102.00% | 100.71% | 97.64% | 96.45% | 102.73% |
| QC CHECK | | | | | |
| FOUND | 0.721 | 1.02 | 1.05 | 1.01 | 1.02 |
| TRUE | 0.668 | 1.00 | 1.00 | 1.00 | 1.00 |
| % RECOVERY | 108.00% | 101.96% | 104.91% | 101.44% | 102.04% |
| PREP BLANK | <0.050 | <0.005 | <0.020 | <0.005 | <0.005 |

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| | | |
|---|-----------|-------------------------|
| CASE FILE NUMBER: | BKS001-45 | PAGE 8 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/28/10 | DATE RECEIVED: 12/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | CALCIUM (mg/l) | POTASSIUM (mg/l) | SODIUM (mg/l) |
|-----------------|-------------------|---------------------|------------------|
| METHOD | EPA 200.7 | EPA 200.7 | EPA 200.7 |
| DATE ANALYZED | 01/04/11 | 01/04/11 | 01/04/11 |
| DETECTION LIMIT | 0.050 | 0.300 | 0.300 |
| REPORTING LIMIT | 0.100 | 0.500 | 0.500 |
| DUPLICATE | | | |
| SAMPLE ID | SW-3 | SW-3 | SW-3 |
| ORIGINAL | 2.55 | <0.500 | 1.71 |
| DUPLICATE | 2.54 | <0.500 | 1.71 |
| RPD | 0.45% | NC | 0.42% |
| SPIKE SAMPLE | | | |
| SAMPLE ID | SW-3 | SW-3 | SW-3 |
| ORIGINAL | 2.55 | <0.500 | 1.71 |
| SPIKED SAMPLE | 12.3 | 11.9 | 12.6 |
| SPIKE ADDED | 10.0 | 10.0 | 10.0 |
| % RECOVERY | 97.47% | 119.47% | 109.16% |
| QC CHECK | | | |
| FOUND | 10.1 | 9.53 | 9.67 |
| TRUE | 10.0 | 10.0 | 10.0 |
| % RECOVERY | 101.20% | 95.28% | 96.71% |
| PREP BLANK | <0.100 | <0.500 | <0.500 |

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Submitted By:

Steven Lazoff
Laboratory Director



Aquatic Research Inc.
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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00145A7 | Matrix: | Water |
| Sample ID No.: | SW-4 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/04/11 | Supervisor's Initials: | |
| Date of Report: | 01/05/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\1401014.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIALOMETHANES (THM) | | | |
|----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00145A7 | Matrix: | Water |
| Sample ID No.: | SW-4 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/28/10 | Dilution Factor: | 1 |
| Date Received: | 12/29/10 | | |
| Date Analyzed: | 01/04/11 | Analyst: | T. Meadows |
| Date of Report: | 01/05/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110104\1401014.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------------------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| Surrogate Recoveries | | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 73% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 94% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 89% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



Aquatic Research Inc.
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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

Case File Number: **BKS00146A4 MS** Matrix: Water
Sample ID No.: **BKS00146A4 MS** Sample Wt/Vol. (gm/ml) **25.0**
Date Collected: **n/a** Dilution Factor: **1**
Date Received: **n/a**
Date Analyzed: **1/5/11** Analyst: **T. Meadows**
Date of Report: **1/6/11** Supervisor's Initials:
Data File Path: **C:\HPCHEM\1\DATA\VOA\110105\ 1101011.D**

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 65% | 66% | 118% |
| Toluene-d8 | 89% | 51% | 143% |
| 4-Bromofluorobenzene | 68% | 46% | 103% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 115% | 75% | 125% |
| Benzene | 83% | 75% | 125% |
| Trichloroethene | 102% | 75% | 125% |
| Toluene | 107% | 75% | 125% |
| Chlorobenzene | 102% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00146A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-8-0 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/29/10 | Dilution Factor: | 1 |
| Date Received: | 12/30/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/05/11 | Supervisor's Initials: | |
| Date of Report: | 01/06/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\1001010.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-3 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-4 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIALOMETHANES (THM) | | | |
|----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00146A4 | Matrix: | Water |
| Sample ID No.: | OL-MW-8-0 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/29/10 | Dilution Factor: | 1 |
| Date Received: | 12/30/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/05/11 | Supervisor's Initials: | |
| Date of Report: | 01/06/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\1001010.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | | QC limits | | | |
| | | | | Water | Soil | | |
| | Dibromofluoromethane | 67% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 87% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 72% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



Aquatic Research Inc.
3927 Aurora Ave. N., Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00146A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-7-0 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/29/10 | Dilution Factor: | 1 |
| Date Received: | 12/30/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/05/11 | Supervisor's Initials: | |
| Date of Report: | 01/06/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\1201012.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIALOMETHANES (THM) | | | |
|----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00146A5 | Matrix: | Water |
| Sample ID No.: | OL-MW-7-0 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/29/10 | Dilution Factor: | 1 |
| Date Received: | 12/30/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/05/11 | Supervisor's Initials: | |
| Date of Report: | 01/06/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\1201012.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 64% | 66-118 | 66-118% | | | |
| | Toluene-d8 | 92% | 51-143 | 51-143% | | | |
| | 4-Bromofluorobenzene | 78% | 46-103 | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00146A3 | Matrix: | Water |
| Sample ID No.: | OL-MW-6-1 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/29/10 | Dilution Factor: | 1 |
| Date Received: | 12/30/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/05/11 | Supervisor's Initials: | |
| Date of Report: | 01/06/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105 0901009.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 70% | 66-1189 | 66-118% | | | |
| | Toluene-d8 | 90% | 51-1439 | 51-143% | | | |
| | 4-Bromofluorobenzene | 79% | 46-1039 | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|------------------------|-------------------|
| Case File Number: | BKS00146A2 | Matrix: | Water |
| Sample ID No.: | OL-MW-6-0 | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/29/10 | Dilution Factor: | 1 |
| Date Received: | 12/30/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/05/11 | Supervisor's Initials: | |
| Date of Report: | 01/06/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\ 0801008.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | | QC limits | | | |
| | | | Water | Soil | | | |
| | Dibromofluoromethane | 64% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 91% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 87% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00146A1 Dup | Matrix: | Water |
| Sample ID No.: | OL-MW-3-0 Duplicate | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/29/10 | Dilution Factor: | 1 |
| Date Received: | 12/30/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/05/11 | Supervisor's Initials: | |
| Date of Report: | 01/06/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\ 0701007.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | | Water | Soil | | |
| | Dibromofluoromethane | 65% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 92% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 70% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in µg/L or µg/KG dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|------------------------|-------------------|
| Case File Number: | BKS00146A1 | Matrix: | Water |
| Sample ID No.: | OL-MW-3-0 | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 12/29/10 | Dilution Factor: | 1 |
| Date Received: | 12/30/10 | Analyst: | T. Meadows |
| Date Analyzed: | 01/05/11 | Supervisor's Initials: | |
| Date of Report: | 01/06/11 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\ 0601006.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 70% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 93% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 77% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---------------------------------------|------------------------|------------|
| Case File Number: | 1/5/11-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed: | 01/05/11 | Analyst: | T. Meadows |
| Date of Report: | 01/06/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\0401004.D | | |

| Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | QC limits | | | | |
| Surrogate Recoveries | %Rec. | Water | Soil | | | |
| Dibromofluoromethane | 102% | 66-118% | 66-118% | | | |
| Toluene-d8 | 91% | 51-143% | 51-143% | | | |
| 4-Bromofluorobenzene | 92% | 46-103% | 46-103% | | | |
| Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| Iodomethane | 0.4 | U | | | | |

Indicates compound was analyzed for, but not detected at the specified detection limit.
 Blank contaminated with this analyte.
 Estimated value - compound positively identified, but below specified detection limit.
 Estimated value - compound exceeded calibration range.
 Compound analyzed at a secondary dilution factor of _____ from data file: _____
 Compound Purges Poorly, requiring elevated detection limit.

ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | |
|-------------|-----------------|-------------------|
| <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-----------------|-------------------|



Aquatic Research Inc.
3927 Aurora Ave. N. , Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|------------------------------|-------------------------|------------|
| Case File Number: | 1/5/11-LCS | Matrix: | Water |
| Sample ID No.: | 1/5/11-LCS | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed: | 1/5/11 | Analyst: | T. Meadows |
| Date of Report: | 1/6/11 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\110105\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 108% | 66% | 118% |
| Toluene-d8 | 96% | 51% | 143% |
| 4-Bromofluorobenzene | 103% | 46% | 103% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 80% | 75% | 125% |
| Benzene | 88% | 75% | 125% |
| Trichloroethene | 92% | 75% | 125% |
| Toluene | 104% | 75% | 125% |
| Chlorobenzene | 102% | 75% | 125% |

Vinyl Chloride by SIM

T. Meadows

1/7/11

| Sample ID | Client ID | Sample Amount (ml) | Dilution Factor | Vinyl Chloride (µg/L) | 1,2- Dichloroethane- d4 (µg/L) | % R Surr | % R VC |
|----------------|--------------|--------------------------|--------------------|--------------------------|--------------------------------------|----------|--------|
| CCV | CCV | 25 | 1 | 0.09 | 0.11 | 110.0% | 90.0% |
| CCV | CCV | 25 | 1 | 0.10 | 0.09 | 90.0% | 100.0% |
| 1/7/11-MB | Method Blank | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00146A1 | OL-MW-3-0 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00146A2 | OL-MW-6-0 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00146A2 Dup | OL-MW-6-0 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00146A3 | OL-MW-6-1 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| BKS00146A4 | OL-MW-8-0 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00146A5 | OL-MW-7-0 | 25 | 1 | <0.02 | 0.08 | 80.0% | |
| BKS00146A6 | OL-MW-3-2 | 25 | 1 | <0.02 | 0.07 | 70.0% | |
| | | | RPD | 0% | | 13% | |



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| | | |
|---|-----------|-------------------------|
| CASE FILE NUMBER: | BKS001-46 | PAGE 1 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/29/10 | DATE RECEIVED: 12/30/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

CASE NARRATIVE

Five water samples were received by the laboratory in good condition. Analysis was performed according to the chain of custody received with the samples. Dissolved arsenic analysis was performed by ICP/MS by concentrating the samples 10 fold. No difficulties were encountered in the preparation or analysis of these samples. Sample data follows while QA/QC data is contained on subsequent pages. VOC data and Vinyl Chloride QC data are included as separate reports.

SAMPLE DATA

| SAMPLE ID | AMMONIA (mg/l) | NITRATE (mg/l) | NITRITE (mg/l) | TOC (mg/l) | COD (mg/l) | VINYL CHLORIDE (ug/l) |
|-----------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| OL-MW-3-0 | <0.010 | 0.373 | <0.002 | 2.31 | <10.0 | <0.02 |
| OL-MW-6-0 | 0.011 | <0.010 | <0.002 | 2.58 | <10.0 | <0.02 |
| OL-MW-6-1 | 0.014 | <0.010 | <0.002 | 2.62 | <10.0 | <0.02 |
| OL-MW-8-0 | 0.017 | 0.424 | 0.003 | 0.607 | <10.0 | <0.02 |
| OL-MW-7-0 | <0.010 | 0.594 | <0.002 | 0.358 | <10.0 | <0.02 |



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| CASE FILE NUMBER: | BKS001-46 | PAGE 2 |
| REPORT DATE: | 01/13/11 | |
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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
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SAMPLE DATA-CONTINUED

| SAMPLE ID | TOTAL COLIFORM (#/100mls) | CHLORIDE (mg/l) | SULFATE (mg/l) |
|-----------|------------------------------|--------------------|-------------------|
| OL-MW-3-0 | <1 | 2.59 | 17.1 |
| OL-MW-6-0 | <1 | 0.98 | 9.90 |
| OL-MW-6-1 | <1 | 1.08 | 10.3 |
| OL-MW-8-0 | <1 | 1.17 | 3.02 |
| OL-MW-7-0 | <1 | 1.37 | 3.50 |

| | pH | ALKALINITY (mgCaCO3/l) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|-----------|------|---------------------------|---------------------|--------------------|
| OL-MW-3-0 | 6.16 | 178 | 69.7 | <1.00 |
| OL-MW-6-0 | 6.60 | 75.0 | 48.0 | <1.00 |
| OL-MW-6-1 | 6.57 | 75.4 | 47.0 | <1.00 |
| OL-MW-8-0 | 6.75 | 50.2 | 35.9 | <1.00 |
| OL-MW-7-0 | 6.70 | 51.8 | 35.8 | <1.00 |



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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-46 | PAGE 3 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/29/10 | 12/30/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

SAMPLE DATA-CONTINUED

| SAMPLE ID | DISSOLVED METALS | | | | |
|-----------|-------------------|------------------|----------------|---------------------|----------------|
| | ARSENIC (ug/l) | BARIUM (mg/l) | IRON (mg/l) | MANGANESE (mg/l) | ZINC (mg/l) |
| OL-MW-3-0 | 0.107 | 0.009 | <0.020 | 1.10 | <0.005 |
| OL-MW-6-0 | 0.983 | <0.005 | 0.298 | 0.713 | <0.005 |
| OL-MW-6-1 | 0.979 | <0.005 | 0.268 | 0.698 | <0.005 |
| OL-MW-8-0 | 1.87 | <0.005 | 0.180 | 0.631 | <0.005 |
| OL-MW-7-0 | 0.318 | <0.005 | <0.020 | <0.005 | <0.005 |

| SAMPLE ID | METALS | | |
|-----------|-------------------|---------------------|------------------|
| | CALCIUM (mg/l) | POTASSIUM (mg/l) | SODIUM (mg/l) |
| OL-MW-3-0 | 35.9 | 0.907 | 8.21 |
| OL-MW-6-0 | 12.3 | 0.991 | 6.51 |
| OL-MW-6-1 | 12.0 | 0.973 | 6.38 |
| OL-MW-8-0 | 9.64 | 0.677 | 4.36 |
| OL-MW-7-0 | 9.30 | 0.806 | 4.29 |



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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-46 | PAGE 4 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/29/10 | 12/30/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | AMMONIA (mg/l) | NITRATE (mg/l) | NITRITE (mg/l) | TOC (mg/l) | COD (mg/l) | VINYL CHLORIDE (ug/l) |
|-----------------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| METHOD | EPA 350.1 | EPA 353.2 | EPA 354.1 | EPA 415.1 | EPA 410.2 | EPA 8260-SIM |
| DATE ANALYZED | 12/30/10 | 12/30/10 | 12/30/10 | 01/11/11 | 01/11/11 | 01/07/11 |
| DETECTION LIMIT | 0.006 | 0.005 | 0.001 | 0.100 | 6.00 | 0.01 |
| REPORTING LIMIT | 0.010 | 0.010 | 0.002 | 0.250 | 10.0 | 0.02 |
| DUPLICATE | | | | | | |
| SAMPLE ID | OL-MW-7-0 | OL-MW-7-0 | OL-MW-7-0 | BATCH | OL-MW-7-0 | OL-MW-6-0 |
| ORIGINAL | <0.010 | 0.594 | <0.002 | 2.86 | <10.0 | <0.02 |
| DUPLICATE | <0.010 | 0.593 | <0.002 | 2.87 | <10.0 | <0.02 |
| RPD | NC | 0.21% | NC | 0.14% | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | OL-MW-7-0 | OL-MW-7-0 | OL-MW-7-0 | BATCH | OL-MW-7-0 | |
| ORIGINAL | <0.010 | 0.594 | <0.002 | 2.86 | <10.0 | |
| SPIKED SAMPLE | 0.194 | 0.796 | 0.041 | 6.87 | 54.3 | |
| SPIKE ADDED | 0.200 | 0.200 | 0.040 | 4.50 | 50.0 | |
| % RECOVERY | 96.87% | 100.64% | 102.50% | 89.11% | 108.66% | NA |
| QC CHECK | | | | | | |
| FOUND | 0.332 | 0.409 | 0.040 | 3.88 | 97.7 | 0.09 |
| TRUE | 0.324 | 0.408 | 0.040 | 4.00 | 100 | 0.10 |
| % RECOVERY | 102.42% | 100.17% | 100.00% | 96.93% | 97.69% | 90.00% |
| BLANK | | | | | | |
| | <0.010 | <0.010 | <0.002 | <0.250 | <10.0 | <0.02 |

RPD = RELATIVE PERCENT DIFFERENCE.
 NA = NOT APPLICABLE OR NOT AVAILABLE.
 NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.
 OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.



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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-46 | PAGE 5 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/29/10 | 12/30/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | TOTAL COLIFORM (#/100mls) | CHLORIDE (mg/l) | SULFATE (mg/l) |
|-----------------|------------------------------|--------------------|-------------------|
| METHOD | SM18 9222B | EPA 325.3 | EPA 375.4 |
| DATE ANALYZED | 12/30/10 | 01/04/11 | 01/05/11 |
| DETECTION LIMIT | 1 | 0.16 | 0.76 |
| REPORTING LIMIT | 1 | 0.50 | 1.00 |
| DUPLICATE | | | |
| SAMPLE ID | OL-MW-7-0 | OL-MW-7-0 | OL-MW-7-0 |
| ORIGINAL | <1 | 1.37 | 3.50 |
| DUPLICATE | <1 | 1.37 | 3.42 |
| RPD | NC | 0.00% | 2.31% |
| SPIKE SAMPLE | | | |
| SAMPLE ID | | OL-MW-7-0 | OL-MW-7-0 |
| ORIGINAL | | 1.37 | 3.50 |
| SPIKED SAMPLE | | 11.0 | 13.2 |
| SPIKE ADDED | | 10.0 | 10.0 |
| % RECOVERY | NA | 96.78% | 97.43% |
| QC CHECK | | | |
| FOUND | | 28.4 | 9.76 |
| TRUE | | 30.0 | 10.0 |
| % RECOVERY | NA | 94.82% | 97.56% |
| PREP BLANK | | | |
| | < 1 | <0.50 | <1.00 |

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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-46 | PAGE 6 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/29/10 | 12/30/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | pH | ALKALINITY (mgCaCO3/l) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|-----------------|-----------|---------------------------|---------------------|--------------------|
| METHOD | EPA 150.1 | EPA 310.1 | EPA 310.1 | EPA 310.1 |
| DATE ANALYZED | 12/30/10 | 01/03/11 | 01/03/11 | 01/03/11 |
| DETECTION LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| REPORTING LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| DUPLICATE | | | | |
| SAMPLE ID | | OL-MW-7-0 | | |
| ORIGINAL | | 51.8 | | |
| DUPLICATE | | 52.2 | | |
| RPD | NA | 0.77% | NA | NA |
| SPIKE SAMPLE | | | | |
| SAMPLE ID | | | | |
| ORIGINAL | | | | |
| SPIKED SAMPLE | | | | |
| SPIKE ADDED | | | | |
| % RECOVERY | NA | NA | NA | NA |
| QC CHECK | | | | |
| FOUND | | 102 | | |
| TRUE | | 100 | | |
| % RECOVERY | NA | 102.20% | NA | NA |
| PREP BLANK | NA | NA | NA | NA |

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| | | |
|---|-----------|----------|
| CASE FILE NUMBER: | BKS001-46 | PAGE 7 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/29/10 | 12/30/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | ARSENIC (ug/l) | BARIUM (mg/l) | IRON (mg/l) | MANGANESE (mg/l) | ZINC (mg/l) |
|-----------------|-------------------|------------------|----------------|---------------------|----------------|
| METHOD | EPA 6020 | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 |
| DATE ANALYZED | 01/05/11 | 01/04/11 | 01/04/11 | 01/04/11 | 01/04/11 |
| DETECTION LIMIT | 0.010 | 0.002 | 0.008 | 0.003 | 0.003 |
| REPORTING LIMIT | 0.050 | 0.005 | 0.020 | 0.005 | 0.005 |
| DUPLICATE | | | | | |
| SAMPLE ID | OL-MW-7-0 | BATCH | BATCH | BATCH | BATCH |
| ORIGINAL | 0.318 | 0.006 | 0.091 | 0.007 | 0.006 |
| DUPLICATE | 0.365 | 0.006 | 0.092 | 0.007 | 0.006 |
| RPD | 13.76% | 1.77% | 1.42% | 2.86% | 3.17% |
| SPIKE SAMPLE | | | | | |
| SAMPLE ID | OL-MW-7-0 | BATCH | BATCH | BATCH | BATCH |
| ORIGINAL | 0.318 | 0.006 | 0.091 | 0.007 | 0.006 |
| SPIKED SAMPLE | 0.522 | 1.01 | 9.86 | 0.972 | 1.03 |
| SPIKE ADDED | 0.200 | 1.00 | 10.0 | 1.00 | 1.00 |
| % RECOVERY | 102.00% | 100.71% | 97.64% | 96.45% | 102.73% |
| QC CHECK | | | | | |
| FOUND | 0.721 | 1.02 | 1.05 | 1.01 | 1.02 |
| TRUE | 0.668 | 1.00 | 1.00 | 1.00 | 1.00 |
| % RECOVERY | 108.00% | 101.96% | 104.91% | 101.44% | 102.04% |
| PREP BLANK | <0.050 | <0.005 | <0.020 | <0.005 | <0.005 |

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| | | |
|--|------------------|--------------------------------|
| CASE FILE NUMBER: | BKS001-46 | PAGE 8 |
| REPORT DATE: | 01/13/11 | |
| DATE SAMPLED: | 12/29/10 | DATE RECEIVED: 12/30/10 |
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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | CALCIUM (mg/l) | POTASSIUM (mg/l) | SODIUM (mg/l) |
|-----------------|-------------------|---------------------|------------------|
| METHOD | EPA 200.7 | EPA 200.7 | EPA 200.7 |
| DATE ANALYZED | 01/04/11 | 01/04/11 | 01/04/11 |
| DETECTION LIMIT | 0.050 | 0.300 | 0.300 |
| REPORTING LIMIT | 0.100 | 0.500 | 0.500 |
| DUPLICATE | | | |
| SAMPLE ID | BATCH | BATCH | BATCH |
| ORIGINAL | 2.55 | <0.500 | 1.71 |
| DUPLICATE | 2.54 | <0.500 | 1.71 |
| RPD | 0.45% | NC | 0.42% |
| SPIKE SAMPLE | | | |
| SAMPLE ID | BATCH | BATCH | BATCH |
| ORIGINAL | 2.55 | <0.500 | 1.71 |
| SPIKED SAMPLE | 12.3 | 11.9 | 12.6 |
| SPIKE ADDED | 10.0 | 10.0 | 10.0 |
| % RECOVERY | 97.47% | 119.47% | 109.16% |
| QC CHECK | | | |
| FOUND | 10.1 | 9.53 | 9.67 |
| TRUE | 10.0 | 10.0 | 10.0 |
| % RECOVERY | 101.20% | 95.28% | 96.71% |
| PREP BLANK | | | |
| | <0.100 | <0.500 | <0.500 |

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Submitted By:

Steven Lazoff
 Laboratory Director



Aquatic Research Inc.
3927 Aurora Ave. N. , Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00143A1 | Matrix: | Water |
| Sample ID No.: | MW-1 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\1011C0601006.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-1 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRihalomethanes (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00143A1 | Matrix: | Water |
| Sample ID No.: | MW-1 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\1011C\0601006.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 74% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 95% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 92% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|------------------------|-------------------|
| Case File Number: | BKS00143A2 | Matrix: | Water |
| Sample ID No.: | MW-2 | Sample Wt/Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\10110 0701007.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------------------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| Surrogate Recoveries | | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 70% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 97% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 87% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



Aquatic Research Inc.
3927 Aurora Ave. N., Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00143A5 | Matrix: | Water |
| Sample ID No.: | MW-3 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109 1001010.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIALOMETHANES (THM) | | | |
|----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00143A5 | Matrix: | Water |
| Sample ID No.: | MW-3 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109 1001010.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 68% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 91% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 85% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00143A3 | Matrix: | Water |
| Sample ID No.: | MW-4 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109\0801008.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-1 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-4 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.4 | U | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIALOMETHANES (THM) | | | |
|----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00143A3 | Matrix: | Water |
| Sample ID No.: | MW-4 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109\0801008.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | | QC limits | | | |
| | | | | Water | Soil | | |
| | Dibromofluoromethane | 70% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 93% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 86% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00143A4 | Matrix: | Water |
| Sample ID No.: | MW-5A | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109\0901009.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 74% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 94% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 92% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00144A2 | Matrix: | Water |
| Sample ID No.: | MW-6 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/29/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109 1301013.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | | Water | Soil | | |
| | Dibromofluoromethane | 73% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 108% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 88% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00144A3 | Matrix: | Water |
| Sample ID No.: | MW-7 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/29/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/10/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101110\ 0601007.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | | Water | Soil | | |
| | Dibromofluoromethane | 78% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 96% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 105% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



Aquatic Research Inc.
3927 Aurora Ave. N., Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00144A4 | Matrix: | Water |
| Sample ID No.: | MW-8 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/29/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/10/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\10111C\0801009.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|------------------------------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | BTEX | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | 71-43-2 | Benzene | 0.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | 108-88-3 | Toluene | 0.4 | U |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | 100-41-4 | Ethylbenzene | 0.4 | U |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | p/m-Xylene | 0.4 | U |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | 95-47-6 | o-Xylene | 0.4 | U |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | TRihalOMETHANES (THM) | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | 67-66-3 | Chloroform | 0.4 | U |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | 75-27-4 | Bromodichloromethane | 0.4 | U |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | 124-48-1 | Dibromochloromethane | 0.4 | U |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | 75-25-2 | Bromoform | 0.4 | U |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.2 | J | | | | |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.4 | U | KETONES, CS2, 2-CEVE | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | 67-64-1 | Acetone | 0.4 | U |
| 98-82-8 | Isopropylbenzene | 0.4 | U | 78-93-3 | 2-Butanone | 0.4 | U |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | 591-78-6 | 2-Hexanone | 0.4 | U |
| 75-09-2 | Methylene chloride | 0.4 | U | 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | 75-15-0 | Carbon Disulfide | 0.4 | U |
| 100-42-5 | Styrene | 0.4 | U | 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00144A4 | Matrix: | Water |
| Sample ID No.: | MW-8 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/29/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/10/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\10111C\0801009.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 78% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 95% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 108% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00143A6 | Matrix: | Water |
| Sample ID No.: | MW-9 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/28/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109\ 1101011.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|---------|--------------|----------------------|--------------|------|
| | Surrogate Recoveries | %Rec. | | QC limits | | | |
| | | | | Water | Soil | | |
| | Dibromofluoromethane | 63% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 91% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 82% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | BKS00144A1 | Matrix: | Water |
| Sample ID No.: | MW-10 | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/29/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109\1201012.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 69% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 94% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 87% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|---|-------------------------|-------------------|
| Case File Number: | BKS00144A2 Dup | Matrix: | Water |
| Sample ID No.: | MW-6 Duplicate | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/29/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\1011091401014.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------------------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| Surrogate Recoveries | | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 72% | 66-118% | 66-118% | | | |
| | Toluene-d8 | 102% | 51-143% | 51-143% | | | |
| | 4-Bromofluorobenzene | 86% | 46-103% | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



Aquatic Research Inc.
3927 Aurora Ave. N. , Seattle, WA 98103 | (206) 632-2715

VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | BKS00144A3 MS | Matrix: | Water |
| Sample ID No.: | BKS00144A3 MS | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | 10/29/10 | Dilution Factor: | 1 |
| Date Received: | 10/29/10 | | |
| Date Analyzed: | 11/10/10 | Analyst: | T. Meadows |
| Date of Report: | 11/10/10 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101110\ | 0701008.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 81% | 66% | 118% |
| Toluene-d8 | 97% | 51% | 143% |
| 4-Bromofluorobenzene | 104% | 46% | 103% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 89% | 75% | 125% |
| Benzene | 99% | 75% | 125% |
| Trichloroethene | 118% | 75% | 125% |
| Toluene | 118% | 75% | 125% |
| Chlorobenzene | 118% | 75% | 125% |



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Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | 11/10/10-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 11/10/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\1011 0401005.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-3 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-4 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 1.4 | | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRihalomethanes (THM) | | | |
|-----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | 11/10/10-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 11/10/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\1011 0401005.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------------|-------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | Water | Soil | | | |
| | Dibromofluoromethane | 112% | 66-11 | 66-118% | | | |
| | Toluene-d8 | 97% | 51-14 | 51-143% | | | |
| | 4-Bromofluorobenzene | 106% | 46-10 | 46-103% | | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|------------|
| Case File Number: | 11/9/10-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\1011\0401004.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|------------|-----------------------------|--------------|------|----------|---------------------------|--------------|------|
| 108-86-1 | Bromobenzene | 0.4 | U | 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.4 | U |
| 74-97-5 | Bromochloromethane | 0.4 | U | 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.4 | U |
| 74-83-9 | Bromomethane | 0.4 | U | 127-18-4 | Tetrachloroethene | 0.4 | U |
| 104-51-8 | n-Butylbenzene | 0.4 | U | 87-61-6 | 1,2,3-Trichlorobenzene | 0.4 | U |
| 135-98-8 | sec-Butylbenzene | 0.4 | U | 120-82-1 | 1,2,4-Trichlorobenzene | 0.4 | U |
| 98-06-6 | tert-Butylbenzene | 0.4 | U | 71-55-6 | 1,1,1-Trichloroethane | 0.4 | U |
| 56-23-5 | Carbon tetrachloride | 0.4 | U | 79-00-5 | 1,1,2-Trichloroethane | 0.4 | U |
| 108-90-7 | Chlorobenzene | 0.4 | U | 79-01-6 | Trichloroethene | 0.4 | U |
| 75-00-3 | Chloroethane | 0.4 | U | 75-69-4 | Trichlorofluoromethane | 0.4 | U |
| 74-87-3 | Chloromethane | 0.4 | U | 96-18-4 | 1,2,3-Trichloropropane | 0.4 | U |
| 95-49-8 | 2-Chlorotoluene | 0.4 | U | 95-63-6 | 1,2,4-Trimethylbenzene | 0.4 | U |
| 106-43-4 | 4-Chlorotoluene | 0.4 | U | 108-67-8 | 1,3,5-Trimethylbenzene | 0.4 | U |
| 96-12-8 | 1,2-Dibromo-3-Chloropropane | 0.4 | U | 75-01-4 | Vinyl chloride | 0.4 | U |
| 106-93-4 | 1,2-Dibromoethane | 0.4 | U | | | | |
| 74-95-3 | Dibromomethane | 0.4 | U | | | | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.4 | U | | | | |
| 541-73-1 | 1,3-Dichlorobenzene | 0.4 | U | | | | |
| 106-46-7 | 1,4-Dichlorobenzene | 0.4 | U | | | | |
| 75-71-8 | Dichlorodifluoromethane | 0.4 | U | | | | |
| 75-34-3 | 1,1-Dichloroethane | 0.4 | U | | | | |
| 107-06-2 | 1,2-Dichloroethane | 0.4 | U | | | | |
| 75-35-4 | 1,1-Dichloroethene | 0.4 | U | | | | |
| 156-59-4 | cis-1,2-Dichloroethene | 0.4 | U | | | | |
| 156-60-5 | trans-1,2-Dichloroethene | 0.4 | U | | | | |
| 78-87-5 | 1,2-Dichloropropane | 0.4 | U | | | | |
| 142-28-9 | 1,3-Dichloropropane | 0.4 | U | | | | |
| 590-20-7 | 2,2-Dichloropropane | 0.4 | U | | | | |
| 563-58-6 | 1,1-Dichloropropene | 0.4 | U | | | | |
| 10061-01-5 | cis-1,3-Dichloropropene | 0.4 | U | | | | |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.4 | U | | | | |
| 87-68-3 | Hexachlorobutadiene | 0.4 | U | | | | |
| 98-82-8 | Isopropylbenzene | 0.4 | U | | | | |
| 99-87-6 | 4-Isopropyltoluene | 0.4 | U | | | | |
| 75-09-2 | Methylene chloride | 0.6 | | | | | |
| 91-20-3 | Naphthalene | 0.4 | U | | | | |
| 103-65-1 | n-Propylbenzene | 0.4 | U | | | | |
| 100-42-5 | Styrene | 0.4 | U | | | | |

| BTEX | | | |
|----------|--------------|-----|---|
| 71-43-2 | Benzene | 0.4 | U |
| 108-88-3 | Toluene | 0.4 | U |
| 100-41-4 | Ethylbenzene | 0.4 | U |
| | p/m-Xylene | 0.4 | U |
| 95-47-6 | o-Xylene | 0.4 | U |

| TRIALOMETHANES (THM) | | | |
|----------------------|----------------------|-----|---|
| 67-66-3 | Chloroform | 0.4 | U |
| 75-27-4 | Bromodichloromethane | 0.4 | U |
| 124-48-1 | Dibromochloromethane | 0.4 | U |
| 75-25-2 | Bromoform | 0.4 | U |

| KETONES, CS2, 2-CEVE | | | |
|----------------------|-------------------------|-----|---|
| 67-64-1 | Acetone | 0.4 | U |
| 78-93-3 | 2-Butanone | 0.4 | U |
| 591-78-6 | 2-Hexanone | 0.4 | U |
| 108-10-1 | 4-Methyl-2-Pentanone | 0.4 | U |
| | | | |
| 75-15-0 | Carbon Disulfide | 0.4 | U |
| 110-75-8 | 2-Chloroethylvinylether | 0.4 | U |



VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|--|-------------------------|-------------------|
| Case File Number: | 11/9/10-MB | Matrix: | Water |
| Sample ID No.: | Method Blank | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed: | 11/ 9/10 | Analyst: | T. Meadows |
| Date of Report: | 11/10/10 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\1011\0401004.D | | |

| CAS# | Name of Compound | Amount (ppb) | Flag | CAS# | Name of Compound | Amount (ppb) | Flag |
|----------|-----------------------------|--------------|--------|--------------|----------------------|--------------|------|
| | | | | QC limits | | | |
| | Surrogate Recoveries | %Rec. | | Water | Soil | | |
| | Dibromofluoromethane | 93% | 66-118 | 66-118 | 66-118% | | |
| | Toluene-d8 | 93% | 51-143 | 51-143 | 51-143% | | |
| | 4-Bromofluorobenzene | 92% | 46-103 | 46-103 | 46-103% | | |
| 107-13-1 | Acrylonitrile | 0.4 | U | 126-98-7 | Methacrylonitrile | 0.4 | U |
| 107-05-1 | Allyl Chloride | 0.4 | U | 96-33-3 | Methyl Acrylate | 0.4 | U |
| 109-69-3 | 1-Chlorobutane | 0.4 | U | 1634-04-4 | Methyl-t-Butyl Ether | 0.4 | U |
| 110-57-6 | trans-1,4-Dichloro-2-Butene | 0.4 | U | 80-62-6 | Methyl methacrylate | 0.4 | U |
| 60-29-7 | Diethyl Ether | 0.4 | U | 98-95-3 | Nitrobenzene | 0.4 | U |
| 97-63-2 | Ethyl Methacrylate | 0.4 | U | 79-46-9 | 2-Nitropropane | 0.4 | U |
| 67-72-1 | Hexachloroethane | 0.4 | U | 76-01-7 | Pentachloroethane | 0.4 | U |
| 74-88-4 | Iodomethane | 0.4 | U | | | | |

FLAGS:

- U Indicates compound was analyzed for, but not detected at the specified detection limit.
- B Blank contaminated with this analyte.
- J Estimated value - compound positively identified, but below specified detection limit.
- E Estimated value - compound exceeded calibration range.
- D Compound analyzed at a secondary dilution factor of _____ from data file: _____
- PP Compound Purges Poorly, requiring elevated detection limit.

NOTE: ppb Amounts are in $\mu\text{g/L}$ or $\mu\text{g/KG}$ dry weight.

Tentatively Identified Compounds

| | | | |
|-------------|-------------|-----------------|-------------------|
| <u>R.T.</u> | <u>CAS#</u> | <u>Compound</u> | <u>Rel. Conc.</u> |
|-------------|-------------|-----------------|-------------------|



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|------------------------------|-------------------------|------------|
| Case File Number: | 11/10/10-LCS | Matrix: | Water |
| Sample ID No.: | 11/10/10-LCS | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | | |
| Date Analyzed: | 11/10/10 | Analyst: | T. Meadows |
| Date of Report: | 11/10/10 | Supervisor's Initials: | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101110\ | 0501006.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 116% | 66% | 118% |
| Toluene-d8 | 99% | 51% | 143% |
| 4-Bromofluorobenzene | 110% | 46% | 103% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 87% | 75% | 125% |
| Benzene | 113% | 75% | 125% |
| Trichloroethene | 110% | 75% | 125% |
| Toluene | 98% | 75% | 125% |
| Chlorobenzene | 97% | 75% | 125% |



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VOLATILE ORGANIC CHEMICAL REPORT

Results of Analysis by EPA Method 8260
Measurement of Purgeable Organic Compounds by Capillary Column
Gas Chromatography/Mass Spectrometry

| | | | |
|-------------------|-------------------------------------|-------------------------|-------------------|
| Case File Number: | 11/9/10-LCS | Matrix: | Water |
| Sample ID No.: | 11/9/10-LCS | Sample Wt./Vol. (gm/ml) | 25.0 |
| Date Collected: | n/a | Dilution Factor: | 1 |
| Date Received: | n/a | Analyst: | T. Meadows |
| Date Analyzed: | 11/ 9/10 | Supervisor's Initials: | |
| Date of Report: | 11/10/10 | | |
| Data File Path: | C:\HPCHEM\1\DATA\VOA\101109\ | 0501005.D | |

| Surrogate Recoveries | %Rec. | QC limits | |
|----------------------|-------|-----------|------|
| | | | |
| Dibromofluoromethane | 91% | 66% | 118% |
| Toluene-d8 | 92% | 51% | 143% |
| 4-Bromofluorobenzene | 96% | 46% | 103% |

| Spike Recoveries | %Rec. | QC limits | |
|--------------------|-------|-----------|------|
| | | | |
| 1,1-Dichloroethene | 105% | 75% | 125% |
| Benzene | 111% | 75% | 125% |
| Trichloroethene | 112% | 75% | 125% |
| Toluene | 109% | 75% | 125% |
| Chlorobenzene | 103% | 75% | 125% |



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PHONE: (206) 632-2715 FAX: (206) 632-2417

| | | |
|--|---------------------|--------------------------------|
| CASE FILE NUMBER: | BKS001-43,44 | PAGE 1 |
| REPORT DATE: | 12/03/10 | |
| DATE SAMPLED: | 10/28,29/10 | DATE RECEIVED: 10/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

CASE NARRATIVE

Ten water samples were received by the laboratory in good condition. Analysis was performed according to the chain of custody received with the samples. Dissolved arsenic analysis was performed by ICP/MS by concentrating the samples 10 fold. No difficulties were encountered in the preparation or analysis of these samples. Sample data follows while QA/QC data is contained on subsequent pages. Organic data are included as separate reports.

SAMPLE DATA

| SAMPLE ID | AMMONIA (mg/L) | NITRATE (mg/L) | NITRITE (mg/L) | TOC (mg/L) | COD (mg/L) | VINYL CHLORIDE (ug/L) |
|-----------|----------------|----------------|----------------|------------|------------|-----------------------|
| MW-1 | <0.010 | 0.126 | 0.003 | 0.338 | <10.0 | <0.02 |
| MW-2 | <0.010 | 0.347 | <0.002 | <0.250 | <10.0 | <0.02 |
| MW-4 | <0.010 | 0.470 | <0.002 | <0.250 | <10.0 | <0.02 |
| MW-5A | <0.010 | 0.341 | <0.002 | <0.250 | <10.0 | <0.02 |
| MW-3 | <0.010 | 0.033 | <0.002 | 1.81 | <10.0 | 0.03 |
| MW-9 | <0.010 | 0.033 | <0.002 | 2.11 | 12.3 | 0.03 |
| MW-10 | 0.024 | <0.010 | <0.002 | 2.83 | <10.0 | 0.06 |
| MW-6 | 0.018 | <0.010 | 0.002 | 1.93 | <10.0 | 0.04 |
| MW-7 | <0.010 | 0.615 | <0.002 | 0.436 | <10.0 | <0.02 |
| MW-8 | 0.016 | <0.010 | 0.003 | 1.42 | <10.0 | 0.16 |



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| | | |
|---|--------------|----------|
| CASE FILE NUMBER: | BKS001-43,44 | PAGE 2 |
| REPORT DATE: | 12/03/10 | |
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| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

SAMPLE DATA-CONTINUED

| SAMPLE ID | CHLORIDE (mg/L) | SULFATE (mg/L) | TDS (mg/L) | CYANIDE (mg/L) | TOTAL COLIFORM (#/100mls) | FECAL COLIFORM (#/100mls) |
|-----------|--------------------|-------------------|---------------|-------------------|------------------------------|------------------------------|
| MW-1 | 2.64 | 3.80 | 90.5 | <0.010 | 30 | <2 |
| MW-2 | 2.01 | 2.83 | 111 | <0.010 | 52 | <2 |
| MW-4 | 2.40 | 2.48 | 112 | <0.010 | 4 | <2 |
| MW-5A | 1.92 | 3.74 | 112 | <0.010 | 4 | <2 |
| MW-3 | 2.64 | 9.56 | 239 | <0.010 | <2 | <2 |
| MW-9 | 2.64 | 10.4 | 288 | <0.010 | <2 | <2 |
| MW-10 | 0.88 | 21.5 | 299 | <0.010 | <2 | <2 |
| MW-6 | 2.01 | 18.7 | 270 | <0.010 | <2 | <2 |
| MW-7 | 1.37 | 4.38 | 97.5 | <0.010 | <2 | <2 |
| MW-8 | 1.96 | 6.35 | 224 | <0.010 | 62 | <2 |

| | pH | ALKALINITY (mgCaCO3/L) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|-------|------|---------------------------|---------------------|--------------------|
| MW-1 | 6.72 | 64.4 | 45.1 | <1.00 |
| MW-2 | 7.27 | 57.5 | 51.3 | <1.00 |
| MW-4 | 7.37 | 63.5 | 57.9 | <1.00 |
| MW-5A | 6.82 | 79.4 | 59.3 | <1.00 |
| MW-3 | 6.58 | 209 | 132 | <1.00 |
| MW-9 | 6.43 | 209 | 114 | <1.00 |
| MW-10 | 6.86 | 257 | 197 | <1.00 |
| MW-6 | 6.85 | 259 | 196 | <1.00 |
| MW-7 | 7.07 | 67.6 | 56.7 | <1.00 |
| MW-8 | 6.93 | 206 | 163 | <1.00 |



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SAMPLE DATA-CONTINUED

| DISSOLVED METALS | | | | | | |
|------------------|----------------|-----------------|---------------|------------------|----------------|-----------------|
| SAMPLE ID | ARSENIC (ug/L) | ANTIMONY (mg/L) | BARIUM (mg/L) | BERYLLIUM (mg/L) | CADMIUM (mg/L) | CHROMIUM (mg/L) |
| MW-1 | 0.094 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0043 |
| MW-2 | 0.687 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0036 |
| MW-4 | 0.226 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0030 |
| MW-5A | 0.153 | <0.0020 | 0.005 | <0.0010 | <0.0010 | 0.0031 |
| MW-3 | 0.184 | <0.0020 | 0.014 | <0.0010 | <0.0010 | <0.0010 |
| MW-9 | 0.179 | <0.0020 | 0.014 | <0.0010 | <0.0010 | <0.0010 |
| MW-10 | 2.37 | <0.0020 | 0.015 | <0.0010 | <0.0010 | <0.0010 |
| MW-6 | 1.17 | <0.0020 | 0.018 | <0.0010 | <0.0010 | 0.0031 |
| MW-7 | 0.345 | <0.0020 | <0.005 | <0.0010 | <0.0010 | 0.0032 |
| MW-8 | 2.77 | <0.0020 | 0.014 | <0.0010 | <0.0010 | <0.0010 |

| DISSOLVED METALS | | | | | | |
|------------------|---------------|---------------|-------------|-----------------|---------------|-----------------|
| SAMPLE ID | COBALT (mg/L) | COPPER (mg/L) | LEAD (mg/L) | SELENIUM (mg/L) | SILVER (mg/L) | THALLIUM (mg/L) |
| MW-1 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-2 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-4 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-5A | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-3 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-9 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-10 | <0.010 | 0.0017 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-6 | 0.038 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-7 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| MW-8 | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |

| DISSOLVED METALS | | | |
|------------------|------------|-----------------|-------------|
| SAMPLE ID | TIN (mg/L) | VANADIUM (mg/L) | ZINC (mg/L) |
| MW-1 | <0.040 | <0.0030 | <0.0050 |
| MW-2 | <0.040 | 0.0038 | <0.0050 |
| MW-4 | <0.040 | <0.0030 | <0.0050 |
| MW-5A | <0.040 | <0.0030 | <0.0050 |
| MW-3 | <0.040 | <0.0030 | <0.0050 |
| MW-9 | <0.040 | <0.0030 | <0.0050 |
| MW-10 | <0.040 | <0.0030 | <0.0050 |
| MW-6 | <0.040 | <0.0030 | <0.0050 |
| MW-7 | <0.040 | <0.0030 | <0.0050 |
| MW-8 | <0.040 | <0.0030 | <0.0050 |



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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

SAMPLE DATA-CONTINUED

| SAMPLE ID | DISSOLVED METALS | | | | | |
|-----------|------------------|------------------|---------------|------------------|-------------|------------------|
| | CALCIUM (mg/L) | POTASSIUM (mg/L) | SODIUM (mg/L) | MAGNESIUM (mg/L) | IRON (mg/L) | MANGANESE (mg/L) |
| MW-1 | 11.5 | 0.712 | 4.50 | 5.71 | 0.025 | <0.010 |
| MW-2 | 8.27 | 1.03 | 4.53 | 6.03 | <0.010 | <0.010 |
| MW-4 | 9.33 | 0.874 | 4.72 | 6.69 | <0.010 | <0.010 |
| MW-5A | 11.6 | 0.821 | 5.03 | 8.74 | <0.010 | <0.010 |
| MW-3 | 43.0 | 0.951 | 10.9 | 14.1 | 0.023 | 1.30 |
| MW-9 | 42.1 | 0.927 | 11.2 | 13.8 | 0.025 | 1.28 |
| MW-10 | 41.0 | 1.36 | 14.8 | 21.9 | 0.037 | 5.31 |
| MW-6 | 39.3 | 1.93 | 11.7 | 23.1 | 2.15 | 0.745 |
| MW-7 | 10.4 | 0.825 | 4.77 | 7.37 | 0.016 | <0.010 |
| MW-8 | 35.4 | 1.76 | 11.0 | 16.7 | 0.215 | 2.16 |

| SAMPLE ID | TOTAL METALS | |
|-----------|----------------|---------------|
| | MERCURY (mg/L) | NICKEL (mg/L) |
| MW-1 | <0.0002 | 0.0288 |
| MW-2 | <0.0002 | <0.0050 |
| MW-4 | <0.0002 | <0.0050 |
| MW-5A | <0.0002 | <0.0050 |
| MW-3 | <0.0002 | <0.0050 |
| MW-9 | <0.0002 | <0.0050 |
| MW-10 | <0.0002 | 0.0199 |
| MW-6 | <0.0002 | <0.0050 |
| MW-7 | <0.0002 | <0.0050 |
| MW-8 | <0.0002 | 0.0087 |



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QA/QC DATA

| QC PARAMETER | AMMONIA (mg/L) | NITRATE (mg/L) | NITRITE (mg/L) | TOC (mg/L) | COD (mg/L) | VINYL CHLORIDE (ug/L) |
|-----------------|-------------------|-------------------|-------------------|---------------|---------------|--------------------------|
| METHOD | EPA 350.1 | EPA 353.2 | EPA 354.1 | EPA 415.1 | EPA 410.2 | EPA 8260-SIM |
| DATE ANALYZED | 10/29/10 | 10/29/10 | 10/29/10 | 11/12/10 | 11/08/10 | 11/08/10 |
| DETECTION LIMIT | 0.006 | 0.005 | 0.001 | 0.100 | 6.00 | 0.01 |
| REPORTING LIMIT | 0.010 | 0.010 | 0.002 | 0.250 | 10.0 | 0.02 |
| DUPLICATE | | | | | | |
| SAMPLE ID | MW-9 | MW-9 | MW-9 | MW-1 | MW-8 | MW-6 |
| ORIGINAL | <0.010 | 0.033 | <0.002 | 0.338 | <10.0 | 0.04 |
| DUPLICATE | <0.010 | 0.033 | <0.002 | <0.250 | <10.0 | 0.04 |
| RPD | NC | 0.40% | NC | NC | NC | 0.00% |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | MW-9 | MW-9 | MW-9 | MW-1 | MW-8 | |
| ORIGINAL | <0.010 | 0.033 | <0.002 | 0.338 | <10.0 | |
| SPIKED SAMPLE | 0.191 | 0.229 | 0.041 | 4.82 | 60.0 | |
| SPIKE ADDED | 0.200 | 0.200 | 0.040 | 4.50 | 50.0 | |
| % RECOVERY | 95.41% | 97.81% | 102.50% | 99.66% | 119.97% | NA |
| QC CHECK | | | | | | |
| FOUND | 0.320 | 0.408 | 0.040 | 4.09 | 98.2 | 0.09 |
| TRUE | 0.324 | 0.408 | 0.040 | 4.00 | 100 | 0.10 |
| % RECOVERY | 98.83% | 99.89% | 100.00% | 102.18% | 98.16% | 90.00% |
| BLANK | | | | | | |
| | <0.010 | <0.010 | <0.002 | <0.250 | <10.0 | <0.02 |

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 OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.



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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | CHLORIDE (mg/l) | SULFATE (mg/l) | TDS (mg/L) | CYANIDE (mg/L) | TOTAL COLIFORM (#/100mls) | FECAL COLIFORM (#/100mls) |
|-----------------|--------------------|-------------------|---------------|-------------------|------------------------------|------------------------------|
| METHOD | EPA 325.3 | EPA 375.4 | SM18 2540C | SM 4500CNE | SM18 9222B | SM18 9222D |
| DATE ANALYZED | 11/01/10 | 11/10/10 | 11/04/10 | 11/09/10 | 10/29/10 | 10/29/10 |
| DETECTION LIMIT | 0.16 | 0.76 | 5.0 | 0.010 | 2 | 2 |
| REPORTING LIMIT | 0.50 | 1.00 | 5.0 | 0.010 | 2 | 2 |
| DUPLICATE | | | | | | |
| SAMPLE ID | BATCH | MW-8 | MW-8 | MW-9 | MW-9 | MW-9 |
| ORIGINAL | 0.98 | 6.35 | 224 | <0.010 | <2 | <2 |
| DUPLICATE | 0.98 | 6.43 | 240 | <0.010 | <2 | <2 |
| RPD | 0.00% | 1.20% | 6.91% | NC | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | BATCH | MW-8 | | | | |
| ORIGINAL | 0.98 | 6.35 | | | | |
| SPIKED SAMPLE | 10.9 | 16.3 | | | | |
| SPIKE ADDED | 10.0 | 10.0 | | | | |
| % RECOVERY | 98.73% | 99.01% | NA | NA | NA | NA |
| QC CHECK | | | | | | |
| FOUND | 28.4 | 10.2 | | 0.182 | | |
| TRUE | 30.0 | 10.0 | | 0.200 | | |
| % RECOVERY | 94.82% | 102.18% | NA | 90.80% | NA | NA |
| PREP BLANK | | | | | | |
| | <0.50 | <1.00 | <5.0 | <0.010 | < 1 | < 1 |

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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA

| QC PARAMETER | pH | ALKALINITY (mgCaCO3/l) | HCO3 (mgCaCO3/l) | CO3 (mgCaCO3/l) |
|-----------------|-----------|---------------------------|---------------------|--------------------|
| METHOD | EPA 150.1 | EPA 310.1 | EPA 310.1 | EPA 310.1 |
| DATE ANALYZED | 10/29/10 | 11/11/10 | 11/11/10 | 11/11/10 |
| DETECTION LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| REPORTING LIMIT | 0.10 | 1.00 | 1.00 | 1.00 |
| DUPLICATE | | | | |
| SAMPLE ID | | MW-8 | | |
| ORIGINAL | | 206 | | |
| DUPLICATE | | 196 | | |
| RPD | NA | 5.10% | NA | NA |
| SPIKE SAMPLE | | | | |
| SAMPLE ID | | | | |
| ORIGINAL | | | | |
| SPIKED SAMPLE | | | | |
| SPIKE ADDED | | | | |
| % RECOVERY | NA | NA | NA | NA |
| QC CHECK | | | | |
| FOUND | | 99.9 | | |
| TRUE | | 100 | | |
| % RECOVERY | NA | 99.90% | NA | NA |
| PREP BLANK | | | | |
| | NA | NA | NA | NA |

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| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | ARSENIC (ug/l) | ANTIMONY (mg/L) | BARIUM (mg/L) | BERYLLIUM (mg/L) | CADMIUM (mg/L) | CHROMIUM (mg/L) |
|-----------------|-------------------|--------------------|------------------|---------------------|-------------------|--------------------|
| METHOD | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 12/02/10 | 11/12/10 | 11/12/10 | 11/12/10 | 11/12/10 | 11/12/10 |
| DETECTION LIMIT | 0.010 | 0.0010 | 0.001 | 0.0003 | 0.0002 | 0.0005 |
| REPORTING LIMIT | 0.050 | 0.0020 | 0.005 | 0.0010 | 0.0010 | 0.0010 |
| DUPLICATE | | | | | | |
| SAMPLE ID | MW-1 | MW-8 | MW-8 | MW-8 | MW-8 | MW-8 |
| ORIGINAL | 0.094 | <0.0020 | 0.014 | <0.0010 | <0.0010 | <0.0010 |
| DUPLICATE | 0.094 | <0.0020 | 0.014 | <0.0010 | <0.0010 | <0.0010 |
| RPD | 0.21% | NC | 0.43% | NC | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | MW-1 | MW-8 | MW-8 | MW-8 | MW-8 | MW-8 |
| ORIGINAL | 0.094 | <0.0020 | 0.0138 | <0.0010 | <0.0010 | <0.0010 |
| SPIKED SAMPLE | 0.268 | 0.0258 | 0.065 | 0.0548 | 0.0535 | 0.0523 |
| SPIKE ADDED | 0.200 | 0.0250 | 0.050 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 86.80% | 103.20% | 102.56% | 109.56% | 107.02% | 104.52% |
| QC CHECK | | | | | | |
| FOUND | 0.668 | 0.0474 | 0.048 | 0.0549 | 0.0483 | 0.0502 |
| TRUE | 0.668 | 0.0500 | 0.050 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 100.00% | 94.74% | 96.48% | 109.74% | 96.68% | 100.40% |
| PREP BLANK | | | | | | |
| | <0.050 | <0.0020 | <0.010 | <0.0010 | <0.0010 | <0.0010 |

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QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | COBALT (mg/L) | COPPER (mg/L) | LEAD (mg/L) | SELENIUM (mg/L) | SILVER (mg/L) | THALLIUM (mg/L) |
|-----------------|------------------|------------------|----------------|--------------------|------------------|--------------------|
| METHOD | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 11/12/10 | 11/12/10 | 11/12/10 | 11/12/10 | 11/12/10 | 11/12/10 |
| DETECTION LIMIT | 0.001 | 0.0004 | 0.0003 | 0.0010 | 0.0004 | 0.0010 |
| REPORTING LIMIT | 0.010 | 0.0010 | 0.0010 | 0.0050 | 0.0010 | 0.0020 |
| DUPLICATE | | | | | | |
| SAMPLE ID | MW-8 | MW-8 | MW-8 | MW-8 | MW-8 | MW-8 |
| ORIGINAL | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| DUPLICATE | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| RPD | NC | NC | NC | NC | NC | NC |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | MW-8 | MW-8 | MW-8 | MW-8 | MW-8 | MW-8 |
| ORIGINAL | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |
| SPIKED SAMPLE | 0.054 | 0.0508 | 0.0506 | 0.0583 | 0.0254 | 0.0511 |
| SPIKE ADDED | 0.050 | 0.0500 | 0.0500 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 107.94% | 101.50% | 101.18% | 116.58% | 50.80% | 102.10% |
| QC CHECK | | | | | | |
| FOUND | 0.050 | 0.0496 | 0.0476 | 0.0488 | 0.0507 | 0.0462 |
| TRUE | 0.050 | 0.0500 | 0.0500 | 0.0500 | 0.0500 | 0.0500 |
| % RECOVERY | 99.14% | 99.20% | 95.20% | 97.52% | 101.30% | 92.38% |
| PREP BLANK | <0.010 | <0.0010 | <0.0010 | <0.0050 | <0.0010 | <0.0020 |

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QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | TIN (mg/L) | VANADIUM (mg/L) | ZINC (mg/L) |
|-----------------|---------------|--------------------|----------------|
| METHOD | EPA 6010 | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 11/07/10 | 11/12/10 | 11/12/10 |
| DETECTION LIMIT | 0.010 | 0.0010 | 0.0010 |
| REPORTING LIMIT | 0.040 | 0.0030 | 0.0050 |
| DUPLICATE | | | |
| SAMPLE ID | MW-8 | MW-8 | MW-8 |
| ORIGINAL | <0.040 | <0.0030 | <0.0050 |
| DUPLICATE | <0.040 | <0.0030 | <0.0050 |
| RPD | NC | NC | NC |
| SPIKE SAMPLE | | | |
| SAMPLE ID | MW-8 | MW-8 | MW-8 |
| ORIGINAL | <0.040 | <0.0030 | <0.0050 |
| SPIKED SAMPLE | 1.00 | 0.0526 | 0.0532 |
| SPIKE ADDED | 1.00 | 0.0500 | 0.0500 |
| % RECOVERY | 100.48% | 105.20% | 106.30% |
| QC CHECK | | | |
| FOUND | 0.942 | 0.0504 | 0.0482 |
| TRUE | 1.00 | 0.0500 | 0.0500 |
| % RECOVERY | 94.15% | 100.78% | 96.34% |
| PREP BLANK | <0.040 | <0.0030 | <0.0050 |

RPD = RELATIVE PERCENT DIFFERENCE.
 NA = NOT APPLICABLE OR NOT AVAILABLE.
 NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.
 OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.



AQUATIC RESEARCH INCORPORATED
LABORATORY & CONSULTING SERVICES
 3927 AURORA AVENUE NORTH, SEATTLE, WA 98103
 PHONE: (206) 632-2715 FAX: (206) 632-2417

| | | |
|--|--------------|--------------------------------|
| CASE FILE NUMBER: | BKS001-43,44 | PAGE 11 |
| REPORT DATE: | 12/03/10 | |
| DATE SAMPLED: | 10/28,29/10 | DATE RECEIVED: 10/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - DISSOLVED METALS

| QC PARAMETER | CALCIUM (mg/L) | POTASSIUM (mg/L) | SODIUM (mg/L) | MAGNESIUM (mg/L) | IRON (mg/L) | MANGANESE (mg/L) |
|-----------------|-------------------|---------------------|------------------|---------------------|----------------|---------------------|
| METHOD | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 | EPA 6010 |
| DATE ANALYZED | 11/16/10 | 11/07/10 | 11/16/10 | 11/07/10 | 11/17/10 | 11/07/10 |
| DETECTION LIMIT | 0.050 | 0.300 | 0.300 | 0.050 | 0.010 | 0.003 |
| REPORTING LIMIT | 0.100 | 0.500 | 0.500 | 0.100 | 0.020 | 0.010 |
| DUPLICATE | | | | | | |
| SAMPLE ID | MW-8 | MW-8 | MW-8 | MW-8 | MW-4 | MW-8 |
| ORIGINAL | 35.4 | 1.76 | 11.0 | 16.7 | <0.010 | 2.16 |
| DUPLICATE | 35.2 | 1.83 | 11.2 | 17.5 | 0.011 | 2.30 |
| RPD | 0.46% | 3.99% | 2.06% | 4.83% | NC | 4.97% |
| SPIKE SAMPLE | | | | | | |
| SAMPLE ID | MW-8 | MW-8 | MW-8 | MW-8 | MW-4 | MW-8 |
| ORIGINAL | 35.4 | 1.76 | 11.0 | 16.7 | <0.010 | 2.16 |
| SPIKED SAMPLE | 83.2 | 12.9 | 59.5 | 25.8 | 0.975 | 3.22 |
| SPIKE ADDED | 50.0 | 10.0 | 50.0 | 10.0 | 1.00 | 1.00 |
| % RECOVERY | 95.71% | 111.61% | 97.12% | 91.44% | 97.52% | 106.58% |
| QC CHECK | | | | | | |
| FOUND | 10.2 | 10.1 | 10.3 | 0.979 | 9.25 | 0.961 |
| TRUE | 10.0 | 10.0 | 10.0 | 1.00 | 10.0 | 1.00 |
| % RECOVERY | 102.38% | 100.99% | 103.37% | 97.90% | 92.51% | 96.05% |
| PREP BLANK | | | | | | |
| | <0.100 | <0.500 | <0.500 | <0.100 | <0.500 | <0.500 |

RPD = RELATIVE PERCENT DIFFERENCE.
 NA = NOT APPLICABLE OR NOT AVAILABLE.
 NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.
 OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.



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PHONE: (206) 632-2715 FAX: (206) 632-2417

| | | |
|--|---------------------|--------------------------------|
| CASE FILE NUMBER: | BKS001-43,44 | PAGE 12 |
| REPORT DATE: | 12/03/10 | |
| DATE SAMPLED: | 10/28,29/10 | DATE RECEIVED: 10/29/10 |
| FINAL REPORT, LABORATORY ANALYSIS OF SELECTED PARAMETERS ON WATER | | |
| SAMPLES FROM KITSAP COUNTY SOLID WASTE / OLALLA LANDFILL | | |

QA/QC DATA - TOTAL METALS

| QC PARAMETER | MERCURY (mg/l) | NICKEL (mg/l) |
|-----------------|-------------------|------------------|
| METHOD | EPA 6020 | EPA 6020 |
| DATE ANALYZED | 11/12/10 | 11/12/10 |
| DETECTION LIMIT | 0.0001 | 0.0010 |
| REPORTING LIMIT | 0.0002 | 0.0050 |
| DUPLICATE | | |
| SAMPLE ID | MW-8 | MW-8 |
| ORIGINAL | <0.0002 | 0.0087 |
| DUPLICATE | <0.0002 | 0.0088 |
| RPD | NC | 0.91% |
| SPIKE SAMPLE | | |
| SAMPLE ID | MW-8 | MW-8 |
| ORIGINAL | <0.0002 | 0.0087 |
| SPIKED SAMPLE | 0.0023 | 0.0600 |
| SPIKE ADDED | 0.0020 | 0.0500 |
| % RECOVERY | 113.40% | 102.53% |
| QC CHECK | | |
| FOUND | 0.0019 | 0.0491 |
| TRUE | 0.0020 | 0.0500 |
| % RECOVERY | 96.40% | 98.16% |
| PREP BLANK | <0.0002 | <0.0050 |

RPD = RELATIVE PERCENT DIFFERENCE.
NA = NOT APPLICABLE OR NOT AVAILABLE.
NC = NOT CALCULABLE DUE TO ONE OR MORE VALUES BEING BELOW THE DETECTION LIMIT.
OR = RECOVERY NOT CALCULABLE DUE TO SPIKE SAMPLE OUT OF RANGE OR SPIKE TOO LOW RELATIVE TO SAMPLE CONCENTRATION.

Submitted By:

Steven Lazoff
Laboratory Director

Vinyl Chloride by SIM

T. Meadows

11/8/10

| Sample ID | Client ID | Sample Amount (ml) | Dilution Factor | Vinyl Chloride (µg/L) | 1,2-Dichloroethane-d4 (µg/L) | % R Surr | % R VC |
|----------------|--------------|--------------------|-----------------|-----------------------|------------------------------|----------|--------|
| CCV | CCV | 25 | 1 | 0.09 | 0.10 | 100.0% | 90.0% |
| CCV | CCV | 25 | 1 | 0.10 | 0.11 | 110.0% | 100.0% |
| 11/8/10-MB | Method Blank | 25 | 1 | <0.02 | 0.10 | 100.0% | |
| BKS00143A1 | MW-1 | 25 | 1 | <0.02 | 0.11 | 110.0% | |
| BKS00143A2 | MW-2 | 25 | 1 | <0.02 | 0.09 | 90.0% | |
| BKS00143A3 | MW-4 | 25 | 1 | <0.02 | 0.11 | 110.0% | |
| BKS00143A4 | MW-5A | 25 | 1 | <0.02 | 0.09 | 90.0% | |
| BKS00143A5 | MW-3 | 25 | 1 | 0.03 | 0.10 | 100.0% | |
| BKS00143A6 | MW-9 | 25 | 1 | 0.03 | 0.10 | 100.0% | |
| BKS00144A1 | MW-10 | 25 | 1 | 0.06 | 0.08 | 80.0% | |
| BKS00144A2 | MW-6 | 25 | 1 | 0.04 | 0.08 | 80.0% | |
| BKS00144A2 Dup | MW-6 | 25 | 1 | 0.04 | 0.09 | 90.0% | |
| BKS00144A3 | MW-7 | 25 | 1 | <0.02 | 0.09 | 90.0% | |
| BKS00144A4 | MW-8 | 25 | 1 | 0.16 | 0.11 | 110.0% | |
| | | | RPD | 0% | | 12% | |

Appendix J

Test Pit Logs and Photos

Olalla Landfill

OL-MW-##-W-0 ^{primary}

- 1 dup
- 2 trip
- 3 equip

0945 onsite

Met Keli (Means McKay)
Jan Brower, KCHD
Eric Caddey, EPI
Mike Warfel, parametrix

Glacier
Cascade
Pact
Noel Phillips
↓
Ecology
Well
construction
Coordinator

10/7/10

MW-2 water level @ 11:12 am 66.89' toC

OL-MW-02-W-0 collected @ "11:00"

MW-4 water level @ 11:59 64.09' toC

OL-MW-04-W-0 collected @ "12:00"

Sampled by

Mike & Mary
Warfel Holder

Samples taken by Mike to
Aquatic Research

with bailers
put on ice

Approximately 13:30 started digging Test Pit 1

~ 14:00

Test Pit 2

~ 14:15

Test Pit 3

~ 14:35

Test Pit 4

~ 14:50 dug out drum & additional area @ TP 4 (North end)

~ 15:20

Test Pit 5 N. Phillips offsite

~ 15:45 started backfilling

M. Warfel & J. Brower offsite

reburied all waste

K. McKay-Means offsite

16:15 M. Holder offsite only E. Caddey & Ben Bohren (Glacier)
onsite - finishing the backfilling of TP 2 & TP 1.

Multi Rae (Oil Ind Safety) used to measure PID readings.

Test Pit Designation:

1

| | |
|---|---|
| <p>Client: Logged By: <i>M. Holder</i> Date of Excavation: <i>10/7/10</i> Site Address: Excavation Contractor: <i>Glacier</i> Method:</p> | <p>Site Personnel: <i>Mary Holder - EPI</i> <i>Mike Warfel - Pavamatrix</i> <i>Keli McKay-Means - KCCo</i> <i>Jan Brower - KCHD</i> <i>Noel Phillips - Ecy</i> <i>Ben Bohren - Glacier</i></p> |
|---|---|

| Length along trench (in feet) | SUBSURFACE PROFILE | | Trash | Liquid | Depth (Feet bgs) | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|--|--------------------------------|-------|--------|------------------|--|
| | DESCRIPTION: Trash, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | | | | | |
| 1 | <i>North</i> | | | | | <p><i>no PID kits</i></p> <p><i>Mostly dirt & wood debris with sm amounts of trash</i></p> |
| 2 | <i>brick plastic</i> | | | | | |
| 3 | <i>tires wood debris</i> | | | | | |
| 4 | <i>2x4</i> | | | | | |
| 5 | | | | | | |
| 6 | | | | | | |
| 7 | | | | | | |
| 8 | | | | | | |
| 9 | <i>Rebar</i> | | | | | |
| 10 | <i>plywood</i> | | | | | |
| 11 | | | | | | |
| 12 | | | | | | |
| 13 | | | | | | |
| 14 | | | | | | |
| 15 | | | | | | |
| 16 | <i>grey dirt</i> | <i>some black dirt</i> | | | | |
| 17 | <i>roof</i> | <i>shingles</i> | | | | |
| 18 | | <i>car bumper</i> | | | | |
| 19 | | <i>~3 inch diam alum? pipe</i> | | | | |
| 20 | | | | | | |
| 21 | | | | | | |
| 22 | | | | | | |
| 23 | | <i>yellow insulation</i> | | | | |
| 24 | | | | | | |
| 25 | | <i>South</i> | | | | |

| | |
|---|--|
| <p>Client: Olalla Landfill - Kitsap County</p> <p>Logged By: Mary Holder</p> <p>Date of Excavation: 10/7/10</p> <p>Site Address:</p> <p>Excavation Contractor: Glacier</p> <p>Method:</p> | <p>Site Personnel:</p> <p>Mary Mike Keli Jan Ben Bohren - Glacier</p> |
|---|--|

| Length along trench (in feet) | SUBSURFACE PROFILE | Trash | Liquid | Depth (Feet bgs) | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|--|-------|--------|------------------|--|
| | DESCRIPTION: Trash, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | | | | |
| 1 | North | | NA | 8' | no PID hits no liquids no perched layers |
| 2 | | | | | |
| 3 | | | | | |
| 4 | grey | | | | |
| 5 | darker brown than TP 1 | | | | |
| 6 | | | | | |
| 7 | | | | | |
| 8 | | | | | |
| 9 | | | | | |
| 10 | | | | | |
| 11 | | | | | |
| 12 | | | | | |
| 13 | | | | | |
| 14 | No trash or wood debris | | | | |
| 15 | | | | | |
| 16 | layers of grey brown and vegetation (grassy) | | | | |
| 17 | | | | | |
| 18 | | | | | |
| 19 | | | | | |
| 20 | | | | | |
| 21 | | | | | |
| 22 | | | | | |
| 23 | | | | | |
| 24 | | | | | |
| 25 | South | | | | |

Client: Kitsap County Olalla Landfill
 Logged By: Mary Holder
 Date of Excavation: 10/7/10
 Site Address:
 Excavation Contractor: Glacier
 Method:

Site Personnel: Mary, Mike, Keli, Jan

started ~ 2:15
 finished ~ 2:30
 4:45pm Finishing Filling in

| Length along trench (in feet) | SUBSURFACE PROFILE | | Depth (Feet bgs) | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|--|----------------|------------------|--|
| | DESCRIPTION: Trash, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | Trash / Liquid | | |
| 1 | North | | NA 8' | |
| 2 | top soil lighter tan color than pit 1 | | | |
| 3 | | | | |
| 4 | | | | |
| 5 | | | | |
| 6 | | | | |
| 7 | | | | |
| 8 | bot copper | | | |
| 9 | brick section of bldg | | | |
| 10 | plastic bag | | | |
| 11 | " landscape planters " | | | |
| 12 | black woody debris @ bottom | | | |
| 13 | | | | |
| 14 | | | | |
| 15 | top of metal bucket | | | |
| 16 | | | | |
| 17 | | | | |
| 18 | netting or strings small boot | | | |
| 19 | wood debris @ bottom | | | |
| 20 | | | | |
| 21 | plastic onion bags | | | |
| 22 | | | | |
| 23 | | | | |
| 24 | | | | |
| 25 | South | | | |

NO PID hits
 No liquid seen perched layer

Client: Olalla Landfill Kitsap county
 Logged By: Mary Holder
 Date of Excavation: 10/7/2010
 Site Address:
 Excavation Contractor: Glacier
 Method:

Site Personnel: Mary, Mike, Keli, Jan, Noel

started @ 2:30pm
 finished @ 3:05
 started filling @ 4pm

| Length along trench (in feet) | SUBSURFACE PROFILE | | Depth (Feet bgs) | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|---|----------------|------------------|--|
| | DESCRIPTION: Trash, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | Trash / Liquid | | |
| 1 | North | | 8' | <p>★ pulled w/ barrel - 55 gal drum @ start of hole & went back another 22'</p> <p>TOTAL Trench 47'</p> <p>★ dirt 55gal drum painted red barrel tar like black subs</p> <p>bicycle tires inner tube bottle styrofoam toilet sink</p> <p>plastic bottles shoe netting possibly cooling coils from back of refrigerator or grill</p> <p>large metal pieces truck</p> <p>22' shingles metal netting 22' further North from starting point</p> <p>NO PID hits except for 2.3 ppm when held up</p> <p>* to tar like substance in 55 gal drum</p> <p>No liquid or perched layers</p> |
| 2 | | NA | | |
| 3 | | | | |
| 4 | | | | |
| 5 | | | | |
| 6 | | | | |
| 7 | | | | |
| 8 | | | | |
| 9 | | | | |
| 10 | | | | |
| 11 | | | | |
| 12 | | | | |
| 13 | | | | |
| 14 | | | | |
| 15 | | | | |
| 16 | | | | |
| 17 | | | | |
| 18 | | | | |
| 19 | | | | |
| 20 | | | | |
| 21 | | | | |
| 22 | | | | |
| 23 | | | | |
| 24 | | | | |
| 25 | South end | | | |

Client: Olalla Landfill / Kitsop County
 Logged By: Mary Holder
 Date of Excavation: 10/2/10
 Site Address:
 Excavation Contractor: Glacier started digging @ 3:30
 Method: back excavator
 Site Personnel: Mary, Mike, Kel, Jan
 digging started @ 3:10pm
 finished @ 4pm

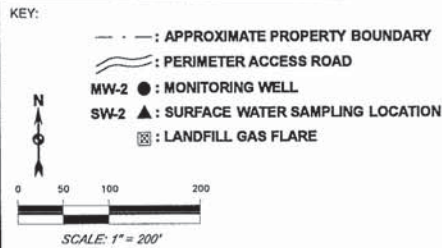
| Length along trench (in feet) | SUBSURFACE PROFILE | | Depth (Feet bgs) | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|---|--------------|------------------|---|
| | DESCRIPTION: Trash, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | Trash Liquid | | |
| 1 | North | | NA | Overall this pit seemed the cleanest of the 5 with very little trash or construction debris. Mostly soil. No liquid or perched layers. |
| 2 | | | | |
| 3 | dirt | | | |
| 4 | | | | |
| 5 | | | | |
| 6 | | | | |
| 7 | | | | |
| 8 | glass bottle | | | |
| 9 | | | | |
| 10 | | | | |
| 11 | | | | |
| 12 | small amt construction debris | | | |
| 13 | | | | |
| 14 | | | | |
| 15 | | | | |
| 16 | | | | |
| 17 | | | | |
| 18 | | | | |
| 19 | | | | |
| 20 | bottles newspaper | | | |
| 21 | plastic | | | |
| 22 | | | | |
| 23 | | | | |
| 24 | | | | |
| 25 | South | | | |

S.E. BURLEY OLALLA ROAD



Approximate Test Pit (TP) locations dug 10/7/2010. Each pit 25' long and 8' deep.

*MW-5 is completed in a shallow perched groundwater zone.



ePi ENVIRONMENTAL PARTNERS INC
 295 NE Gilman Boulevard, Suite 201
 Issaquah, Washington 98027

FIGURE 1

OLALLA LANDFILL
 MONITORING LOCATIONS
 KITSAP COUNTY, WASHINGTON

| | | | |
|------------------------|--|---------------------------|-------------------------|
| PROJECT | 45401.6 OLALLA, WASHINGTON | | |
| PREPARED FOR | KITSAP COUNTY | | |
| LOCATION | OLALLA LANDFILL KITSAP COUNTY, WASHINGTON | | |
| SHEET 1 of 1 | DRAWN BY ARM | REVIEWED BY DCK | DATE 07/06/10 |

| | | | |
|-------------------------------|---|------------------------|---|
| Client: | Kitsap County Dept. of Public Works | Site Personnel: | Mary Holder - EPI, Ben Bohren - Glacier, |
| Logged By: | Mary Holder | | Mike Warfel - Parametrix, Keli McKay-Means - KCDPW, |
| Date of Excavation: | October 7, 2010 | | Jan Brower - KCHD, Noel Phillips - WADOE |
| Site Address: | 0.75 mi E of Hwy 16 on Olalla-Burley Rd | | |
| Excavation Contractor: | Glacier Environmental Services, Inc. | | |
| Method: | Excavator | | |

| Length along trench (in feet) | SUBSURFACE PROFILE | | Liquid | Trench Depth | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|---|--|--------|--------------|--|
| | DESCRIPTION: Refuse, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | | | | |
| 1 | North End | | No | 8' bgs | No PID hits. |
| 2 | | | No | 8' bgs | No perched water layers. |
| 3 | Soil with no refuse | | No | 8' bgs | Predominately soil with very little refuse and construction debris |
| 4 | | | No | 8' bgs | (See photos IMG_0890.jpg, 0893 and 0899) |
| 5 | | | No | 8' bgs | Thin layers of light brown and grey soils |
| 6 | | | No | 8' bgs | (See photo IMG_0889.jpg and 0897) |
| 7 | | | No | 8' bgs | Walls of pit collapsed |
| 8 | Glass bottle | | No | 8' bgs | (See photo IMG_0898.jpg) |
| 9 | | | No | 8' bgs | |
| 10 | | | No | 8' bgs | |
| 11 | | | No | 8' bgs | |
| 12 | | | No | 8' bgs | |
| 13 | A couple pieces of construction debris | | No | 8' bgs | |
| 14 | | | No | 8' bgs | |
| 15 | | | No | 8' bgs | |
| 16 | | | No | 8' bgs | |
| 17 | | | No | 8' bgs | |
| 18 | | | No | 8' bgs | |
| 19 | | | No | 8' bgs | |
| 20 | | | No | 8' bgs | |
| 21 | Woody debris | | No | 8' bgs | |
| 22 | | | No | 8' bgs | |
| 23 | Newspaper | | No | 8' bgs | |
| 24 | Grouping of glass bottles and plastic bags or sheeting | | No | 8' bgs | |
| 25 | | | No | 8' bgs | |
| | South End | | | | |

| | | | |
|-------------------------------|---|------------------------|---|
| Client: | Kitsap County Dept. of Public Works | Site Personnel: | Mary Holder - EPI, Ben Bohren - Glacier, Mike Warfel - Parametrix, Keli McKay-Means - KCDPW, Jan Brower - KCHD, Noel Phillips - WADOE |
| Logged By: | Mary Holder | | |
| Date of Excavation: | October 7, 2010 | | |
| Site Address: | 0.75 mi E of Hwy 16 on Olalla-Burley Rd | | |
| Excavation Contractor: | Glacier Environmental Services, Inc. | | |
| Method: | Excavator | | |

| Length along trench (in feet) | SUBSURFACE PROFILE | | Liquid Observed? | Trench Depth | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|--|--|------------------|--------------|--|
| | DESCRIPTION: Trash, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | | | | |
| 1 | North End | | No | 8' bgs | No organic vapors were measured with PID |
| 2 | Nearly empty red 55-gallon drum containing a small quantity of tar-like material | | No | 8' bgs | No perched water layers were observed |
| 3 | (See photos IMG_0864.jpg and 0865) | | No | 8' bgs | Mostly soil with construction debris and refuse |
| 4 | | | No | 8' bgs | (See photos IMG_0859.jpg, 0861, 0862 and 0871) |
| 5 | Bicycle tire or innertube | | No | 8' bgs | Light brown and grey soil in thick layers |
| 6 | Bottle, styrofoam | | No | 8' bgs | Soils are predominately sand and gravel |
| 7 | Piece of a toilet | | No | 8' bgs | Walls of pit collapsed while digging out drum |
| 8 | | | No | 8' bgs | (See photo IMG_0866.jpg) |
| 9 | Small amounts of random refuse | | No | 8' bgs | |
| 10 | Metal, glass, foam, plastic, lumber, brick, etc. | | No | 8' bgs | |
| 11 | | | No | 8' bgs | |
| 12 | | | No | 8' bgs | |
| 13 | | | No | 8' bgs | |
| 14 | | | No | 8' bgs | |
| 15 | | | No | 8' bgs | |
| 16 | | | No | 8' bgs | |
| 17 | Automobile tire | | No | 8' bgs | |
| 18 | | | No | 8' bgs | |
| 19 | | | No | 8' bgs | |
| 20 | | | No | 8' bgs | |
| 21 | Small amount of metal debris and wood | | No | 8' bgs | |
| 22 | | | No | 8' bgs | |
| 23 | glass bottles | | No | 8' bgs | |
| 24 | | | No | 8' bgs | |
| 25 | South End | | No | 8' bgs | |

| | | | |
|-----------------------------|---|------------------------|---|
| Client: | Kitsap County Dept. of Public Works | Site Personnel: | Mary Holder - EPI, Ben Bohren - Glacier, Mike Warfel - Parametrix, Keli McKay-Means - KCDPW, Jan Brower - KCHD, Noel Phillips - WADOE |
| Logged By: | Mary Holder | | |
| Date of Excavation: | October 7, 2010 | | |
| Site Address: | 0.75 mi E of Hwy 16 on Olalla-Burley Rd | | |
| Excavation Contract: | Glacier Environmental Services, Inc. | | |
| Method: | Excavator | | |

| Length along trench (in feet) | SUBSURFACE PROFILE | | Liquid Observed? | Trench Depth | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|---|--|------------------|--------------|---|
| | DESCRIPTION: Refuse, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | | | | |
| 22 | North end of TP-4 extension | | No | 8' bgs | In order to check for the possibility of more buried drums, additional soils were excavated to the north of the red drum. |
| 21 | Heavy duty woven metal screen | | No | 8' bgs | |
| 20 | Large quantity of roofing shingles | | No | 8' bgs | No measured PID detections other than from the small volume of tar-like substance from red |
| 19 | | | No | 8' bgs | No perched water layers were observed |
| 18 | Lumber and wood debris | | No | 8' bgs | Soil, construction debris and metal objects |
| 17 | | | No | 8' bgs | (See photos IMG_0874.jpg, 0875, 0877, 0880, 0883, and 0885) |
| 16 | Grey soil, lumber and wood debris and netting | | No | 8' bgs | Soil layered light brown and grey |
| 15 | | | No | 8' bgs | (See photo IMG_0881.jpg) |
| 14 | Automobile tire | | No | 8' bgs | Soils are predominately sand and gravel |
| 13 | | | No | 8' bgs | Walls of pit collapsed |
| 12 | Large pieces of thick metal | | No | 8' bgs | (See photo IMG_0886.jpg) |
| 11 | | | No | 8' bgs | |
| 10 | | | No | 8' bgs | |
| 9 | | | No | 8' bgs | |
| 8 | Netting | | No | 8' bgs | |
| 7 | | | No | 8' bgs | |
| 6 | Metal - possibly cooling coils from back or refrigerator or grill | | No | 8' bgs | |
| 5 | | | No | 8' bgs | |
| 4 | Shoe | | No | 8' bgs | |
| 3 | Plastic bottles | | No | 8' bgs | |
| 2 | Cloth | | No | 8' bgs | |
| 1 | Drum was surrounded by soil and contained approximately one liter of a black tar like substance | | No | 8' bgs | PID reading of 2.3 ppm when tar-like substance was kneaded |

South end of additional excavation
North end of 25' test pit and location of red drum

(See photos IMG_0867.jpg through 0870)

| | | | |
|-------------------------------|---|------------------------|---|
| Client: | Kitsap County Dept. of Public Works | Site Personnel: | Mary Holder - EPI, Ben Bohren - Glacier, |
| Logged By: | Mary Holder | | Mike Warfel - Parametrix, Keli McKay-Means - KCDPW, |
| Date of Excavation: | October 7, 2010 | | Jan Brower - KCHD, Noel Phillips - WADOE |
| Site Address: | 0.75 mi E of Hwy 16 on Olalla-Burley Rd | | |
| Excavation Contractor: | Glacier Environmental Services, Inc. | | |
| Method: | Excavator | | |

| Length along trench (in feet) | SUBSURFACE PROFILE | | Liquid Observed? | Trench Depth | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|--|--|------------------|--------------|--|
| | DESCRIPTION: Refuse Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | | | | |
| 1 | North End | | No | 8' bgs | No organic vapors measured with PID |
| 2 | Soil with no refuse | | No | 8' bgs | No perched water layers were observed |
| 3 | | | No | 8' bgs | Mostly soil and construction debris |
| 4 | | | No | 8' bgs | Very little refuse observed |
| 5 | | | No | 8' bgs | (See photos IMG_0840.jpg, 0842 and 0850) |
| 6 | | | No | 8' bgs | Top 1-2' of soil tan in color with grey soil below |
| 7 | | | No | 8' bgs | Blackened soil at bottom of excavation |
| 8 | Strip of copper | | No | 8' bgs | (See photos IMG_0850.jpg and 0851) |
| 9 | Section of brick building | | No | 8' bgs | Soils are predominately sand and gravel |
| 10 | Plastic bag printed with the words "landscape planters" | | No | 8' bgs | Walls of pit started collapsing |
| 11 | Blackened lumber and woody debris toward bottom of trench | | No | 8' bgs | (See photo IMG_0855.jpg) |
| 12 | | | No | 8' bgs | |
| 13 | | | No | 8' bgs | For location see photo IMG_0852.jpg |
| 14 | | | No | 8' bgs | |
| 15 | Top of gallon sized metal bucket | | No | 8' bgs | |
| 16 | | | No | 8' bgs | |
| 17 | | | No | 8' bgs | |
| 18 | Netting or string | | No | 8' bgs | |
| 19 | Small boot | | No | 8' bgs | |
| 20 | Lumber and wood debris at bottom of excavation | | No | 8' bgs | |
| 21 | | | No | 8' bgs | |
| 22 | | | No | 8' bgs | |
| 23 | | | No | 8' bgs | |
| 24 | Plastic onion bags | | No | 8' bgs | |
| 25 | South End | | No | 8' bgs | |

| | | | |
|-------------------------------|---|------------------------|---|
| Client: | Kitsap County Dept. of Public Works | Site Personnel: | Mary Holder - EPI, Ben Bohren - Glacier, Mike Warfel - Parametrix, Keli McKay-Means - KCDPW, Jan Brower - KCHD, Noel Phillips - WADOE |
| Logged By: | Mary Holder | | |
| Date of Excavation: | October 7, 2010 | | |
| Site Address: | 0.75 mi E of Hwy 16 on Olalla-Burley Rd | | |
| Excavation Contractor: | Glacier Environmental Services, Inc. | | |
| Method: | Excavator | | |

| Length along trench (in feet) | SUBSURFACE PROFILE | | Liquid Observed? | Trench Depth | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|--|--|------------------|--------------|--|
| | DESCRIPTION: Refuse Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | | | | |
| 1 | North End | | No | 8' bgs | No organic vapors were measured with PID |
| 2 | Grey and brown soils | | No | 8' bgs | No perched water layers were observed |
| 3 | | | No | 8' bgs | No refuse was observed |
| 4 | | | No | 8' bgs | (See photos IMG_0837.jpg and 0838) |
| 5 | | | No | 8' bgs | Light brown and grey soils layered with grasses |
| 6 | | | No | 8' bgs | (See photo IMG_0835.jpg) |
| 7 | | | No | 8' bgs | |
| 8 | | | No | 8' bgs | Soils consisted predominately of |
| 9 | | | No | 8' bgs | sand and gravel with silt |
| 10 | | | No | 8' bgs | (See photo IMG_0831.jpg) |
| 11 | No trash or wood debris were observed | | No | 8' bgs | |
| 12 | Soil layered with grassy vegetation | | No | 8' bgs | |
| 13 | | | No | 8' bgs | |
| 14 | | | No | 8' bgs | |
| 15 | | | No | 8' bgs | |
| 16 | | | No | 8' bgs | |
| 17 | | | No | 8' bgs | |
| 18 | | | No | 8' bgs | |
| 19 | | | No | 8' bgs | |
| 20 | | | No | 8' bgs | |
| 21 | | | No | 8' bgs | |
| 22 | | | No | 8' bgs | |
| 23 | | | No | 8' bgs | |
| 24 | Soil with no refuse | | No | 8' bgs | |
| 25 | South End | | No | 8' bgs | |

| | | | |
|-------------------------------|---|------------------------|---|
| Client: | Kitsap County Dept. of Public Works | Site Personnel: | Mary Holder - EPI, Ben Bohren - Glacier, |
| Logged By: | Mary Holder | | Mike Warfel - Parametrix, Keli McKay-Means - KCDPW, |
| Date of Excavation: | October 7, 2010 | | Jan Brower - KCHD, Noel Phillips - WADOE |
| Site Address: | 0.75 mi E of Hwy 16 on Olalla-Burley Rd | | |
| Excavation Contractor: | Glacier Environmental Services, Inc. | | |
| Method: | Excavator | | |

| Length along trench (in feet) | SUBSURFACE PROFILE | | Liquid Observed? | Trench Depth | Comments (Including location(s) of additional excavation along side of trench) |
|-------------------------------|---|--|------------------|--------------|--|
| | DESCRIPTION: Refuse, Soil, Color, Moisture, Description. If liquid: pH, conductivity, temperature, DO, ORP | | | | |
| 1 | North End | | No | 8' bgs | No organic vapors were measured with PID. |
| 2 | Brick, plastic | | No | 8' bgs | No perched water layers were observed. |
| 3 | tire, wood debris | | No | 8' bgs | Mostly soil and wood debris with |
| 4 | 2 x 4 | | No | 8' bgs | small amounts of refuse. |
| 5 | Rebar | | No | 8' bgs | (See photos IMG_0811.jpg, 0812, 0823 and 0828) |
| 6 | Plywood | | No | 8' bgs | Soils predominately sand and gravel with |
| 7 | | | No | 8' bgs | layers of light brown soil, grey soil, and refuse |
| 8 | | | No | 8' bgs | (See photos IMG_0822.jpg, 0825 and 0826) |
| 9 | | | No | 8' bgs | |
| 10 | | | No | 8' bgs | |
| 11 | | | No | 8' bgs | |
| 12 | Wood debris | | No | 8' bgs | For location see photo IMG_0830.jpg |
| 13 | | | No | 8' bgs | |
| 14 | | | No | 8' bgs | |
| 15 | | | No | 8' bgs | |
| 16 | Grey and black soil at bottom | | No | 8' bgs | |
| 17 | Possibly burned materials | | No | 8' bgs | |
| 18 | | | No | 8' bgs | |
| 19 | | | No | 8' bgs | |
| 20 | ~ 3" diameter metal pipe | | No | 8' bgs | |
| 21 | Automobile tire | | No | 8' bgs | |
| 22 | Roof shingles | | No | 8' bgs | |
| 23 | Automobile bumper | | No | 8' bgs | |
| 24 | | | No | 8' bgs | |
| 25 | Piece of yellow insulation | | No | 8' bgs | |
| | South End | | | | |

TEST PIT NO. 5



IMG_0888.JPG



IMG_0889.JPG



IMG_0890.JPG



IMG_0891.JPG

TEST PIT NO. 5



IMG_0892.JPG



IMG_0893.JPG



IMG_0894.JPG



IMG_0895.JPG

TEST PIT NO. 5



IMG_0896.JPG



IMG_0897.JPG



IMG_0898.JPG



IMG_0899.JPG

TEST PIT NO. 5



IMG_0900.JPG



IMG_0901.JPG



IMG_0902.JPG



IMG_0903.JPG

TEST PIT NO. 4



IMG_0858.JPG



IMG_0858.JPEG

TEST PIT NO. 3



IMG_0840.JPG



IMG_0841.JPG



IMG_0842.JPG



IMG_0843.JPG

TEST PIT NO. 3



IMG_0844.JPG



IMG_0845.JPG



IMG_0846.JPG



IMG_0847.JPG

TEST PIT NO. 3



IMG_0848.JPG



IMG_0849.JPG



IMG_0850.JPG



IMG_0851.JPG

TEST PIT NO. 3



IMG_0852.JPG



IMG_0853.JPG



IMG_0854.JPG



IMG_0855.JPG

TEST PIT NO. 3



IMG_0856.JPG



IMG_0857.JPG

TEST PIT NO. 2



IMG_0831.JPG



IMG_0832.JPG



IMG_0833.JPG



IMG_0834.JPG

TEST PIT NO. 2



IMG_0835.JPG



IMG_0836.JPG



IMG_0837.JPG



IMG_0838.JPG

TEST PIT NO. 2



IMG_0839.JPG

TEST PIT NO. 1



IMG_0811.JPG



IMG_0812.JPG



IMG_0813.JPG



IMG_0814.JPG

TEST PIT NO. 1



IMG_0815.JPG



IMG_0816.JPG



IMG_0817.JPG



IMG_0818.JPG

TEST PIT NO. 1



IMG_0819.JPG



IMG_0820.JPG



IMG_0821.JPG



IMG_0822.JPG

TEST PIT NO. 1



IMG_0823.JPG



IMG_0824.JPG



IMG_0825.JPG



IMG_0826.JPG

TEST PIT NO. 1



IMG_0827.JPG



IMG_0828.JPG



IMG_0829.JPG



IMG_0830.JPG

Appendix K

Landfill Gas Laboratory Data Sheets



CHAIN-OF-CUSTODY RECORD

Sample Transportation Notice

Relinquishing signature on this document indicates that sample is being shipped in compliance with all applicable local, State, Federal, national, and international laws, regulations and ordinances of any kind. Air Toxics Limited assumes no liability with respect to the collection, handling or shipping of these samples. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Air Toxics Limited against any claim, demand, or action, of any kind, related to the collection, handling, or shipping of samples. D.O.T. Hotline (800) 467-4922

180 BLUE RAVINE ROAD, SUITE B
FOLSOM, CA 95630-4719
(916) 985-1000 FAX (916) 985-1020

Page ___ of ___

Project Manager MIKE WARFEL
 Collected by: (Print and Sign) STEVE ENGEL
 Company PARAMETRIX Email engel@parametrix.com
 Address 411 108th AVE NE City Belleve State WA Zip 98004
 Phone 425-458-6200 Fax 425-458-6363

| | | |
|-------------------------------------|---|--|
| Project Info: | Turn Around Time: | Lab Use Only |
| | <input checked="" type="checkbox"/> Normal <input type="checkbox"/> Rush <small>specify</small> | Pressurized by: Date: Pressurization Gas: N ₂ He |
| P.O. # | | |
| Project # <u>215-1578-121(03)</u> | | |
| Project Name <u>OLALLA LF 0303A</u> | | |

| Lab I.D. | Field Sample I.D. (Location) | Can # | Date of Collection | Time of Collection | Analyses Requested | Canister Pressure/Vacuum | | | |
|----------|------------------------------|-------|--------------------|--------------------|--------------------|--------------------------|-------|---------|-------------|
| | | | | | | Initial | Final | Receipt | Final (psi) |
| 01A | FL-1 | 36570 | 10/15/10 | 1025 | T0-15 | -29"vc | | | |
| 02A | FL-2 | 34170 | 10/15/10 | 1045 | T0-15 | -730"vc | | | |
| 03A | FL-3 | 34582 | 10/15/10 | 1110 | T0-15 | -730"vc | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |

| | | |
|--|--|--------|
| Relinquished by: (signature) <u>[Signature]</u> Date/Time <u>10/15/10 1230</u> | Received by: (signature) <u>[Signature]</u> Date/Time <u>10-15-10 0905</u> | Notes: |
| Relinquished by: (signature) _____ Date/Time _____ | Received by: (signature) _____ Date/Time _____ | |
| Relinquished by: (signature) _____ Date/Time _____ | Received by: (signature) _____ Date/Time _____ | |

| | | | | | | |
|--------------|--------------|------------|-----------|-------------|-----------------------|----------------|
| Lab Use Only | Shipper Name | Air Bill # | Temp (°C) | Condition | Custody Seals Intact? | Work Order # |
| | <u>FedEx</u> | | <u>NA</u> | <u>GOOD</u> | Yes No <u>None</u> | <u>1010352</u> |

10/28/2010

Mr. Michael Warfel
Parametrix, Inc.
5814 Graham Avenue, Suite 201
PO Box 460
Sumner WA 98390

Project Name: Olalla LF
Project #: 215-1578-121(03/0303P)
Workorder #: 1010352

Dear Mr. Michael Warfel

The following report includes the data for the above referenced project for sample(s) received on 10/18/2010 at Air Toxics Ltd.

The data and associated QC analyzed by Modified TO-15 are compliant with the project requirements or laboratory criteria with the exception of the deviations noted in the attached case narrative.

Thank you for choosing Air Toxics Ltd. for your air analysis needs. Air Toxics Ltd. is committed to providing accurate data of the highest quality. Please feel free to contact the Project Manager: Karen Lopez at 916-985-1000 if you have any questions regarding the data in this report.

Regards,



Karen Lopez
Project Manager

WORK ORDER #: 1010352

Work Order Summary

| | | | |
|------------------------|---|------------------|---|
| CLIENT: | Mr. Michael Warfel Parametrix, Inc. 5814 Graham Avenue, Suite 201 PO Box 460 Sumner, WA 98390 | BILL TO: | Mr. Stephen Emge Parametrix, Inc. 411 108th Avenue Suite 1800 Bellevue, WA 98004-5571 |
| PHONE: | 253-863-5128 | P.O. # | |
| FAX: | 253-826-2873 | PROJECT # | 215-1578-121(03/0303P) Olalla LF |
| DATE RECEIVED: | 10/18/2010 | CONTACT: | Karen Lopez |
| DATE COMPLETED: | 10/28/2010 | | |

| <u>FRACTION #</u> | <u>NAME</u> | <u>TEST</u> | <u>RECEIPT VAC./PRES.</u> | <u>FINAL PRESSURE</u> |
|-------------------|-------------|----------------|-------------------------------|---------------------------|
| 01A | FL-1 | Modified TO-15 | 0.6 psi | 15 psi |
| 02A | FL-2 | Modified TO-15 | 0.6 "Hg | 15 psi |
| 03A | FL-3 | Modified TO-15 | 1.0 "Hg | 15 psi |
| 04A | Lab Blank | Modified TO-15 | NA | NA |
| 05A | CCV | Modified TO-15 | NA | NA |
| 06A | LCS | Modified TO-15 | NA | NA |
| 06AA | LCS D | Modified TO-15 | NA | NA |

CERTIFIED BY: 

DATE: 10/28/10

Laboratory Director

Certification numbers: CA NELAP - 02110CA, LA NELAP/LELAP- AI 30763,
NY NELAP - 11291, UT NELAP - 9166389892, AZ Licensure AZ0719

Name of Accrediting Agency: NELAP/Florida Department of Health, Scope of Application: Clean Air Act,
Accreditation number: E87680, Effective date: 07/01/09, Expiration date: 06/30/11

Air Toxics Ltd. certifies that the test results contained in this report meet all requirements of the NELAC standards

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

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(916) 985-1000 . (800) 985-5955 . FAX (916) 985-1020

**LABORATORY NARRATIVE
EPA Method TO-15
Parametrix, Inc.
Workorder# 1010352**

Three 1 Liter Summa Canister samples were received on October 18, 2010. The laboratory performed analysis via modified EPA Method TO-15 using GC/MS in the full scan mode.

This workorder was independently validated prior to submittal using 'USEPA National Functional Guidelines' as generally applied to the analysis of volatile organic compounds in air. A rules-based, logic driven, independent validation engine was employed to assess completeness, evaluate pass/fail of relevant project quality control requirements and verification of all quantified amounts.

Receiving Notes

There were no receiving discrepancies.

Analytical Notes

All Quality Control Limit exceedences and affected sample results are noted by flags. Each flag is defined at the bottom of this Case Narrative and on each Sample Result Summary page. Target compound non-detects in the samples that are associated with high bias in QC analyses have not been flagged.

Definition of Data Qualifying Flags

Eight qualifiers may have been used on the data analysis sheets and indicates as follows:

B - Compound present in laboratory blank greater than reporting limit (background subtraction not performed).

J - Estimated value.

E - Exceeds instrument calibration range.

S - Saturated peak.

Q - Exceeds quality control limits.

U - Compound analyzed for but not detected above the reporting limit.

UJ- Non-detected compound associated with low bias in the CCV

N - The identification is based on presumptive evidence.

File extensions may have been used on the data analysis sheets and indicates as follows:

a-File was requantified

b-File was quantified by a second column and detector

r1-File was requantified for the purpose of reissue

Summary of Detected Compounds
MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

Client Sample ID: FL-1

Lab ID#: 1010352-01A

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|-----------------|--------------------------|----------------------|---------------------------|-----------------------|
| Freon 12 | 0.97 | 1.3 | 4.8 | 6.6 |
| Freon 11 | 0.97 | 2.0 | 5.4 | 11 |
| Acetone | 3.9 | 12 | 9.2 | 28 |
| Heptane | 0.97 | 11 | 4.0 | 44 |
| Toluene | 0.97 | 5.8 | 3.6 | 22 |

Client Sample ID: FL-2

Lab ID#: 1010352-02A

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|-----------------|--------------------------|----------------------|---------------------------|-----------------------|
| Freon 12 | 1.0 | 39 | 5.1 | 190 |
| Freon 114 | 1.0 | 25 | 7.2 | 180 |
| Freon 11 | 1.0 | 17 | 5.8 | 94 |
| Acetone | 4.1 | 47 | 9.8 | 110 |
| Hexane | 1.0 | 34 | 3.6 | 120 |
| Heptane | 1.0 | 63 | 4.2 | 260 |
| Toluene | 1.0 | 12 | 3.9 | 45 |

Client Sample ID: FL-3

Lab ID#: 1010352-03A

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|-----------------|--------------------------|----------------------|---------------------------|-----------------------|
| Freon 12 | 1.0 | 4.9 | 5.2 | 24 |
| Freon 114 | 1.0 | 3.7 | 7.3 | 26 |
| Freon 11 | 1.0 | 6.5 | 5.9 | 36 |
| Acetone | 4.2 | 8.0 | 9.9 | 19 |
| Heptane | 1.0 | 9.1 | 4.3 | 37 |
| Toluene | 1.0 | 5.7 | 3.9 | 22 |

Client Sample ID: FL-1

Lab ID#: 1010352-01A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102324 | Date of Collection: 10/15/10 10:25:00 A |
| Dil. Factor: | 1.94 | Date of Analysis: 10/23/10 07:29 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|----------------------------------|--------------------------|----------------------|---------------------------|-----------------------|
| Freon 12 | 0.97 | 1.3 | 4.8 | 6.6 |
| Freon 114 | 0.97 | Not Detected | 6.8 | Not Detected |
| Chloromethane | 3.9 | Not Detected | 8.0 | Not Detected |
| Vinyl Chloride | 0.97 | Not Detected | 2.5 | Not Detected |
| 1,3-Butadiene | 0.97 | Not Detected | 2.1 | Not Detected |
| Bromomethane | 0.97 | Not Detected | 3.8 | Not Detected |
| Chloroethane | 0.97 | Not Detected | 2.6 | Not Detected |
| Freon 11 | 0.97 | 2.0 | 5.4 | 11 |
| Ethanol | 3.9 | Not Detected | 7.3 | Not Detected |
| Freon 113 | 0.97 | Not Detected | 7.4 | Not Detected |
| 1,1-Dichloroethene | 0.97 | Not Detected | 3.8 | Not Detected |
| Acetone | 3.9 | 12 | 9.2 | 28 |
| 2-Propanol | 3.9 | Not Detected | 9.5 | Not Detected |
| Carbon Disulfide | 0.97 | Not Detected | 3.0 | Not Detected |
| 3-Chloropropene | 3.9 | Not Detected | 12 | Not Detected |
| Methylene Chloride | 0.97 | Not Detected | 3.4 | Not Detected |
| Methyl tert-butyl ether | 0.97 | Not Detected | 3.5 | Not Detected |
| trans-1,2-Dichloroethene | 0.97 | Not Detected | 3.8 | Not Detected |
| Hexane | 0.97 | Not Detected | 3.4 | Not Detected |
| 1,1-Dichloroethane | 0.97 | Not Detected | 3.9 | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 0.97 | Not Detected | 2.9 | Not Detected |
| cis-1,2-Dichloroethene | 0.97 | Not Detected | 3.8 | Not Detected |
| Tetrahydrofuran | 0.97 | Not Detected | 2.9 | Not Detected |
| Chloroform | 0.97 | Not Detected | 4.7 | Not Detected |
| 1,1,1-Trichloroethane | 0.97 | Not Detected | 5.3 | Not Detected |
| Cyclohexane | 0.97 | Not Detected | 3.3 | Not Detected |
| Carbon Tetrachloride | 0.97 | Not Detected | 6.1 | Not Detected |
| 2,2,4-Trimethylpentane | 0.97 | Not Detected | 4.5 | Not Detected |
| Benzene | 0.97 | Not Detected | 3.1 | Not Detected |
| 1,2-Dichloroethane | 0.97 | Not Detected | 3.9 | Not Detected |
| Heptane | 0.97 | 11 | 4.0 | 44 |
| Trichloroethene | 0.97 | Not Detected | 5.2 | Not Detected |
| 1,2-Dichloropropane | 0.97 | Not Detected | 4.5 | Not Detected |
| 1,4-Dioxane | 3.9 | Not Detected | 14 | Not Detected |
| Bromodichloromethane | 0.97 | Not Detected | 6.5 | Not Detected |
| cis-1,3-Dichloropropene | 0.97 | Not Detected | 4.4 | Not Detected |
| 4-Methyl-2-pentanone | 0.97 | Not Detected | 4.0 | Not Detected |
| Toluene | 0.97 | 5.8 | 3.6 | 22 |
| trans-1,3-Dichloropropene | 0.97 | Not Detected | 4.4 | Not Detected |
| 1,1,2-Trichloroethane | 0.97 | Not Detected | 5.3 | Not Detected |
| Tetrachloroethene | 0.97 | Not Detected | 6.6 | Not Detected |

Client Sample ID: FL-1

Lab ID#: 1010352-01A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102324 | Date of Collection: 10/15/10 10:25:00 A |
| Dil. Factor: | 1.94 | Date of Analysis: 10/23/10 07:29 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|---------------------------|--------------------------|----------------------|---------------------------|-----------------------|
| 2-Hexanone | 3.9 | Not Detected | 16 | Not Detected |
| Dibromochloromethane | 0.97 | Not Detected | 8.3 | Not Detected |
| 1,2-Dibromoethane (EDB) | 0.97 | Not Detected | 7.4 | Not Detected |
| Chlorobenzene | 0.97 | Not Detected | 4.5 | Not Detected |
| Ethyl Benzene | 0.97 | Not Detected | 4.2 | Not Detected |
| m,p-Xylene | 0.97 | Not Detected | 4.2 | Not Detected |
| o-Xylene | 0.97 | Not Detected | 4.2 | Not Detected |
| Styrene | 0.97 | Not Detected | 4.1 | Not Detected |
| Bromoform | 0.97 | Not Detected | 10 | Not Detected |
| Cumene | 0.97 | Not Detected | 4.8 | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.97 | Not Detected | 6.6 | Not Detected |
| Propylbenzene | 0.97 | Not Detected | 4.8 | Not Detected |
| 4-Ethyltoluene | 0.97 | Not Detected | 4.8 | Not Detected |
| 1,3,5-Trimethylbenzene | 0.97 | Not Detected | 4.8 | Not Detected |
| 1,2,4-Trimethylbenzene | 0.97 | Not Detected | 4.8 | Not Detected |
| 1,3-Dichlorobenzene | 0.97 | Not Detected | 5.8 | Not Detected |
| 1,4-Dichlorobenzene | 0.97 | Not Detected | 5.8 | Not Detected |
| alpha-Chlorotoluene | 0.97 | Not Detected | 5.0 | Not Detected |
| 1,2-Dichlorobenzene | 0.97 | Not Detected | 5.8 | Not Detected |
| 1,2,4-Trichlorobenzene | 3.9 | Not Detected | 29 | Not Detected |
| Hexachlorobutadiene | 3.9 | Not Detected | 41 | Not Detected |

Container Type: 1 Liter Summa Canister

| Surrogates | %Recovery | Method Limits |
|-----------------------|------------------|----------------------|
| Toluene-d8 | 94 | 70-130 |
| 1,2-Dichloroethane-d4 | 112 | 70-130 |
| 4-Bromofluorobenzene | 94 | 70-130 |

Client Sample ID: FL-2

Lab ID#: 1010352-02A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102326 | Date of Collection: 10/15/10 10:40:00 A |
| Dil. Factor: | 2.06 | Date of Analysis: 10/23/10 08:22 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|----------------------------------|--------------------------|----------------------|---------------------------|-----------------------|
| Freon 12 | 1.0 | 39 | 5.1 | 190 |
| Freon 114 | 1.0 | 25 | 7.2 | 180 |
| Chloromethane | 4.1 | Not Detected | 8.5 | Not Detected |
| Vinyl Chloride | 1.0 | Not Detected | 2.6 | Not Detected |
| 1,3-Butadiene | 1.0 | Not Detected | 2.3 | Not Detected |
| Bromomethane | 1.0 | Not Detected | 4.0 | Not Detected |
| Chloroethane | 1.0 | Not Detected | 2.7 | Not Detected |
| Freon 11 | 1.0 | 17 | 5.8 | 94 |
| Ethanol | 4.1 | Not Detected | 7.8 | Not Detected |
| Freon 113 | 1.0 | Not Detected | 7.9 | Not Detected |
| 1,1-Dichloroethene | 1.0 | Not Detected | 4.1 | Not Detected |
| Acetone | 4.1 | 47 | 9.8 | 110 |
| 2-Propanol | 4.1 | Not Detected | 10 | Not Detected |
| Carbon Disulfide | 1.0 | Not Detected | 3.2 | Not Detected |
| 3-Chloropropene | 4.1 | Not Detected | 13 | Not Detected |
| Methylene Chloride | 1.0 | Not Detected | 3.6 | Not Detected |
| Methyl tert-butyl ether | 1.0 | Not Detected | 3.7 | Not Detected |
| trans-1,2-Dichloroethene | 1.0 | Not Detected | 4.1 | Not Detected |
| Hexane | 1.0 | 34 | 3.6 | 120 |
| 1,1-Dichloroethane | 1.0 | Not Detected | 4.2 | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 1.0 | Not Detected | 3.0 | Not Detected |
| cis-1,2-Dichloroethene | 1.0 | Not Detected | 4.1 | Not Detected |
| Tetrahydrofuran | 1.0 | Not Detected | 3.0 | Not Detected |
| Chloroform | 1.0 | Not Detected | 5.0 | Not Detected |
| 1,1,1-Trichloroethane | 1.0 | Not Detected | 5.6 | Not Detected |
| Cyclohexane | 1.0 | Not Detected | 3.5 | Not Detected |
| Carbon Tetrachloride | 1.0 | Not Detected | 6.5 | Not Detected |
| 2,2,4-Trimethylpentane | 1.0 | Not Detected | 4.8 | Not Detected |
| Benzene | 1.0 | Not Detected | 3.3 | Not Detected |
| 1,2-Dichloroethane | 1.0 | Not Detected | 4.2 | Not Detected |
| Heptane | 1.0 | 63 | 4.2 | 260 |
| Trichloroethene | 1.0 | Not Detected | 5.5 | Not Detected |
| 1,2-Dichloropropane | 1.0 | Not Detected | 4.8 | Not Detected |
| 1,4-Dioxane | 4.1 | Not Detected | 15 | Not Detected |
| Bromodichloromethane | 1.0 | Not Detected | 6.9 | Not Detected |
| cis-1,3-Dichloropropene | 1.0 | Not Detected | 4.7 | Not Detected |
| 4-Methyl-2-pentanone | 1.0 | Not Detected | 4.2 | Not Detected |
| Toluene | 1.0 | 12 | 3.9 | 45 |
| trans-1,3-Dichloropropene | 1.0 | Not Detected | 4.7 | Not Detected |
| 1,1,2-Trichloroethane | 1.0 | Not Detected | 5.6 | Not Detected |
| Tetrachloroethene | 1.0 | Not Detected | 7.0 | Not Detected |

Client Sample ID: FL-2

Lab ID#: 1010352-02A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102326 | Date of Collection: 10/15/10 10:40:00 A |
| Dil. Factor: | 2.06 | Date of Analysis: 10/23/10 08:22 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|---------------------------|--------------------------|----------------------|---------------------------|-----------------------|
| 2-Hexanone | 4.1 | Not Detected | 17 | Not Detected |
| Dibromochloromethane | 1.0 | Not Detected | 8.8 | Not Detected |
| 1,2-Dibromoethane (EDB) | 1.0 | Not Detected | 7.9 | Not Detected |
| Chlorobenzene | 1.0 | Not Detected | 4.7 | Not Detected |
| Ethyl Benzene | 1.0 | Not Detected | 4.5 | Not Detected |
| m,p-Xylene | 1.0 | Not Detected | 4.5 | Not Detected |
| o-Xylene | 1.0 | Not Detected | 4.5 | Not Detected |
| Styrene | 1.0 | Not Detected | 4.4 | Not Detected |
| Bromoform | 1.0 | Not Detected | 11 | Not Detected |
| Cumene | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,1,2,2-Tetrachloroethane | 1.0 | Not Detected | 7.1 | Not Detected |
| Propylbenzene | 1.0 | Not Detected | 5.1 | Not Detected |
| 4-Ethyltoluene | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,3,5-Trimethylbenzene | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,2,4-Trimethylbenzene | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,3-Dichlorobenzene | 1.0 | Not Detected | 6.2 | Not Detected |
| 1,4-Dichlorobenzene | 1.0 | Not Detected | 6.2 | Not Detected |
| alpha-Chlorotoluene | 1.0 | Not Detected | 5.3 | Not Detected |
| 1,2-Dichlorobenzene | 1.0 | Not Detected | 6.2 | Not Detected |
| 1,2,4-Trichlorobenzene | 4.1 | Not Detected | 30 | Not Detected |
| Hexachlorobutadiene | 4.1 | Not Detected | 44 | Not Detected |

Container Type: 1 Liter Summa Canister

| Surrogates | %Recovery | Method Limits |
|-----------------------|------------------|----------------------|
| Toluene-d8 | 99 | 70-130 |
| 1,2-Dichloroethane-d4 | 106 | 70-130 |
| 4-Bromofluorobenzene | 100 | 70-130 |

Client Sample ID: FL-3

Lab ID#: 1010352-03A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102325 | Date of Collection: 10/15/10 11:10:00 A |
| Dil. Factor: | 2.09 | Date of Analysis: 10/23/10 07:52 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|----------------------------------|--------------------------|----------------------|---------------------------|-----------------------|
| Freon 12 | 1.0 | 4.9 | 5.2 | 24 |
| Freon 114 | 1.0 | 3.7 | 7.3 | 26 |
| Chloromethane | 4.2 | Not Detected | 8.6 | Not Detected |
| Vinyl Chloride | 1.0 | Not Detected | 2.7 | Not Detected |
| 1,3-Butadiene | 1.0 | Not Detected | 2.3 | Not Detected |
| Bromomethane | 1.0 | Not Detected | 4.0 | Not Detected |
| Chloroethane | 1.0 | Not Detected | 2.8 | Not Detected |
| Freon 11 | 1.0 | 6.5 | 5.9 | 36 |
| Ethanol | 4.2 | Not Detected | 7.9 | Not Detected |
| Freon 113 | 1.0 | Not Detected | 8.0 | Not Detected |
| 1,1-Dichloroethene | 1.0 | Not Detected | 4.1 | Not Detected |
| Acetone | 4.2 | 8.0 | 9.9 | 19 |
| 2-Propanol | 4.2 | Not Detected | 10 | Not Detected |
| Carbon Disulfide | 1.0 | Not Detected | 3.2 | Not Detected |
| 3-Chloropropene | 4.2 | Not Detected | 13 | Not Detected |
| Methylene Chloride | 1.0 | Not Detected | 3.6 | Not Detected |
| Methyl tert-butyl ether | 1.0 | Not Detected | 3.8 | Not Detected |
| trans-1,2-Dichloroethene | 1.0 | Not Detected | 4.1 | Not Detected |
| Hexane | 1.0 | Not Detected | 3.7 | Not Detected |
| 1,1-Dichloroethane | 1.0 | Not Detected | 4.2 | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 1.0 | Not Detected | 3.1 | Not Detected |
| cis-1,2-Dichloroethene | 1.0 | Not Detected | 4.1 | Not Detected |
| Tetrahydrofuran | 1.0 | Not Detected | 3.1 | Not Detected |
| Chloroform | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,1,1-Trichloroethane | 1.0 | Not Detected | 5.7 | Not Detected |
| Cyclohexane | 1.0 | Not Detected | 3.6 | Not Detected |
| Carbon Tetrachloride | 1.0 | Not Detected | 6.6 | Not Detected |
| 2,2,4-Trimethylpentane | 1.0 | Not Detected | 4.9 | Not Detected |
| Benzene | 1.0 | Not Detected | 3.3 | Not Detected |
| 1,2-Dichloroethane | 1.0 | Not Detected | 4.2 | Not Detected |
| Heptane | 1.0 | 9.1 | 4.3 | 37 |
| Trichloroethene | 1.0 | Not Detected | 5.6 | Not Detected |
| 1,2-Dichloropropane | 1.0 | Not Detected | 4.8 | Not Detected |
| 1,4-Dioxane | 4.2 | Not Detected | 15 | Not Detected |
| Bromodichloromethane | 1.0 | Not Detected | 7.0 | Not Detected |
| cis-1,3-Dichloropropene | 1.0 | Not Detected | 4.7 | Not Detected |
| 4-Methyl-2-pentanone | 1.0 | Not Detected | 4.3 | Not Detected |
| Toluene | 1.0 | 5.7 | 3.9 | 22 |
| trans-1,3-Dichloropropene | 1.0 | Not Detected | 4.7 | Not Detected |
| 1,1,2-Trichloroethane | 1.0 | Not Detected | 5.7 | Not Detected |
| Tetrachloroethene | 1.0 | Not Detected | 7.1 | Not Detected |

Client Sample ID: FL-3

Lab ID#: 1010352-03A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102325 | Date of Collection: 10/15/10 11:10:00 A |
| Dil. Factor: | 2.09 | Date of Analysis: 10/23/10 07:52 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|---------------------------|--------------------------|----------------------|---------------------------|-----------------------|
| 2-Hexanone | 4.2 | Not Detected | 17 | Not Detected |
| Dibromochloromethane | 1.0 | Not Detected | 8.9 | Not Detected |
| 1,2-Dibromoethane (EDB) | 1.0 | Not Detected | 8.0 | Not Detected |
| Chlorobenzene | 1.0 | Not Detected | 4.8 | Not Detected |
| Ethyl Benzene | 1.0 | Not Detected | 4.5 | Not Detected |
| m,p-Xylene | 1.0 | Not Detected | 4.5 | Not Detected |
| o-Xylene | 1.0 | Not Detected | 4.5 | Not Detected |
| Styrene | 1.0 | Not Detected | 4.4 | Not Detected |
| Bromoform | 1.0 | Not Detected | 11 | Not Detected |
| Cumene | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,1,2,2-Tetrachloroethane | 1.0 | Not Detected | 7.2 | Not Detected |
| Propylbenzene | 1.0 | Not Detected | 5.1 | Not Detected |
| 4-Ethyltoluene | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,3,5-Trimethylbenzene | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,2,4-Trimethylbenzene | 1.0 | Not Detected | 5.1 | Not Detected |
| 1,3-Dichlorobenzene | 1.0 | Not Detected | 6.3 | Not Detected |
| 1,4-Dichlorobenzene | 1.0 | Not Detected | 6.3 | Not Detected |
| alpha-Chlorotoluene | 1.0 | Not Detected | 5.4 | Not Detected |
| 1,2-Dichlorobenzene | 1.0 | Not Detected | 6.3 | Not Detected |
| 1,2,4-Trichlorobenzene | 4.2 | Not Detected | 31 | Not Detected |
| Hexachlorobutadiene | 4.2 | Not Detected | 44 | Not Detected |

Container Type: 1 Liter Summa Canister

| Surrogates | %Recovery | Method Limits |
|-----------------------|------------------|----------------------|
| Toluene-d8 | 96 | 70-130 |
| 1,2-Dichloroethane-d4 | 107 | 70-130 |
| 4-Bromofluorobenzene | 98 | 70-130 |

Client Sample ID: Lab Blank

Lab ID#: 1010352-04A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102309 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 10/23/10 12:42 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|----------------------------------|--------------------------|----------------------|---------------------------|-----------------------|
| Freon 12 | 0.50 | Not Detected | 2.5 | Not Detected |
| Freon 114 | 0.50 | Not Detected | 3.5 | Not Detected |
| Chloromethane | 2.0 | Not Detected | 4.1 | Not Detected |
| Vinyl Chloride | 0.50 | Not Detected | 1.3 | Not Detected |
| 1,3-Butadiene | 0.50 | Not Detected | 1.1 | Not Detected |
| Bromomethane | 0.50 | Not Detected | 1.9 | Not Detected |
| Chloroethane | 0.50 | Not Detected | 1.3 | Not Detected |
| Freon 11 | 0.50 | Not Detected | 2.8 | Not Detected |
| Ethanol | 2.0 | Not Detected | 3.8 | Not Detected |
| Freon 113 | 0.50 | Not Detected | 3.8 | Not Detected |
| 1,1-Dichloroethene | 0.50 | Not Detected | 2.0 | Not Detected |
| Acetone | 2.0 | Not Detected | 4.8 | Not Detected |
| 2-Propanol | 2.0 | Not Detected | 4.9 | Not Detected |
| Carbon Disulfide | 0.50 | Not Detected | 1.6 | Not Detected |
| 3-Chloropropene | 2.0 | Not Detected | 6.3 | Not Detected |
| Methylene Chloride | 0.50 | Not Detected | 1.7 | Not Detected |
| Methyl tert-butyl ether | 0.50 | Not Detected | 1.8 | Not Detected |
| trans-1,2-Dichloroethene | 0.50 | Not Detected | 2.0 | Not Detected |
| Hexane | 0.50 | Not Detected | 1.8 | Not Detected |
| 1,1-Dichloroethane | 0.50 | Not Detected | 2.0 | Not Detected |
| 2-Butanone (Methyl Ethyl Ketone) | 0.50 | Not Detected | 1.5 | Not Detected |
| cis-1,2-Dichloroethene | 0.50 | Not Detected | 2.0 | Not Detected |
| Tetrahydrofuran | 0.50 | Not Detected | 1.5 | Not Detected |
| Chloroform | 0.50 | Not Detected | 2.4 | Not Detected |
| 1,1,1-Trichloroethane | 0.50 | Not Detected | 2.7 | Not Detected |
| Cyclohexane | 0.50 | Not Detected | 1.7 | Not Detected |
| Carbon Tetrachloride | 0.50 | Not Detected | 3.1 | Not Detected |
| 2,2,4-Trimethylpentane | 0.50 | Not Detected | 2.3 | Not Detected |
| Benzene | 0.50 | Not Detected | 1.6 | Not Detected |
| 1,2-Dichloroethane | 0.50 | Not Detected | 2.0 | Not Detected |
| Heptane | 0.50 | Not Detected | 2.0 | Not Detected |
| Trichloroethene | 0.50 | Not Detected | 2.7 | Not Detected |
| 1,2-Dichloropropane | 0.50 | Not Detected | 2.3 | Not Detected |
| 1,4-Dioxane | 2.0 | Not Detected | 7.2 | Not Detected |
| Bromodichloromethane | 0.50 | Not Detected | 3.4 | Not Detected |
| cis-1,3-Dichloropropene | 0.50 | Not Detected | 2.3 | Not Detected |
| 4-Methyl-2-pentanone | 0.50 | Not Detected | 2.0 | Not Detected |
| Toluene | 0.50 | Not Detected | 1.9 | Not Detected |
| trans-1,3-Dichloropropene | 0.50 | Not Detected | 2.3 | Not Detected |
| 1,1,2-Trichloroethane | 0.50 | Not Detected | 2.7 | Not Detected |
| Tetrachloroethene | 0.50 | Not Detected | 3.4 | Not Detected |

Client Sample ID: Lab Blank

Lab ID#: 1010352-04A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102309 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 10/23/10 12:42 PM |

| Compound | Rpt. Limit (ppbv) | Amount (ppbv) | Rpt. Limit (ug/m3) | Amount (ug/m3) |
|---------------------------|------------------------------|--------------------------|-------------------------------|---------------------------|
| 2-Hexanone | 2.0 | Not Detected | 8.2 | Not Detected |
| Dibromochloromethane | 0.50 | Not Detected | 4.2 | Not Detected |
| 1,2-Dibromoethane (EDB) | 0.50 | Not Detected | 3.8 | Not Detected |
| Chlorobenzene | 0.50 | Not Detected | 2.3 | Not Detected |
| Ethyl Benzene | 0.50 | Not Detected | 2.2 | Not Detected |
| m,p-Xylene | 0.50 | Not Detected | 2.2 | Not Detected |
| o-Xylene | 0.50 | Not Detected | 2.2 | Not Detected |
| Styrene | 0.50 | Not Detected | 2.1 | Not Detected |
| Bromoform | 0.50 | Not Detected | 5.2 | Not Detected |
| Cumene | 0.50 | Not Detected | 2.4 | Not Detected |
| 1,1,2,2-Tetrachloroethane | 0.50 | Not Detected | 3.4 | Not Detected |
| Propylbenzene | 0.50 | Not Detected | 2.4 | Not Detected |
| 4-Ethyltoluene | 0.50 | Not Detected | 2.4 | Not Detected |
| 1,3,5-Trimethylbenzene | 0.50 | Not Detected | 2.4 | Not Detected |
| 1,2,4-Trimethylbenzene | 0.50 | Not Detected | 2.4 | Not Detected |
| 1,3-Dichlorobenzene | 0.50 | Not Detected | 3.0 | Not Detected |
| 1,4-Dichlorobenzene | 0.50 | Not Detected | 3.0 | Not Detected |
| alpha-Chlorotoluene | 0.50 | Not Detected | 2.6 | Not Detected |
| 1,2-Dichlorobenzene | 0.50 | Not Detected | 3.0 | Not Detected |
| 1,2,4-Trichlorobenzene | 2.0 | Not Detected | 15 | Not Detected |
| Hexachlorobutadiene | 2.0 | Not Detected | 21 | Not Detected |

Container Type: NA - Not Applicable

| Surrogates | %Recovery | Method Limits |
|-----------------------|------------------|--------------------------|
| Toluene-d8 | 97 | 70-130 |
| 1,2-Dichloroethane-d4 | 103 | 70-130 |
| 4-Bromofluorobenzene | 90 | 70-130 |

Client Sample ID: CCV

Lab ID#: 1010352-05A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102304 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 10/23/10 09:46 AM |

| Compound | %Recovery |
|----------------------------------|------------------|
| Freon 12 | 123 |
| Freon 114 | 112 |
| Chloromethane | 130 |
| Vinyl Chloride | 120 |
| 1,3-Butadiene | 117 |
| Bromomethane | 118 |
| Chloroethane | 125 |
| Freon 11 | 114 |
| Ethanol | 124 |
| Freon 113 | 105 |
| 1,1-Dichloroethene | 109 |
| Acetone | 117 |
| 2-Propanol | 117 |
| Carbon Disulfide | 116 |
| 3-Chloropropene | 110 |
| Methylene Chloride | 123 |
| Methyl tert-butyl ether | 103 |
| trans-1,2-Dichloroethene | 115 |
| Hexane | 110 |
| 1,1-Dichloroethane | 112 |
| 2-Butanone (Methyl Ethyl Ketone) | 113 |
| cis-1,2-Dichloroethene | 106 |
| Tetrahydrofuran | 112 |
| Chloroform | 109 |
| 1,1,1-Trichloroethane | 107 |
| Cyclohexane | 106 |
| Carbon Tetrachloride | 109 |
| 2,2,4-Trimethylpentane | 112 |
| Benzene | 115 |
| 1,2-Dichloroethane | 115 |
| Heptane | 116 |
| Trichloroethene | 110 |
| 1,2-Dichloropropane | 112 |
| 1,4-Dioxane | 106 |
| Bromodichloromethane | 115 |
| cis-1,3-Dichloropropene | 112 |
| 4-Methyl-2-pentanone | 119 |
| Toluene | 110 |
| trans-1,3-Dichloropropene | 121 |
| 1,1,2-Trichloroethane | 116 |
| Tetrachloroethene | 114 |

Client Sample ID: CCV

Lab ID#: 1010352-05A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102304 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 10/23/10 09:46 AM |

| Compound | %Recovery |
|---------------------------|------------------|
| 2-Hexanone | 126 |
| Dibromochloromethane | 123 |
| 1,2-Dibromoethane (EDB) | 118 |
| Chlorobenzene | 114 |
| Ethyl Benzene | 119 |
| m,p-Xylene | 122 |
| o-Xylene | 117 |
| Styrene | 122 |
| Bromoform | 123 |
| Cumene | 120 |
| 1,1,2,2-Tetrachloroethane | 116 |
| Propylbenzene | 119 |
| 4-Ethyltoluene | 124 |
| 1,3,5-Trimethylbenzene | 126 |
| 1,2,4-Trimethylbenzene | 124 |
| 1,3-Dichlorobenzene | 117 |
| 1,4-Dichlorobenzene | 118 |
| alpha-Chlorotoluene | 133 Q |
| 1,2-Dichlorobenzene | 118 |
| 1,2,4-Trichlorobenzene | 111 |
| Hexachlorobutadiene | 111 |

Q = Exceeds Quality Control limits.

Container Type: NA - Not Applicable

| Surrogates | %Recovery | Method Limits |
|-----------------------|------------------|----------------------|
| Toluene-d8 | 106 | 70-130 |
| 1,2-Dichloroethane-d4 | 103 | 70-130 |
| 4-Bromofluorobenzene | 98 | 70-130 |

Client Sample ID: LCS

Lab ID#: 1010352-06A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102305 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 10/23/10 10:28 AM |

| Compound | %Recovery |
|----------------------------------|------------------|
| Freon 12 | 104 |
| Freon 114 | 110 |
| Chloromethane | 108 |
| Vinyl Chloride | 101 |
| 1,3-Butadiene | 114 |
| Bromomethane | 94 |
| Chloroethane | 109 |
| Freon 11 | 117 |
| Ethanol | 94 |
| Freon 113 | 102 |
| 1,1-Dichloroethene | 93 |
| Acetone | 103 |
| 2-Propanol | 107 |
| Carbon Disulfide | 102 |
| 3-Chloropropene | 101 |
| Methylene Chloride | 98 |
| Methyl tert-butyl ether | 112 |
| trans-1,2-Dichloroethene | 109 |
| Hexane | 116 |
| 1,1-Dichloroethane | 109 |
| 2-Butanone (Methyl Ethyl Ketone) | 114 |
| cis-1,2-Dichloroethene | 103 |
| Tetrahydrofuran | 117 |
| Chloroform | 109 |
| 1,1,1-Trichloroethane | 115 |
| Cyclohexane | 118 |
| Carbon Tetrachloride | 119 |
| 2,2,4-Trimethylpentane | 114 |
| Benzene | 114 |
| 1,2-Dichloroethane | 109 |
| Heptane | 119 |
| Trichloroethene | 110 |
| 1,2-Dichloropropane | 113 |
| 1,4-Dioxane | 111 |
| Bromodichloromethane | 113 |
| cis-1,3-Dichloropropene | 113 |
| 4-Methyl-2-pentanone | 120 |
| Toluene | 109 |
| trans-1,3-Dichloropropene | 121 |
| 1,1,2-Trichloroethane | 116 |
| Tetrachloroethene | 113 |

Client Sample ID: LCS

Lab ID#: 1010352-06A

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102305 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 10/23/10 10:28 AM |

| Compound | %Recovery |
|---------------------------|------------------|
| 2-Hexanone | 127 |
| Dibromochloromethane | 123 |
| 1,2-Dibromoethane (EDB) | 123 |
| Chlorobenzene | 116 |
| Ethyl Benzene | 122 |
| m,p-Xylene | 127 |
| o-Xylene | 122 |
| Styrene | 125 |
| Bromoform | 126 |
| Cumene | 122 |
| 1,1,2,2-Tetrachloroethane | 120 |
| Propylbenzene | 120 |
| 4-Ethyltoluene | 125 |
| 1,3,5-Trimethylbenzene | 130 |
| 1,2,4-Trimethylbenzene | 130 |
| 1,3-Dichlorobenzene | 122 |
| 1,4-Dichlorobenzene | 124 |
| alpha-Chlorotoluene | 137 Q |
| 1,2-Dichlorobenzene | 124 |
| 1,2,4-Trichlorobenzene | 110 |
| Hexachlorobutadiene | 111 |

Q = Exceeds Quality Control limits.

Container Type: NA - Not Applicable

| Surrogates | %Recovery | Method Limits |
|-----------------------|------------------|----------------------|
| Toluene-d8 | 105 | 70-130 |
| 1,2-Dichloroethane-d4 | 106 | 70-130 |
| 4-Bromofluorobenzene | 100 | 70-130 |

Client Sample ID: LCS D

Lab ID#: 1010352-06AA

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102306 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 10/23/10 10:52 AM |

| Compound | %Recovery |
|----------------------------------|------------------|
| Freon 12 | 104 |
| Freon 114 | 105 |
| Chloromethane | 108 |
| Vinyl Chloride | 102 |
| 1,3-Butadiene | 110 |
| Bromomethane | 93 |
| Chloroethane | 111 |
| Freon 11 | 132 Q |
| Ethanol | 98 |
| Freon 113 | 98 |
| 1,1-Dichloroethene | 92 |
| Acetone | 102 |
| 2-Propanol | 104 |
| Carbon Disulfide | 103 |
| 3-Chloropropene | 102 |
| Methylene Chloride | 100 |
| Methyl tert-butyl ether | 107 |
| trans-1,2-Dichloroethene | 109 |
| Hexane | 114 |
| 1,1-Dichloroethane | 106 |
| 2-Butanone (Methyl Ethyl Ketone) | 112 |
| cis-1,2-Dichloroethene | 103 |
| Tetrahydrofuran | 115 |
| Chloroform | 107 |
| 1,1,1-Trichloroethane | 110 |
| Cyclohexane | 112 |
| Carbon Tetrachloride | 114 |
| 2,2,4-Trimethylpentane | 112 |
| Benzene | 112 |
| 1,2-Dichloroethane | 108 |
| Heptane | 117 |
| Trichloroethene | 109 |
| 1,2-Dichloropropane | 112 |
| 1,4-Dioxane | 108 |
| Bromodichloromethane | 113 |
| cis-1,3-Dichloropropene | 113 |
| 4-Methyl-2-pentanone | 119 |
| Toluene | 107 |
| trans-1,3-Dichloropropene | 122 |
| 1,1,2-Trichloroethane | 116 |
| Tetrachloroethene | 114 |

Client Sample ID: LCSD

Lab ID#: 1010352-06AA

MODIFIED EPA METHOD TO-15 GC/MS FULL SCAN

| | | |
|---------------------|----------------|--|
| File Name: | 3102306 | Date of Collection: NA |
| Dil. Factor: | 1.00 | Date of Analysis: 10/23/10 10:52 AM |

| Compound | %Recovery |
|---------------------------|------------------|
| 2-Hexanone | 128 |
| Dibromochloromethane | 123 |
| 1,2-Dibromoethane (EDB) | 124 |
| Chlorobenzene | 115 |
| Ethyl Benzene | 121 |
| m,p-Xylene | 126 |
| o-Xylene | 120 |
| Styrene | 125 |
| Bromoform | 123 |
| Cumene | 120 |
| 1,1,2,2-Tetrachloroethane | 118 |
| Propylbenzene | 120 |
| 4-Ethyltoluene | 125 |
| 1,3,5-Trimethylbenzene | 128 |
| 1,2,4-Trimethylbenzene | 129 |
| 1,3-Dichlorobenzene | 121 |
| 1,4-Dichlorobenzene | 123 |
| alpha-Chlorotoluene | 132 Q |
| 1,2-Dichlorobenzene | 124 |
| 1,2,4-Trichlorobenzene | 114 |
| Hexachlorobutadiene | 115 |

Q = Exceeds Quality Control limits.

Container Type: NA - Not Applicable

| Surrogates | %Recovery | Method Limits |
|-----------------------|------------------|----------------------|
| Toluene-d8 | 105 | 70-130 |
| 1,2-Dichloroethane-d4 | 105 | 70-130 |
| 4-Bromofluorobenzene | 100 | 70-130 |

Appendix L

Opinion of Probable Costs for Alternatives

**Opinion of Probable Cost for Alternative 1
Monitored Natural Attenuation and Land Use Controls**

| Item | Quantity | Units | Unit Cost | Capital Cost | O&M Cost | | Source |
|---|----------|-------|--------------------------|-----------------|-----------|----------------------------|-----------------------------|
| | | | | | Annual | Present Worth ¹ | |
| Low Permeability Caps Low Permeability Soil Cap | | | | | \$2,500 | \$63,586 | Similar Projects |
| Monitoring Well Decommissioning Well Decommissioning | 10 | ea | \$2,000 | | | \$14,560 | Similar Project |
| Subtotal | | | | \$0 | | | |
| Contingency | | | 25% of Capital Cost | \$0 | | | |
| Construction/Project Management | | | 20% of Capital Cost | \$0 | | | |
| Engineering (PS&E) | | | 15% of Capital Cost | \$0 | | | |
| Construction Cost Subtotal | | | | \$0 | | | |
| Sales Tax | | | 8.6% | \$0 | | | |
| Environmental Oversight | | | | | | | |
| General Reporting | | | | | | | |
| Draft Groundwater Monitoring and Well Maintenance Plan | 1 | Each | \$14,000 | \$14,000 | | | Engineer's Estimate |
| Final Groundwater Monitoring and Well Maintenance Plan | 1 | Each | \$6,500 | \$6,500 | | | Engineer's Estimate |
| Annual Groundwater Monitoring Reports | 30 | Each | \$9,000 | | \$231,407 | | Engineer's Estimate |
| Quarterly Groundwater Monitoring Reports | 90 | Each | \$6,500 | | \$501,383 | | Engineer's Estimate |
| Periodic Review Report (every 5 years) | 6 | Each | \$27,700 | | \$139,342 | | Engineer's Estimate |
| Project Management | 1 | LS | \$6,370 | \$6,370 | | | Engineer's Estimate |
| Land Use Controls | | | | | | | |
| Environmental Covenant | 1 | LS | \$6,000 | \$6,000 | \$500 | \$12,717 | Engineer's Estimate |
| Draft Land Use Control Implementation Plan (LUCIP) | 1 | LS | \$5,000 | \$5,000 | | | Engineer's Estimate |
| Final LUCIP | 1 | LS | \$2,200 | \$2,200 | \$500 | \$12,717 | Engineer's Estimate |
| Notice of Conveyance or Other Transfer of an Interest | 1 | LS | \$2,000 | | | \$4,893 | Engineer's Estimate |
| Land Use Control Maintenance | 1 | LS | \$1,000 | | \$1,000 | \$25,435 | Engineer's Estimate |
| Long Term Groundwater Monitoring (yrs 1-30) | | | | | | | |
| Sample Collection (Quarterly) | 120 | ea | \$3,346 | | | \$779,612 | Actual Costs; Annual Costs |
| Sample Analysis (Quarterly) | 120 | ea | \$4,234 | | | | Eng. Estimate; Annual Costs |
| SWHP Permit Fees | 30 | ea | \$8,000 | | | \$203,476 | |
| Environmental Oversight Subtotal | | | | \$40,070 | | | |
| Operation and Maintenance Subtotal | | | | | | \$1,989,128 | |
| O&M Project Management and Support | | | 10% of O&M Present Worth | | | \$198,912.85 | |
| O&M Contingency | | | 25% of O&M Present Worth | | | \$497,282.12 | |
| Operation and Maintenance Total | | | | | | \$2,685,323 | |
| NET PRESENT WORTH | | | | | | \$2,725,393 | |

Notes:

1 - Discount rate used for all present worth calculations per EPA Guidance = 1.1%

**Opinion of Probable Cost for Alternative 2
Low Permeability Geomembrane Cap Phase I Area with MNA and Land Use Controls**

| Item | Quantity | Units | Unit Cost | Capital Cost | O&M Cost | | Source |
|--|----------|----------------------|-------------|--------------------|----------|----------------------------|-----------------------------|
| | | | | | Annual | Present Worth ¹ | |
| Low Permeability Caps | | | | | | | |
| Landfill Phase I Area | | | | | | | |
| Low Permeability Geomembrane Cap | 6.5 | Acre | \$153,500 | \$997,750 | | | Similar Projects |
| Mobilization | 8% | LS | \$997,750 | \$79,820 | \$7,500 | \$190,759 | Similar Projects |
| Monitoring Well Decommissioning | | | | | | | |
| Well Decommissioning | 10 | ea | \$2,000 | | | \$14,560 | Similar Project |
| Subtotal | | | | \$1,077,570 | | | |
| Contingency | 25% | of Capital Cost | | \$269,393 | | | |
| Construction/Project Management | 20% | of Capital Cost | | \$215,514 | | | |
| Engineering (PS&E) | 15% | of Capital Cost | | \$161,636 | | | |
| Construction Cost Subtotal | | | | \$1,724,112 | | | |
| Sales Tax | | | 8.6% | \$148,274 | | | |
| Environmental Oversight | | | | | | | |
| General Reporting | | | | | | | |
| Draft Groundwater Monitoring and Well Maintenance Plan | 1 | Each | \$14,000 | \$14,000 | | | Engineer's Estimate |
| Final Groundwater Monitoring and Well Maintenance Plan | 1 | Each | \$6,500 | \$6,500 | | | Engineer's Estimate |
| Annual Groundwater Monitoring Reports | 30 | Each | \$9,000 | | | \$231,407 | Engineer's Estimate |
| Quarterly Groundwater Monitoring Reports | 70 | Each | \$6,500 | | | \$401,916 | Engineer's Estimate |
| Periodic Review Report (every 5 years) | 6 | Each | \$27,700 | | | \$139,342 | Engineer's Estimate |
| Project Management | 1 | LS | \$6,370 | \$6,370 | | | Engineer's Estimate |
| Land Use Controls | | | | | | | |
| Environmental Covenant | 1 | LS | \$6,000 | \$6,000 | \$500 | \$12,717 | Engineer's Estimate |
| Draft Land Use Control Implementation Plan (LUCIP) | 1 | LS | \$5,000 | \$5,000 | | | Engineer's Estimate |
| Final LUCIP | 1 | LS | \$2,200 | \$2,200 | \$500 | \$12,717 | Engineer's Estimate |
| Notice of Conveyance or Other Transfer of an Interest | 1 | LS | \$2,000 | | | \$4,893 | Engineer's Estimate |
| Land Use Control Maintenance | 1 | LS | \$1,000 | | \$1,000 | \$25,435 | Engineer's Estimate |
| Long Term Groundwater Monitoring (yrs 1-30) | | | | | | | |
| Sample Collection (Quarterly) | 100 | ea | \$3,346 | | | \$663,614 | Actual Costs; Annual Costs |
| Sample Analysis (Quarterly) | 100 | ea | \$4,234 | | | | Eng. Estimate; Annual Costs |
| SWHP Permit Fees | 30 | ea | \$8,000 | | | \$203,476 | |
| Environmental Oversight Subtotal | | | | \$40,070 | | | |
| Operation and Maintenance Subtotal | | | | | | \$1,900,836 | |
| O&M Project Management and Support | 10% | of O&M Present Worth | | | | \$190,084 | |
| O&M Contingency | 25% | of O&M Present Worth | | | | \$475,209 | |
| Operation and Maintenance Total | | | | | | \$2,566,129 | |
| NET PRESENT WORTH | | | | | | \$4,478,585 | |

Notes:

1 - Discount rate used for all present worth calculations per EPA Guidance = 1.1%

**Opinion of Probable Cost for Alternative 3
In-Situ Treatment, Monitored Natural Attenuation, and Land Use Controls**

| Item | Quantity | Units | Unit Cost | Capital Cost | O&M Cost | | Source | |
|--|----------|-------|--------------------------|------------------|----------|----------------------------|--------------------|------------------------------------|
| | | | | | Annual | Present Worth ¹ | | |
| Low Permeability Caps Low Permeability Soil Cap | | | | | | \$2,500 | \$63,586 | Similar Projects |
| MRC™ Injection | | | | | | | | |
| Injection Wells | 21 | ea | \$11,600 | \$243,600 | | | | MW-8 & MW-10 Costs |
| Field Pilot Study | | | | | | | | |
| Injection Well | 1 | ea | \$11,600 | \$11,600 | | | | |
| MRC™ | 100 | lb | \$8.50 | \$850 | | | | Similar Projects; Vendor Quote |
| Freight | 1 | LS | \$500 | \$500 | | | | Similar Project |
| Dosing Labor | 1 | Day | \$2,750 | \$2,750 | | | | Similar Project |
| Water | 600 | gal | \$0.002 | \$1 | | | | Similar Project |
| Mobilization | 8% | LS | \$4,101 | \$328 | | | | Similar Project; % of Capital Cost |
| MRC™ Injection Implementation | | | | | | | | |
| MRC™ | 2,100 | lb | \$8.50 | \$17,850 | | | | Similar Projects; Vendor Quote |
| Freight | 1 | LS | \$3,900 | \$3,900 | | | | Similar Project |
| Dosing Labor | 10 | Day | \$2,750 | \$27,500 | | \$260,357 | | Similar Project |
| Water | 12,600 | gal | \$0.002 | \$25 | | | | Similar Project |
| Mobilization | 8% | LS | \$49,275 | \$3,942 | | | | Similar Project; % of Capital Cost |
| Air Sparging | | | | | | | | |
| Sparge Wells | 10 | ea | \$8,000 | \$80,000 | | | | MW-8 & MW-10 Costs |
| Field Pilot Study | 1 | LS | \$30,000 | \$30,000 | | | | Similar Project |
| Mobilization | 8% | LS | \$30,000 | \$2,400 | | | | Similar Project; % of Capital Cost |
| Air Sparging Implementation | | | | | | | | |
| Air Sparge Equipment | 1 | LS | \$40,000 | \$40,000 | | | | Similar Projects |
| Equipment Building | 400 | SF | \$100 | \$40,000 | | | | |
| Site Work | 1 | LS | \$10,000 | \$10,000 | \$12,000 | \$305,214 | | |
| Well Head Completions | 10 | ea | \$2,000 | \$20,000 | | | | |
| Mobilization | 8% | LS | \$110,000 | \$8,800 | | | | Similar Project; % of Capital Cost |
| Decommissioning | 1 | LS | \$10,000 | | | | \$7,202 | |
| Well Decommissioning | | | | | | | | |
| Well Decommissioning | 42 | ea | \$2,000 | | | | \$61,151 | Similar Project |
| Subtotal | | | | | | | | |
| Contingency | | | 25% of Capital Cost | \$136,012 | | | | |
| Construction/Project Management | | | 20% of Capital Cost | \$108,809 | | | | |
| Engineering (PS&E) | | | 15% of Capital Cost | \$81,607 | | | | |
| Construction Cost Subtotal | | | | \$870,474 | | | | |
| Sales Tax | | | 8.6% | \$74,861 | | | | |
| Environmental Oversight | | | | | | | | |
| General Reporting | | | | | | | | |
| Draft Groundwater Monitoring and Well Maintenance Plan | 1 | Each | \$14,000 | \$14,000 | | | | Engineer's Estimate |
| Final Groundwater Monitoring and Well Maintenance Plan | 1 | Each | \$6,500 | \$6,500 | | | | Engineer's Estimate |
| Annual Groundwater Monitoring Reports | 30 | Each | \$9,000 | | | \$231,407 | | Engineer's Estimate |
| Quarterly Groundwater Monitoring Reports | 90 | Each | \$6,500 | | | \$501,383 | | Engineer's Estimate |
| Annual Cleanup Action Activity Reports | 30 | Each | \$10,700 | | | \$275,118 | | Engineer's Estimate |
| Periodic Review Report (every 5 years) | 6 | Each | \$27,700 | | | \$139,342 | | Engineer's Estimate |
| Project Management | 1 | LS | \$7,440 | \$7,440 | | | | Engineer's Estimate |
| Land Use Controls ¹ | | | | | | | | |
| Environmental Covenant | 1 | LS | \$6,000 | \$6,000 | | \$500 | \$12,717 | Engineer's Estimate |
| Draft Land Use Control Implementation Plan (LUCIP) | 1 | LS | \$5,000 | \$5,000 | | | | Engineer's Estimate |
| Final LUCIP | 1 | LS | \$2,200 | \$2,200 | | \$500 | \$12,717 | Engineer's Estimate |
| Notice of Conveyance or Other Transfer of Interest | 1 | LS | \$2,000 | | | | \$4,893 | Engineer's Estimate |
| Land Use Control Maintenance | 1 | LS | \$1,000 | | | \$1,000 | \$25,435 | Engineer's Estimate |
| Short Term Groundwater Monitoring (yrs 1-5) | | | | | | | | |
| Sample Collection | 48 | ea | \$26,768 | | | | \$577,542 | Actual Costs; Annual Costs |
| Sample Analysis | 48 | ea | \$33,872 | | | | | Engineer's Estimate; Annual Costs |
| Long Term Groundwater Monitoring (yrs 1-30) | | | | | | | | |
| Sample Collection (Quarterly) | 120 | ea | \$3,346 | | | | \$779,612 | Actual Costs; Annual Costs |
| Sample Analysis (Quarterly) | 120 | ea | \$4,234 | | | | | Engineer's Estimate; Annual Costs |
| SWHP Permit Fees | 30 | ea | \$8,000 | | | | \$203,476 | |
| Environmental Oversight Subtotal | | | | \$41,140 | | | | |
| Operation and Maintenance Subtotal | | | | | | | \$3,461,153 | |
| O&M Project Management and Support | | | 10% of O&M Present Worth | | | | \$346,115 | |
| O&M Contingency | | | 25% of O&M Present Worth | | | | \$865,288 | |
| Operation and Maintenance Total | | | | | | | \$4,672,556 | |
| NET PRESENT WORTH | | | | | | | \$5,659,031 | |

Notes:

1 - Discount rate used for all present worth calculations per EPA Guidance =

1.1%