

April 21, 2015

Project No. 923-1000-002.R273

Mr. Bill Kombol  
Palmer Coking Coal Company  
31407 Highway 169  
PO Box 10  
Black Diamond, WA 98010

**RE: LANDSBURG MINE SITE INTERIM GROUNDWATER MONITORING REPORT – NOVEMBER 2014**

Dear Bill:

Golder Associates Inc. (Golder) completed an interim groundwater monitoring event at the Landsburg Mine Site during November 2014. Groundwater samples were collected from monitoring wells LMW-2, LMW-3, LMW-4, LMW-5, LMW-6, LMW-8, LMW-9, LMW-10, and LMW-11 (Figure 1). Monitoring wells LMW-2, LMW-4 and LMW-10 are completed to monitor shallow and deeper zones within the Rogers Coal Seam north of the Rogers Coal Mine subsidence trench. Monitoring wells LMW-3 and LMW-5 are completed to monitor the shallow (~40 feet depth) and deeper zone (~250 feet depth), respectively, within the Rogers Coal Seam at the south end of the mine. Figure 2 presents a cross-section along the strike at the coal seam that also depicts the location of the monitoring wells. Monitoring well LMW-8 is receiving groundwater before discharge from Portal 3 and the mine access incline at the south end of the Rogers Coal Mine. These wells lay along the primary pathways for detection of a chemical release from the mine, were one to occur. Groundwater samples were also collected from well LMW-9 and the deep well LMW-11, which monitor groundwater from within the Rogers Coal Mine near its south end. Wells LMW-9 and LMW-11 are receiving groundwater from near the top of the water table and near the bottom of the mine, respectively. Wells LMW-6 and LMW-7 monitor groundwater from the Frasier and Landsburg Coal Mines to the west and east of the Rogers Coal Mine, respectively.

Groundwater sampling was conducted in accordance with the *Draft Interim Groundwater Monitoring Plan, Landsburg Mine Site* (Golder 1997)<sup>1</sup>, and included the following activities:

- Measurement of static water levels at monitoring wells.
- Well purging to insure sample representativeness with the currently installed dedicated pumping systems.
- Measurement of field parameters including: pH, specific conductance, temperature, dissolved oxygen, Eh, and turbidity.
- Collection of representative samples in appropriate containers; dissolved metals samples were field filtered (total metals were not). The dissolved metals samples were not analyzed.
- Analyses of groundwater for volatile organic compounds (VOCs; United States Environmental Protection Agency [EPA] Method 8260C), semi-volatile organic compounds (SVOCs, EPA Method 8270D), polychlorinated biphenyls (PCBs; EPA 8082A), pesticides (EPA 8081B), priority pollutant metals (EPA Method 6010C/200.8/7470A Series), and a petroleum hydrocarbon identification scan (NWTPH-HCID).

<sup>1</sup> Golder Associates Inc. (Golder). 1997. *Draft Interim Groundwater Monitoring Plan, Landsburg Mine Site*. Prepared for the Landsburg PLP Steering Committee, Redmond, Washington.



Appendix A presents the laboratory analytical reports for all analyses. Sampling activities were documented on Sample Integrity Data Sheets (SIDS). Copies of the completed SIDS are provided in Appendix B. Appendix C shows the validated data with added qualifiers. Table 1 presents water depth measurements and elevations that were collected from wells prior to sampling activities. Groundwater levels are similar to previous monitoring periods and indicate that groundwater is discharging out both ends of the Rogers Coal Mine.

Following sample collection, all bottles were sealed, labeled, and placed in an iced cooler until delivery to the laboratory. All groundwater samples from monitoring wells were transported under chain-of-custody procedures to Analytical Resources Incorporated (ARI), of Tukwila, Washington, for analyses. Screening levels are based on maximum contaminant levels (MCLs) or State of Washington Model Toxics Control Act (MTCA) Method B groundwater cleanup levels, whichever value is less. In cases where an established MCL or Method B Cleanup Level does not exist, a similar (surrogate) compound regulatory screening level is identified for comparison.

The analytical results indicate no significant changes in groundwater conditions from those observed during the remedial investigation (RI) and on-going interim groundwater monitoring. Table 2 presents the field parameter measurements and laboratory analytical results for each groundwater sample. Laboratory analyses did not detect any VOCs, SVOCs, PCBs, pesticides, or petroleum hydrocarbon (HCID) in any of the groundwater samples except as discussed in the following paragraphs.

Carbon disulfide was detected at 0.14 J micrograms per liter ( $\mu\text{g/L}$ ) in the sample collected from LMW-10. This detection was considerably below the MTCA Method B groundwater cleanup level for carbon disulfide (800  $\mu\text{g/L}$ ). Data validation and inspection of the gas chromatographs of the carbon disulfide detection could not eliminate this analyte. Carbon disulfide has been detected at these low levels in site groundwater in previous sampling events. The detection of carbon disulfide is attributed to being present in the coal bed material as a natural constituent. Research has confirmed the presence of carbon disulfide in coal bed materials<sup>2</sup> and is similar to results previously detected at LMW-10.

One SVOC was detected for bis (2-ethylhexyl) phthalate at 12.0  $\mu\text{g/L}$  for LMW-4. This result is above the MTCA level of 6.3  $\mu\text{g/L}$ . This compound is very insoluble in water and is extremely immobile in a groundwater flow system (Koc over 100,000 milliliters per gram [ml/g]). It is very unlikely that this compound would be detected at this well without having other, more mobile contaminants being detected as well. The laboratory was asked to investigate the detected result. Upon evaluation of the laboratory data, it was determined that there was chronic laboratory contamination above the instrument baseline in all samples from the same sample data group including the laboratory method blank. Golder believes the result is due to contamination that could have been introduced during sampling, transport, or at the laboratory, since the compound is ubiquitous in many products used to handle samples and sample solvent extracts. A technical review of the laboratory raw data was performed and more detail is provided in a memo in Appendix C.

Even though the archived groundwater sample from well LMW-4 was beyond the quality assurance holding time, the laboratory re-extracted an aliquot with a different batch of extraction solvent and reanalyzed the sample for bis (2-ethylhexyl) phthalate. The result was non-detect with a reporting limit of 3.0  $\mu\text{g/L}$  and was qualified as estimated (UJ) since it was analyzed outside of the acceptable holding time. The data from the reanalysis are in Appendix A.

Since the re-analysis of the original groundwater sample from LMW-4 was beyond the acceptable holding time, Golder resampled the groundwater from LMW-4 on February 12, 2015 and had the laboratory analyze the sample for bis (2-ethylhexyl) phthalate. The results were non-detect with a reporting limit of 3.0  $\mu\text{g/L}$ . The laboratory data for the resampling of LMW-4 are found in Appendix A. The final results for bis (2-ethylhexyl) phthalate in groundwater from LMW-4 is non-detect at a reporting limit of 3.0  $\mu\text{g/L}$  as presented in Table 2.

<sup>2</sup> Kozinc, J., Treeby, M., and Zupancic-Kralj, L. 2004. Determination of Sulfur Gases from Velenje Coal Stockpile. Acta Chim. Slov., Volume 51, pages 529-536.

A sample was not collected at LMW-7 for the November 2014 round of sampling due to a malfunctioning pump. The pump was removed from the well and taken for repairs. It is expected that the pump will be repaired in time for the next round of sampling. Sampling results at LMW-7 have been consistent for the last 20 years of sampling since the pump was installed.

The laboratory data packages underwent a sample data validation. Items of note are provided in a validation memorandum in Appendix C.

The primary parameters detected in groundwater samples during this sampling event were metals that are naturally occurring. The method reporting limits (MRLs) and method detection limits (MDLs) for all analytes were at or below acceptable concentrations under the MTCA.

Several groundwater samples from site wells contained iron and manganese concentrations above State of Washington secondary drinking water levels (SMCLs) of 0.3 milligrams per liter (mg/L) and 0.05 mg/L, respectively, which are not health-based standards, but are protective of aesthetic qualities of water. Iron and manganese have been detected in mine groundwater above MTCA cleanup levels in every monitoring event at the site and are naturally occurring metals that are typically associated with groundwater from coal mines (Fuste et al. 1983)<sup>3</sup>. The concentrations of iron and manganese detected during the November 2014 sampling event are similar to concentrations detected during the RI (Golder 1996)<sup>4</sup> and the Interim Groundwater Sampling events previously conducted at the site.

The groundwater sample from the deep well (LMW-11) contained total arsenic at a concentration of 6.6 µg/L (0.0066 mg/L), which is less than the Washington State primary drinking water MCL and greater than the MTCA groundwater cleanup level of 10 µg/L and 5 µg/L, respectively. Arsenic also has been detected in groundwater from LMW-11 near or above MTCA cleanup levels during every monitoring event since LMW-11 was installed. Arsenic is also a naturally occurring metal commonly detectable in groundwater, especially in older more stagnant groundwater having low reduction-oxidation (REDOX) and dissolved oxygen levels. The MTCA groundwater cleanup level is based on typical groundwater background levels in the State of Washington. It is believed that the arsenic concentrations are naturally occurring deep within the mine where groundwater is more stagnant and its geochemistry may be different than shallow groundwater within the mine.

If you have any questions or require any additional information, please contact Douglas Morell at (425) 883-0777.

Sincerely,

**GOLDER ASSOCIATES INC.**



Jill S. Lamberts  
Project Environmental Scientist



Douglas J. Morell, PhD, LHG  
Principal

<sup>3</sup> Fuste, L.A., F.A. Packard, M.O.Fretwell, and D.P. Garland. 1983. Data Supplement To: Quality of Coal Mine Drainage in Washington, 1975-77. Open-File Report 83-205. Tacoma, Washington: US Geological Survey.

<sup>4</sup> Golder Associates Inc. (Golder). 1996. Remedial Investigation and Feasibility Study for the Landsburg Mine Site. Landsburg PLP Steering Committee.

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JSL/DJM/cl

## TABLES

Table 1: Groundwater Elevation Data Collection November 17, 2014 Landsburg Mine Site

|                             | UNITS  | LMW-1    | LMW-1a  | LMW-2   | LMW-3    | LMW-4 <sup>1</sup> | LMW-4 <sup>1,2</sup> | LMW-5    | LMW-6    | LMW-7 <sup>1</sup> | LMW-8    | LMW-9    | LMW-10   | LMW-11   | P-2      | Water Drainage | Frazier Seam Tunnel |
|-----------------------------|--------|----------|---------|---------|----------|--------------------|----------------------|----------|----------|--------------------|----------|----------|----------|----------|----------|----------------|---------------------|
| Water Depths                |        |          |         |         |          |                    |                      |          |          |                    |          |          |          |          |          |                |                     |
| Time of data collection     | ft bgs | 10:06 AM | 9:55 AM | 9:19 AM | 10:48 AM | 9:22 AM            | 11:51 AM             | 10:38 AM | 10:16 AM | 9:34 AM            | 10:45 AM | 10:54 AM | 11:09 AM | 10:26 AM | 10:42 AM | NA             | NA                  |
| Measured to Top of PVC      | ft bgs | 138.33   | 133.57  | 6.94    | 12.89    | 8.41               | 7.19                 | 14.43    | 28.06    | 212.11             | 4.70     | 100.18   | 1.67     | 158.02   | 7.49     | NA             | NA                  |
| Measured to Top of Monument | ft bgs | 139.11   | 133.77  | 7.62    | 13.65    | 9.09               | 7.86                 | 15.08    | 28.80    | 212.65             | 5.68     | 100.47   | NC       | 158.38   | 7.87     | NA             | NA                  |
|                             |        |          |         |         |          |                    |                      |          |          |                    |          |          |          |          |          |                |                     |
| Surveyed Elevation          |        |          |         |         |          |                    |                      |          |          |                    |          |          |          |          |          |                |                     |
| Top of PVC                  | ft asl | 765.16   | 759.51  | 617.73  | 656.75   | 619.26             | 619.26               | 658.27   | 632.33   | 771.51             | 646.97   | 743.99   | 618.87   | 801.87   | 651.37   | NA             | NA                  |
| Top of Monument             | ft asl | 765.89   | NC      | 618.29  | 657.48   | 619.85             | 619.85               | 658.87   | 633.00   | 771.88             | NC       | NC       | NC       | 802.20   | NC       | NA             | NA                  |
| Ground Level                | ft asl | 762.90   | 756.59  | 615.35  | 654.40   | 617.09             | 617.09               | 655.63   | 629.95   | 768.79             | 645.25   | 741.13   | 615.75   | 799.50   | 648.54   | 551.38         | 542.15              |
|                             |        |          |         |         |          |                    |                      |          |          |                    |          |          |          |          |          |                |                     |
| Corrected Water Elevation   |        |          |         |         |          |                    |                      |          |          |                    |          |          |          |          |          |                |                     |
| Using PVC elevation         | ft asl | 626.83   | 625.94  | 610.79  | 643.86   | 610.85             | 612.07               | 643.84   | 604.27   | 559.40             | 642.27   | 643.81   | 617.20   | 643.85   | 643.88   | NA             | NA                  |
| Using Monument elevation    | ft asl | 626.78   | NA      | 610.67  | 643.83   | 610.76             | 611.99               | 643.79   | 604.20   | 559.23             | NA       | NA       | NA       | 643.82   | NA       | NA             | NA                  |

Notes:  
<sup>1</sup> Data corrected to accommodate well inclination of 20° from vertical  
<sup>2</sup> Groundwater Level collected on 2/12/2015 as part of resampling effort  
NA = Not applicable  
NC = Data not collected  
ft bgs = feet below ground surface  
ft asl = feet above sea level

Table 2: November 2014 Groundwater Analytical Results Landsburg Mine Site

| ANALYTE                           | UNITS  | LMW-2      | LMW-3      | LMW-4      | LMW-4 <sup>2</sup> | LMW-5      | LMW-6      | LMW-7*     | LMW-8      | LMW-9      | LMW-10     | LMW-11     | Equipment Blank | Trip Blank | Trip Blank | Trip Blank |
|-----------------------------------|--------|------------|------------|------------|--------------------|------------|------------|------------|------------|------------|------------|------------|-----------------|------------|------------|------------|
|                                   |        | 11/18/2014 | 11/19/2014 | 11/18/2014 | 2/12/2015          | 11/19/2014 | 11/20/2014 | 11/20/2014 | 11/19/2014 | 11/19/2014 | 11/18/2014 | 11/20/2014 | 11/19/2014      | 11/18/2014 | 11/19/2014 | 11/20/2014 |
| Field Parameter                   |        |            |            |            |                    |            |            |            |            |            |            |            |                 |            |            |            |
| pH                                | std    | 7.04       | 7.83       | 7.05       | 7.00               | 7.03       | 6.95       | NA         | 6.96       | 7.11       | 8.75       | 7.27       | NA              | NA         | NA         | NA         |
| Conductivity                      | uS/cm  | 894        | 294.2      | 901        | 410                | 691        | 235.3      | NA         | 499        | 629        | 342        | 495        | NA              | NA         | NA         | NA         |
| Dissolved Oxygen                  | mg/L   | 0.13       | 0.00       | 0.01       | 0.00               | 0.00       | 0.00       | NA         | 0.60       | 0.00       | 0.04       | 0.64       | NA              | NA         | NA         | NA         |
| Temperature                       | °C     | 10.9       | 11.1       | 11.2       | 10.7               | 11.2       | 10.3       | NA         | 12.3       | 12.9       | 10.0       | 11.1       | NA              | NA         | NA         | NA         |
| E <sub>h</sub>                    | Rel mV | -90.0      | -88.0      | -97.4      | -78.0              | -95.5      | -84.5      | NA         | -87.7      | -95.1      | -85.2      | -60.0      | NA              | NA         | NA         | NA         |
| Turbidity                         | NTU    | 0.76       | 0.65       | 0.81       | 0.39               | 0.82       | 0.92       | NA         | 5.14       | 0.74       | 6.83       | 0.72       | NA              | NA         | NA         | NA         |
| Metals (Total)                    |        |            |            |            |                    |            |            |            |            |            |            |            |                 |            |            |            |
| Aluminum                          | mg/L   | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Antimony                          | mg/L   | 0.003 U    | 0.003 U    | 0.003 U    | NA                 | 0.003 U    | 0.003 U    | NA         | 0.003 U    | 0.003 U    | 0.003 U    | 0.003 U    | 0.003 U         | NA         | NA         | NA         |
| Arsenic                           | mg/L   | 0.003 U    | 0.003 U    | 0.003 U    | NA                 | 0.003 U    | 0.003 U    | NA         | 0.003 U    | 0.003 U    | 0.003 U    | 0.0066     | 0.003 U         | NA         | NA         | NA         |
| Barium                            | mg/L   | 0.500 U    | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 U      | NA         | 0.5 U      | 0.500 U    | 0.5 U      | 0.5 U      | 0.5 U           | NA         | NA         | NA         |
| Beryllium                         | mg/L   | 0.002 U    | 0.002 U    | 0.002 U    | NA                 | 0.002 U    | 0.002 U    | NA         | 0.002 U    | 0.002 U    | 0.002 U    | 0.002 U    | 0.002 U         | NA         | NA         | NA         |
| Cadmium                           | mg/L   | 0.002 U    | 0.002 U    | 0.002 U    | NA                 | 0.002 U    | 0.002 U    | NA         | 0.002 U    | 0.002 U    | 0.002 U    | 0.002 U    | 0.002 U         | NA         | NA         | NA         |
| Calcium                           | mg/L   | 113        | 38.3       | 110        | NA                 | 93.3       | 27.3       | NA         | 63.2       | 81.3       | 7.84       | 58.2       | 0.5 U           | NA         | NA         | NA         |
| Chromium                          | mg/L   | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Cobalt                            | mg/L   | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |
| Copper                            | mg/L   | 0.003 U    | 0.003 U    | 0.003 U    | NA                 | 0.003 U    | 0.003 U    | NA         | 0.003 U    | 0.003 U    | 0.003 U    | 0.003 U    | 0.003 U         | NA         | NA         | NA         |
| Iron                              | mg/L   | 0.250      | 0.2 U      | 1.09       | NA                 | 0.2 U      | 2.41       | NA         | 12.4       | 1.51       | 0.3        | 1.41       | 0.2 U           | NA         | NA         | NA         |
| Lead                              | mg/L   | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |
| Magnesium                         | mg/L   | 69.2       | 16         | 66.3       | NA                 | 52.3       | 14.0       | NA         | 33.7       | 44.3       | 2.95       | 26.3       | 1 U             | NA         | NA         | NA         |
| Manganese                         | mg/L   | 0.216      | 0.059      | 0.163      | NA                 | 0.228      | 0.034      | NA         | 0.511      | 0.164      | 0.02 U     | 0.165      | 0.02 U          | NA         | NA         | NA         |
| Mercury                           | mg/L   | 0.00002 U  | 0.00002 U  | 0.00002 U  | NA                 | 0.00002 U  | 0.00002 U  | NA         | 0.00002 U  | 0.00002 U  | 0.00002 U  | 0.00002 U  | 0.00002 U       | NA         | NA         | NA         |
| Nickel                            | mg/L   | 0.02000 U  | 0.02000 U  | 0.02000 U  | NA                 | 0.02000 U  | 0.02000 U  | NA         | 0.02000 U  | 0.02 U     | 0.02 U     | 0.02 U     | 0.02 U          | NA         | NA         | NA         |
| Potassium                         | mg/L   | 3.59       | 1.72       | 3.73       | NA                 | 2.77       | 0.72       | NA         | 2.12       | 2.49       | 1.33       | 2.1        | 0.5 U           | NA         | NA         | NA         |
| Selenium                          | mg/L   | 0.005 U    | 0.005 U    | 0.005 U    | NA                 | 0.005 U    | 0.005 U    | NA         | 0.005 U    | 0.005 U    | 0.005 U    | 0.005 U    | 0.005 U         | NA         | NA         | NA         |
| Silver                            | mg/L   | 0.003 U    | 0.003 U    | 0.003 U    | NA                 | 0.003 U    | 0.003 U    | NA         | 0.003 U    | 0.003 U    | 0.003 U    | 0.003 U    | 0.003 U         | NA         | NA         | NA         |
| Sodium                            | mg/L   | 20.7       | 10.2       | 28.4       | NA                 | 16.4       | 7.07       | NA         | 11.5       | 15.5       | 81.3       | 28.8       | 0.500 U         | NA         | NA         | NA         |
| Thallium                          | mg/L   | 0.002 U    | 0.002 U    | 0.002 U    | NA                 | 0.002 U    | 0.002 U    | NA         | 0.002 U    | 0.002 U    | 0.002 U    | 0.002 U    | 0.002 U         | NA         | NA         | NA         |
| Vanadium                          | mg/L   | 0.003 U    | 0.003 U    | 0.003 U    | NA                 | 0.003 U    | 0.003 U    | NA         | 0.003 U    | 0.003 U    | 0.003 U    | 0.003 U    | 0.003 U         | NA         | NA         | NA         |
| Zinc                              | mg/L   | 0.02 U     | 0.02 U     | 0.02 U     | NA                 | 0.02 U     | 0.02 U     | NA         | 0.02 U     | 0.02 U     | 0.02 U     | 0.02 U     | 0.02 U          | NA         | NA         | NA         |
| Volatile Organic Compounds (VOCs) |        |            |            |            |                    |            |            |            |            |            |            |            |                 |            |            |            |
| Acetone                           | µg/L   | 5 U        | 5 U        | 5 U        | NA                 | 5 U        | 5 U        | NA         | 5 U        | 5 U        | 5 U        | 5 U        | 5 U             | 5 U        | NA         | 5 U        |
| Acrolein                          | µg/L   | 2.5 U      | 2.5 U      | 2.5 U      | NA                 | 2.5 U      | 2.5 U      | NA         | 2.5 U      | 2.5 U      | 2.5 U      | 2.5 U      | 2.5 U           | 2.5 U      | NA         | 2.5 U      |
| Acrylonitrile                     | µg/L   | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | 1 U        | NA         | 1 U        |
| Benzene                           | µg/L   | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Bromobenzene                      | µg/L   | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Bromochloromethane                | µg/L   | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Bromodichloromethane              | µg/L   | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Bromoform                         | µg/L   | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Bromomethane                      | µg/L   | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | 1 U        | NA         | 1 U        |
| 2-Butanone                        | µg/L   | 5 U        | 5 U        | 5 U        | NA                 | 5 U        | 5 U        | NA         | 5 U        | 5 U        | 5 U        | 5 U        | 5 U             | 5 U        | NA         | 5 U        |
| n-Butylbenzene                    | µg/L   | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| sec-Butylbenzene                  | µg/L   | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| tert-Butylbenzene                 | µg/L   | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.200 U         | 0.2 U      | NA         | 0.2 U      |

Table 2: November 2014 Groundwater Analytical Results Landsburg Mine Site

| ANALYTE                     | UNITS | LMW-2      | LMW-3      | LMW-4      | LMW-4 <sup>2</sup> | LMW-5      | LMW-6      | LMW-7*     | LMW-8      | LMW-9      | LMW-10     | LMW-11     | Equipment Blank | Trip Blank | Trip Blank | Trip Blank |
|-----------------------------|-------|------------|------------|------------|--------------------|------------|------------|------------|------------|------------|------------|------------|-----------------|------------|------------|------------|
|                             |       | 11/18/2014 | 11/19/2014 | 11/18/2014 | 2/12/2015          | 11/19/2014 | 11/20/2014 | 11/20/2014 | 11/19/2014 | 11/19/2014 | 11/18/2014 | 11/20/2014 | 11/19/2014      | 11/18/2014 | 11/19/2014 | 11/20/2014 |
| Carbon disulfide            | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.14 J     | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Carbon tetrachloride        | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Chlorobenzene               | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Chloroethane                | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 2-Chloroethylvinylether     | µg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 UJ     | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 UJ     | 0.5 U           | 0.5 U      | NA         | 0.5 UJ     |
| Chloroform                  | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Chloromethane               | µg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 U      | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U           | 0.5 U      | NA         | 0.5 U      |
| 2-Chlorotoluene             | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | 0.1 U      | NA         | 0.1 U      |
| 4-Chlorotoluene             | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Chlorodibromomethane        | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,2-Dibromo-3-Chloropropane | µg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 U      | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U           | 0.5 U      | NA         | 0.5 U      |
| 1,2-Dibromoethane           | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | 0.1 U      | NA         | 0.1 U      |
| Dibromomethane              | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,2-Dichlorobenzene         | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,3-Dichlorobenzene         | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,4-Dichlorobenzene         | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| trans-1,4-Dichloro-2-butene | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | 1 U        | NA         | 1 U        |
| 1,1-Dichloroethane          | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,2-Dichloroethane          | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,1-Dichloroethene          | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| cis-1,2-Dichloroethene      | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| trans-1,2-Dichloroethene    | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,2-Dichloropropane         | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,3-Dichloropropane         | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | 0.1 U      | NA         | 0.1 U      |
| 2,2-Dichloropropane         | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | 0.1 U      | NA         | 0.1 U      |
| 1,1-Dichloropropene         | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | 0.1 U      | NA         | 0.1 U      |
| cis-1,3-Dichloropropene     | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| trans-1,3-Dichloropropene   | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Ethylbenzene                | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Hexachlorobutadiene         | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.33 U     | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 2-Hexanone                  | µg/L  | 5 U        | 5 U        | 5 U        | NA                 | 5 U        | 5 U        | NA         | 5 U        | 5 U        | 5 U        | 5 U        | 5 U             | 5 U        | NA         | 5 U        |
| Iodomethane                 | µg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 U      | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U           | 0.5 U      | NA         | 0.5 U      |
| Isopropylbenzene            | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 4-Isopropyltoluene          | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | 0.1 U      | NA         | 0.1 U      |
| Methylene Chloride          | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1.4 U           | 2.3 J+     | NA         | 2.6 J+     |
| 4-Methyl-2-pentanone        | µg/L  | 2.5 U      | 2.5 U      | 2.5 U      | NA                 | 2.5 U      | 2.5 U      | NA         | 2.5 U      | 2.5 U      | 2.5 U      | 2.5 U      | 2.5 U           | 2.5 U      | NA         | 2.5 U      |
| Naphthalene                 | µg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 UJ     | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 UJ     | 0.5 U           | 0.5 U      | NA         | 0.5 UJ     |
| N-Propylbenzene             | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Styrene                     | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,2,3-Trichlorobenzene      | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,2,4-Trichlorobenzene      | µg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 U      | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U           | 0.5 U      | NA         | 0.5 U      |
| 1,3,5-Trichlorobenzene      | µg/L  | NA         | NA         | NA         | NA                 | NA         | NA         | NA         | NA         | NA         | NA         | NA         | NA              | NA         | NA         | NA         |
| 1,1,1,2-Tetrachloroethane   | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,1,2,2-Tetrachloroethane   | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | 0.1 U      | NA         | 0.1 U      |
| Tetrachloroethene           | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Toluene                     | µg/L  | 0.2 U      | 0.2 U      | 0.2 U      | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |

Table 2: November 2014 Groundwater Analytical Results Landsburg Mine Site

| ANALYTE                                         | UNITS | LMW-2      | LMW-3      | LMW-4             | LMW-4 <sup>2</sup> | LMW-5      | LMW-6      | LMW-7*     | LMW-8      | LMW-9      | LMW-10     | LMW-11     | Equipment Blank | Trip Blank | Trip Blank | Trip Blank |
|-------------------------------------------------|-------|------------|------------|-------------------|--------------------|------------|------------|------------|------------|------------|------------|------------|-----------------|------------|------------|------------|
|                                                 |       | 11/18/2014 | 11/19/2014 | 11/18/2014        | 2/12/2015          | 11/19/2014 | 11/20/2014 | 11/20/2014 | 11/19/2014 | 11/19/2014 | 11/18/2014 | 11/20/2014 | 11/19/2014      | 11/18/2014 | 11/19/2014 | 11/20/2014 |
| 1,1,1-Trichloroethane                           | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,1,2-Trichloroethane                           | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Trichloroethene                                 | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Trichlorofluoromethane                          | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Total Benzofluoranthenes                        | µg/L  | 2 U        | 2 U        | 2 U               | NA                 | 2 U        | 2 U        | NA         | 2 U        | 2 U        | 2 U        | 2 U        | 2 U             | 2 U        | NA         | NA         |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (CFC-113) | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,2,3-Trichloropropane                          | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,2,4-Trimethylbenzene                          | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| 1,3,5-Trimethylbenzene                          | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Vinyl acetate                                   | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 UJ     | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 UJ     | 0.2 U           | 0.2 U      | NA         | 0.2 UJ     |
| Vinyl chloride                                  | µg/L  | 0.1 U      | 0.1 U      | 0.1 U             | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | 0.1 U      | NA         | 0.1 U      |
| m-Xylene & p-Xylene                             | µg/L  | 0.4 U      | 0.4 U      | 0.4 U             | NA                 | 0.4 U      | 0.4 U      | NA         | 0.4 U      | 0.4 U      | 0.4 U      | 0.4 U      | 0.4 U           | 0.4 U      | NA         | 0.4 U      |
| o-Xylene                                        | µg/L  | 0.2 U      | 0.2 U      | 0.2 U             | NA                 | 0.2 U      | 0.2 U      | NA         | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U      | 0.2 U           | 0.2 U      | NA         | 0.2 U      |
| Xylenes, Total                                  | µg/L  | 0.4 U      | 0.4 U      | 0.4 U             | NA                 | 0.4 U      | 0.4 U      | NA         | 0.4 U      | 0.4 U      | 0.4 U      | 0.4 U      | 0.4 U           | 0.4 U      | NA         | 0.4 U      |
| Semi-Volatile Organic Compounds (SVOCs)         |       |            |            |                   |                    |            |            |            |            |            |            |            |                 |            |            |            |
| Acenaphthene                                    | µg/L  | 1 UJ       | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Acenaphthylene                                  | µg/L  | 1 UJ       | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Anthracene                                      | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Benzo(a)anthracene                              | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Benzo(a)pyrene                                  | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Benzo(b)fluoranthene                            | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Benzo(ghi)perylene                              | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Benzo(k)fluoranthene                            | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Benzoic Acid                                    | µg/L  | 20 U       | 20 U       | 20 U              | NA                 | 20 U       | 20 U       | NA         | 20 U       | 20 U       | 20 U       | 20 U       | 20 U            | NA         | NA         | NA         |
| Benzyl Alcohol                                  | µg/L  | 2 U        | 2 U        | 2 U               | NA                 | 2 U        | 2 U        | NA         | 2 U        | 2 U        | 2 U        | 2 U        | 2 U             | NA         | NA         | NA         |
| Bis(2-chloroethoxy)methane                      | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Bis(2-chloroethyl)ether                         | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Bis(2-chloroisopropyl)ether                     | µg/L  | NA         | NA         | NA                | NA                 | NA         | NA         | NA         | NA         | NA         | NA         | NA         | NA              | NA         | NA         | NA         |
| Bis(2-ethylhexyl)phthalate                      | µg/L  | 3 U        | 3 U        | 3 UJ <sup>1</sup> | 3 U                | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| 4-Bromophenyl phenyl ether                      | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Butyl benzyl phthalate                          | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Carbazole                                       | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 4-Chloroaniline                                 | µg/L  | 5 U        | 5 U        | 5 U               | NA                 | 5 U        | 5 U        | NA         | 5 U        | 5 U        | 5 U        | 5 U        | 5 U             | NA         | NA         | NA         |
| 4-Chloro-3-methylphenol                         | µg/L  | 3 U        | 3 U        | 3 U               | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| 2-Chloronaphthalene                             | µg/L  | 1 UJ       | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 2-Chlorophenol                                  | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 4-Chlorophenyl phenyl ether                     | µg/L  | 1 UJ       | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 3 & 4-Methylphenol (m,p-Cresols)                | µg/L  | 2 U        | 2 U        | 2 U               | NA                 | 2 U        | 2 U        | NA         | 2 U        | 2 U        | 2 U        | 2 U        | 2 U             | NA         | NA         | NA         |
| 2-Methylphenol (o-Cresol)                       | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Chrysene                                        | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Di-n-butyl phthalate                            | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Dibenz(a,h)anthracene                           | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Dibenzofuran                                    | µg/L  | 1 UJ       | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 1,2-Dichlorobenzene                             | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 1,3-Dichlorobenzene                             | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 1,4-Dichlorobenzene                             | µg/L  | 1 U        | 1 U        | 1 U               | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |

Table 2: November 2014 Groundwater Analytical Results Landsburg Mine Site

| ANALYTE                          | UNITS | LMW-2      | LMW-3      | LMW-4      | LMW-4 <sup>2</sup> | LMW-5      | LMW-6      | LMW-7*     | LMW-8      | LMW-9      | LMW-10     | LMW-11     | Equipment Blank | Trip Blank | Trip Blank | Trip Blank |
|----------------------------------|-------|------------|------------|------------|--------------------|------------|------------|------------|------------|------------|------------|------------|-----------------|------------|------------|------------|
|                                  |       | 11/18/2014 | 11/19/2014 | 11/18/2014 | 2/12/2015          | 11/19/2014 | 11/20/2014 | 11/20/2014 | 11/19/2014 | 11/19/2014 | 11/18/2014 | 11/20/2014 | 11/19/2014      | 11/18/2014 | 11/19/2014 | 11/20/2014 |
| 3,3'-Dichlorobenzidine           | µg/L  | 5 U        | 5 U        | 5 U        | NA                 | 5 U        | 5 U        | NA         | 5 U        | 5 U        | 5 U        | 5 U        | 5 U             | NA         | NA         | NA         |
| 2,4-Dichlorophenol               | µg/L  | 3 U        | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| Diethyl phthalate                | µg/L  | 1 UJ       | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 2,4-Dimethylphenol               | µg/L  | 3 U        | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| Dimethyl phthalate               | µg/L  | 1 UJ       | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 4,6-Dinitro-2-methylphenol       | µg/L  | 10 U       | 10 U       | 10 U       | NA                 | 10 U       | 10 U       | NA         | 10 U       | 10 U       | 10 U       | 10 U       | 10 U            | NA         | NA         | NA         |
| 2,4-Dinitrophenol                | µg/L  | 20 UJ      | 20 U       | 20 U       | NA                 | 20 U       | 20 U       | NA         | 20 U       | 20 U       | 20 U       | 20 U       | 20 U            | NA         | NA         | NA         |
| 2,4-Dinitrotoluene               | µg/L  | 3 UJ       | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| 2,6-Dinitrotoluene               | µg/L  | 3 UJ       | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| N-Nitrosodiphenylamine           | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Fluoranthene                     | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Fluorene                         | µg/L  | 1 UJ       | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Hexachlorobenzene                | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Hexachlorobutadiene              | µg/L  | 3 U        | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| Hexachlorocyclopentadiene        | µg/L  | 5 UJ       | 5 U        | 5 U        | NA                 | 5 U        | 5 U        | NA         | 5 U        | 5 U        | 5 U        | 5 U        | 5 U             | NA         | NA         | NA         |
| Hexachloroethane                 | µg/L  | 2 U        | 2 U        | 2 U        | NA                 | 2 U        | 2 U        | NA         | 2 U        | 2 U        | 2 U        | 2 U        | 2 U             | NA         | NA         | NA         |
| Indeno(1,2,3-cd)pyrene           | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Isophorone                       | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 1-Methylnaphthalene              | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 2-Methylnaphthalene              | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 4-Methylphenol                   | µg/L  | 2 U        | 2 U        | 2 U        | NA                 | 2 U        | 2 U        | NA         | 2 U        | 2 U        | 2 U        | 2 U        | 2 U             | NA         | NA         | NA         |
| Naphthalene                      | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 2-Nitroaniline                   | µg/L  | 3 UJ       | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| 3-Nitroaniline                   | µg/L  | 3 UJ       | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| 4-Nitroaniline                   | µg/L  | 3 UJ       | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| Nitrobenzene                     | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 2-Nitrophenol                    | µg/L  | 3 U        | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| 4-Nitrophenol                    | µg/L  | 10 UJ      | 10 U       | 10 U       | NA                 | 10 U       | 10 U       | NA         | 10 U       | 10 U       | 10 U       | 10 U       | 10 U            | NA         | NA         | NA         |
| N-Nitrosodi-n-propylamine        | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 2,2'-Oxybis(1-Chloropropane)     | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Di-n-octyl phthalate             | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Pentachlorophenol                | µg/L  | 10 U       | 10 U       | 10 U       | NA                 | 10 U       | 10 U       | NA         | 10 U       | 10 U       | 10 U       | 10 U       | 10 U            | NA         | NA         | NA         |
| Phenanthrene                     | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Phenol                           | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| Pyrene                           | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 1,2,4-Trichlorobenzene           | µg/L  | 1 U        | 1 U        | 1 U        | NA                 | 1 U        | 1 U        | NA         | 1 U        | 1 U        | 1 U        | 1 U        | 1 U             | NA         | NA         | NA         |
| 2,4,5-Trichlorophenol            | µg/L  | 5 UJ       | 5 U        | 5 U        | NA                 | 5 U        | 5 U        | NA         | 5 U        | 5 U        | 5 U        | 5 U        | 5 U             | NA         | NA         | NA         |
| 2,4,6-Trichlorophenol            | µg/L  | 3 UJ       | 3 U        | 3 U        | NA                 | 3 U        | 3 U        | NA         | 3 U        | 3 U        | 3 U        | 3 U        | 3 U             | NA         | NA         | NA         |
| Polychlorinated Biphenyls (PCBs) |       |            |            |            |                    |            |            |            |            |            |            |            |                 |            |            |            |
| Aroclor 1016                     | µg/L  | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |
| Aroclor 1221                     | µg/L  | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |
| Aroclor 1232                     | µg/L  | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |
| Aroclor 1242                     | µg/L  | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |
| Aroclor 1248                     | µg/L  | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |
| Aroclor 1254                     | µg/L  | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |
| Aroclor 1260                     | µg/L  | 0.01 U     | 0.01 U     | 0.01 U     | NA                 | 0.01 U     | 0.01 U     | NA         | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U     | 0.01 U          | NA         | NA         | NA         |

Table 2: November 2014 Groundwater Analytical Results Landsburg Mine Site

| ANALYTE                    | UNITS | LMW-2      | LMW-3      | LMW-4      | LMW-4 <sup>2</sup> | LMW-5      | LMW-6      | LMW-7*     | LMW-8      | LMW-9      | LMW-10     | LMW-11     | Equipment Blank | Trip Blank | Trip Blank | Trip Blank |
|----------------------------|-------|------------|------------|------------|--------------------|------------|------------|------------|------------|------------|------------|------------|-----------------|------------|------------|------------|
|                            |       | 11/18/2014 | 11/19/2014 | 11/18/2014 | 2/12/2015          | 11/19/2014 | 11/20/2014 | 11/20/2014 | 11/19/2014 | 11/19/2014 | 11/18/2014 | 11/20/2014 | 11/19/2014      | 11/18/2014 | 11/19/2014 | 11/20/2014 |
| Pesticides                 |       |            |            |            |                    |            |            |            |            |            |            |            |                 |            |            |            |
| Aldrin (2C)                | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| alpha-BHC (2C)             | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| beta-BHC (2C)              | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| delta-BHC (2C)             | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| gamma-BHC (2C)             | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| cis-Chlordane              | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| trans-Chlordane            | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| 4,4'-DDD (2C)              | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| 4,4'-DDE (2C)              | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| 4,4'-DDT (2C)              | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| Dieldrin (2C)              | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| Endosulfan I (2C)          | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| Endosulfan II (2C)         | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| Endosulfan sulfate (2C)    | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| Endrin                     | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| Endrin aldehyde (2C)       | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| Endrin ketone (2C)         | µg/L  | 0.1 U      | 0.1 U      | 0.1 U      | NA                 | 0.1 U      | 0.1 U      | NA         | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U      | 0.1 U           | NA         | NA         | NA         |
| Heptachlor (2C)            | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| Heptachlor epoxide (2C)    | µg/L  | 0.05 U     | 0.05 U     | 0.05 U     | NA                 | 0.05 U     | 0.05 U     | NA         | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U     | 0.05 U          | NA         | NA         | NA         |
| Methoxychlor (2C)          | µg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 U      | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U           | NA         | NA         | NA         |
| Toxaphene                  | µg/L  | 5 U        | 5 U        | 5 U        | NA                 | 5 U        | 5 U        | NA         | 5 U        | 5 U        | 5 U        | 5 U        | 5 U             | NA         | NA         | NA         |
| Hydrocarbon Identification |       |            |            |            |                    |            |            |            |            |            |            |            |                 |            |            |            |
| Diesel Range               | mg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 U      | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U           | NA         | NA         | NA         |
| Gas Range                  | mg/L  | 0.25 U     | 0.25 U     | 0.25 U     | NA                 | 0.25 U     | 0.25 U     | NA         | 0.25 U     | 0.25 U     | 0.25 U     | 0.25 U     | 0.25 U          | NA         | NA         | NA         |
| Lube Oil                   | mg/L  | 0.5 U      | 0.5 U      | 0.5 U      | NA                 | 0.5 U      | 0.5 U      | NA         | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U      | 0.5 U           | NA         | NA         | NA         |

Notes:

NA = Not Analyzed

U - The analyte was not detected above the level of the reporting limit.

UJ - The analyte was not detected above the reporting limit and is estimated.

J+ - The analyte was detected and is estimated with a high bias.

\* LMW-7 was not sampled due to pump malfunction. No Field Duplicate sample was collected.

µS/cm = microsiemens per centimeter

mg/L = milligrams per liter

Rel mV = relative millivolts

NTU = nephelometric turbidity unit

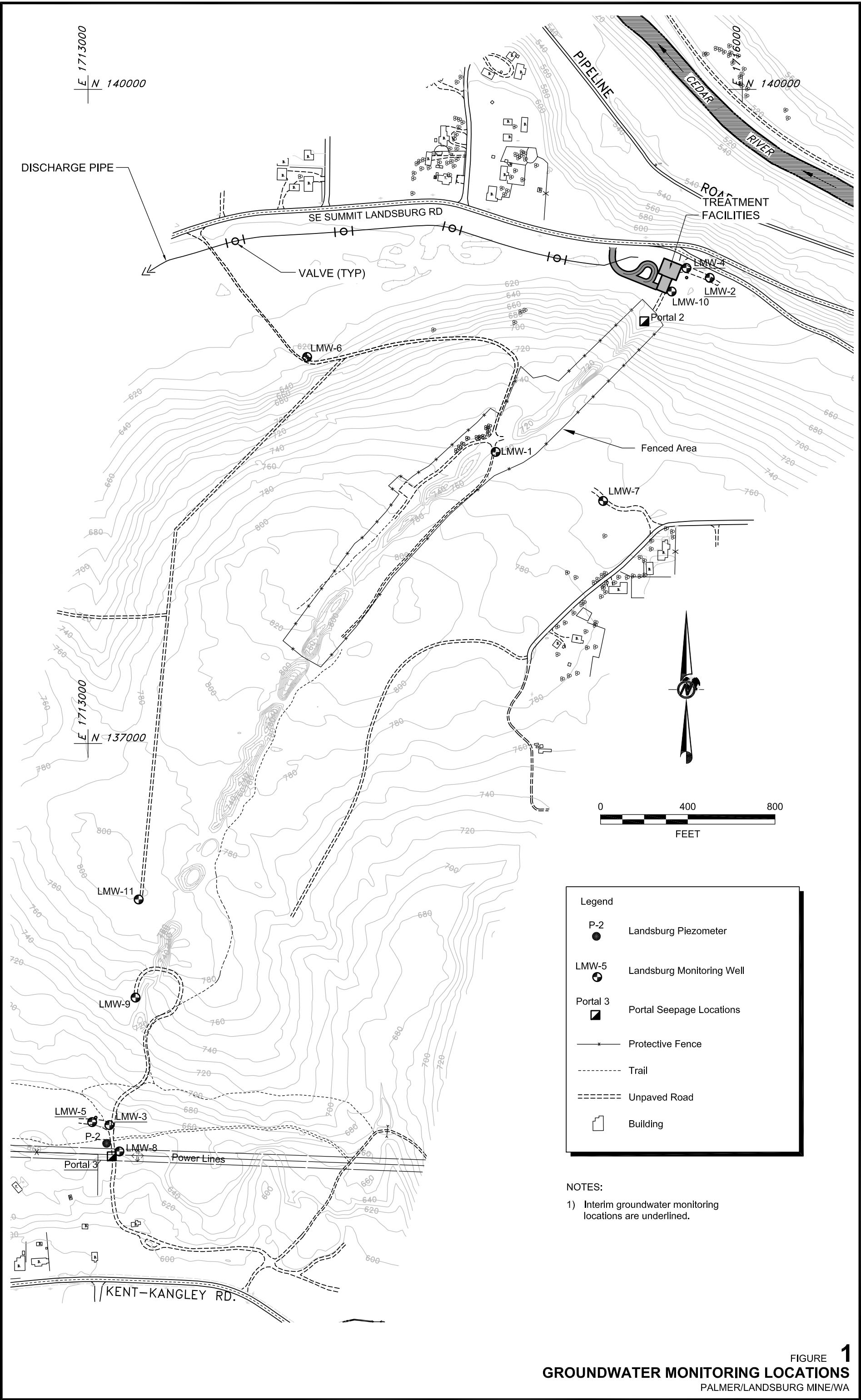
µg/L = micrograms per liter

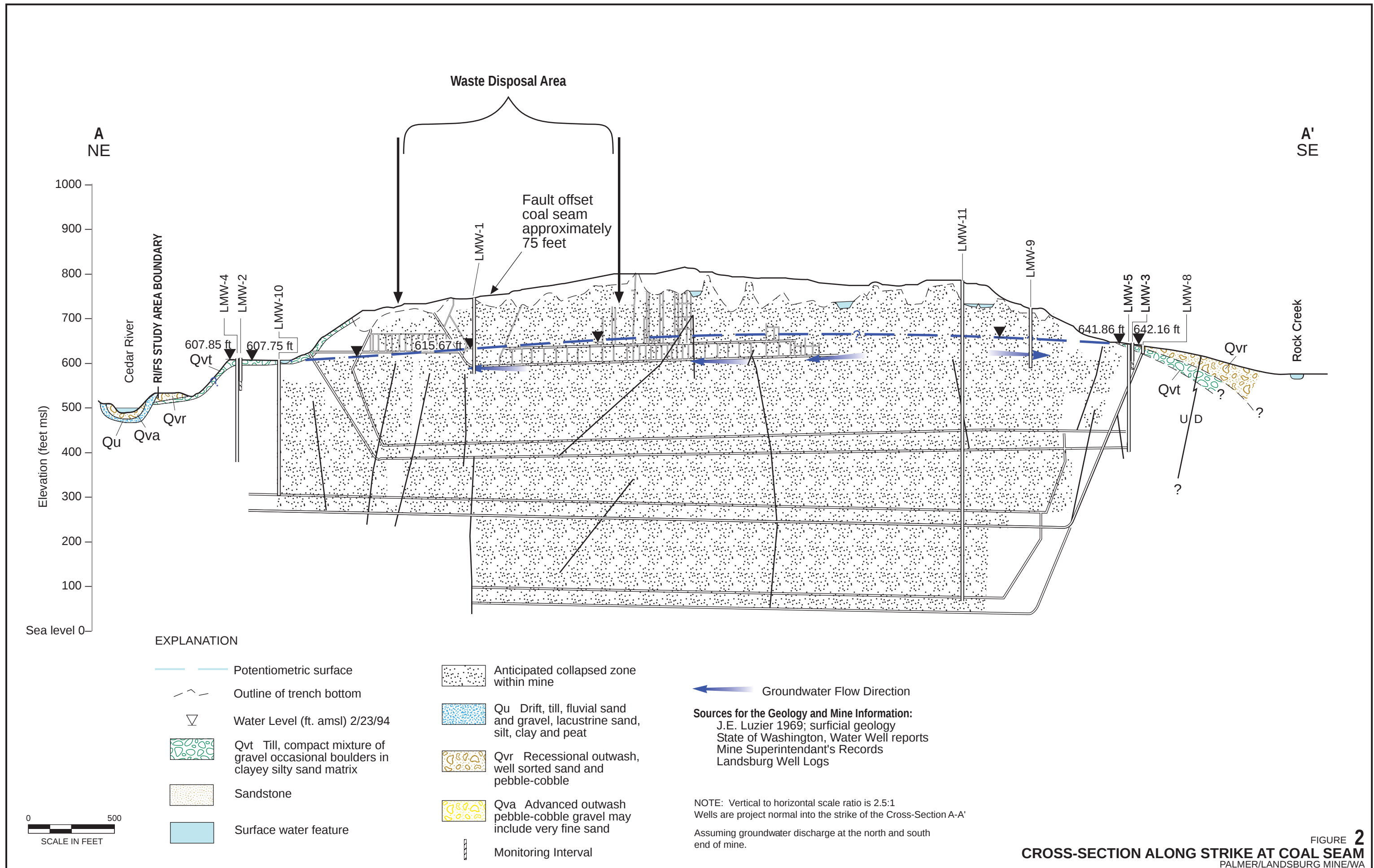
<sup>1</sup> Initial analysis had detection of 12 µg/L. Sample was re-extracted out of hold time and then reanalyzed. Reanalysis was non-detect and was qualified as UJ.

<sup>2</sup> Location was resampled on 2/12/2015 to confirm result.



## FIGURES





**FIGURE 2**  
**CROSS-SECTION ALONG STRIKE AT COAL SEAM**  
 PALMER/LANDBURG MINE/WA  
**Golder Associates**

**APPENDIX A**  
**LABORATORY ANALYTICAL REPORTS**



## **Data Reporting Qualifiers**

**Effective 12/31/13**

### **Inorganic Data**

- U Indicates that the target analyte was not detected at the reported concentration
- \* Duplicate RPD is not within established control limits
- B Reported value is less than the CRDL but  $\geq$  the Reporting Limit
- N Matrix Spike recovery not within established control limits
- NA Not Applicable, analyte not spiked
- H The natural concentration of the spiked element is so much greater than the concentration spiked that an accurate determination of spike recovery is not possible
- L Analyte concentration is  $\leq 5$  times the Reporting Limit and the replicate control limit defaults to  $\pm 1$  RL instead of the normal 20% RPD

### **Organic Data**

- U Indicates that the target analyte was not detected at the reported concentration
- \* Flagged value is not within established control limits
- B Analyte detected in an associated Method Blank at a concentration greater than one-half of ARI's Reporting Limit or 5% of the regulatory limit or 5% of the analyte concentration in the sample.
- J Estimated concentration when the value is less than ARI's established reporting limits
- D The spiked compound was not detected due to sample extract dilution
- E Estimated concentration calculated for an analyte response above the valid instrument calibration range. A dilution is required to obtain an accurate quantification of the analyte.



- Q Indicates a detected analyte with an initial or continuing calibration that does not meet established acceptance criteria (<20%RSD, <20%Drift or minimum RRF).
- S Indicates an analyte response that has saturated the detector. The calculated concentration is not valid; a dilution is required to obtain valid quantification of the analyte
- NA The flagged analyte was not analyzed for
- NR Spiked compound recovery is not reported due to chromatographic interference
- NS The flagged analyte was not spiked into the sample
- M Estimated value for an analyte detected and confirmed by an analyst but with low spectral match parameters. This flag is used only for GC-MS analyses
- N The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification"
- Y The analyte is not detected at or above the reported concentration. The reporting limit is raised due to chromatographic interference. The Y flag is equivalent to the U flag with a raised reporting limit.
- EMPC Estimated Maximum Possible Concentration (EMPC) defined in EPA Statement of Work DLM02.2 as a value "calculated for 2,3,7,8-substituted isomers for which the quantitation and /or confirmation ion(s) has signal to noise in excess of 2.5, but does not meet identification criteria" **(Dioxin/Furan analysis only)**
- C The analyte was positively identified on only one of two chromatographic columns. Chromatographic interference prevented a positive identification on the second column
- P The analyte was detected on both chromatographic columns but the quantified values differ by  $\geq 40\%$  RPD with no obvious chromatographic interference
- X Analyte signal includes interference from polychlorinated diphenyl ethers. **(Dioxin/Furan analysis only)**
- Z Analyte signal includes interference from the sample matrix or perfluorokerosene ions. **(Dioxin/Furan analysis only)**



## **Geotechnical Data**

- A      The total of all fines fractions. This flag is used to report total fines when only sieve analysis is requested and balances total grain size with sample weight.
- F      Samples were frozen prior to particle size determination
- SM     Sample matrix was not appropriate for the requested analysis. This normally refers to samples contaminated with an organic product that interferes with the sieving process and/or moisture content, porosity and saturation calculations
- SS     Sample did not contain the proportion of “fines” required to perform the pipette portion of the grain size analysis
- W      Weight of sample in some pipette aliquots was below the level required for accurate weighting



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants

December 19, 2014

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Redmond, WA 98052-3333

**Client Project Name: Landsburg Mine**  
**Client Project Number: 923-1000-002.R273**  
**ARI ID: ZL38 and ZL39**

Dear Mr. Morell:

Please find enclosed Chain-of-Custody (COC) record, sample receipt documentation, and the final results for the project referenced above. Analytical Resources, Inc. (ARI) accepted eight water samples and two trip blanks in good condition on December 4, 2014. There were no discrepancies between the COC and the sample containers' labels.

The samples were analyzed for VOCs, PCBs, HCID, Pesticides, SVOCs, Total Metals, as requested on the COC. Quality control analyses are included for your review.

The VOCs method blank contained methylene chloride, naphthalene, n-Butylbenzene, 1,2,3-Trichlorobenzene and hexachlorobutadiene. All associated samples that contain analyte have been flagged with a "B" qualifier.

The SVOCs surrogate TBP is out of control low in association with sample LMW-2-1114. All other surrogate recoveries are in control.

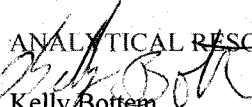
No other analytical complications were noted.

Per client request, the metals reporting limits were raised to meet client required limits.

An electronic copy of this report and all supporting raw data will remain on file at ARI. Please feel free to contact me if you have any questions or require any additional information.

Respectfully,

ANALYTICAL RESOURCES, INC.

  
Kelly Bottem  
Client Services Manager  
(206) 695-6211  
[kellyb@arilabs.com](mailto:kellyb@arilabs.com)

# Chain of Custody Record & Laboratory Analysis Request

|                                                      |                                           |                             |                                  |
|------------------------------------------------------|-------------------------------------------|-----------------------------|----------------------------------|
| ARI Assigned Number:<br><b>ZL38</b>                  | Turn-around Requested:<br><b>Standard</b> | Page: <b>1</b> of <b>1</b>  |                                  |
| ARI Client Company:<br><b>Golder Associates</b>      |                                           | Date:<br><b>11/18/2014</b>  | Ice Present?<br><b>Y</b>         |
| Client Contact:<br><b>Doug Morell, Jill Lamberts</b> |                                           | No. of Coolers:<br><b>7</b> | Cooler Temps:<br><b>0.41-4.7</b> |
| Client Project Name:<br><b>Landsburg</b>             |                                           |                             |                                  |
| Client Project #:<br><b>923-1000-002.R273</b>        | Samplers:<br><b>Rydecki, Lamberts</b>     |                             |                                  |



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants  
4611 South 134th Place, Suite 100  
Tukwila, WA 98168  
206-695-6200 206-695-6201 (fax)  
www.arilabs.com

| Sample ID                                         | Date     | Time | Matrix | No. Containers | Analysis Requested                                 |                               |            |            |            |                                                |          |       |               |       | Notes/Comments                  |             |  |  |  |                             |  |  |  |  |
|---------------------------------------------------|----------|------|--------|----------------|----------------------------------------------------|-------------------------------|------------|------------|------------|------------------------------------------------|----------|-------|---------------|-------|---------------------------------|-------------|--|--|--|-----------------------------|--|--|--|--|
|                                                   |          |      |        |                | VOCs                                               | Client list                   | PCBs (CLL) | Pesticides | SVOCs 8270 | Client list                                    | TPH-HClD | TAM/L | -Total Metals | TAM/L |                                 | Diss Metals |  |  |  |                             |  |  |  |  |
| Trip Blank-111814                                 | 11/18/14 | -    | W      | 9-6            | X                                                  |                               |            |            |            |                                                |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-10-1114                                       |          | 945  | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              | HOLD     |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-2-1114                                        |          | 1140 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-4-1114                                        |          | 1310 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| Trip Blank-111914                                 | 11/19/14 | -    | W      | 6              | X                                                  | HOLD unless there are detects |            |            |            |                                                |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-3-1114                                        |          | 0920 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              | HOLD     |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| EB-1114                                           |          | 0900 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-8-1114                                        |          | 1005 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-5-1114                                        |          | 1130 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-9-1114                                        |          | 1315 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| Comments/Special Instructions<br>-Ecology ETM EDD |          |      |        |                | Relinquished by:<br>(Signature) <i>[Signature]</i> |                               |            |            |            | Received by:<br>(Signature) <i>[Signature]</i> |          |       |               |       | Relinquished by:<br>(Signature) |             |  |  |  | Received by:<br>(Signature) |  |  |  |  |
| Printed Name:<br><b>J. Lamberts</b>               |          |      |        |                | Printed Name:<br><b>J. Lamberts</b>                |                               |            |            |            | Printed Name:<br><b>Taylor Stricker</b>        |          |       |               |       | Printed Name:                   |             |  |  |  | Printed Name:               |  |  |  |  |
| Company:<br><b>Golder</b>                         |          |      |        |                | Company:<br><b>Golder</b>                          |                               |            |            |            | Company:<br><b>ARL</b>                         |          |       |               |       | Company:                        |             |  |  |  | Company:                    |  |  |  |  |
| Date & Time:<br><b>11/19/14 1442</b>              |          |      |        |                | Date & Time:<br><b>11/19/14 1442</b>               |                               |            |            |            | Date & Time:                                   |          |       |               |       | Date & Time:                    |             |  |  |  |                             |  |  |  |  |

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, notwithstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.



# Cooler Receipt Form

ARI Client: Bolder

Project Name: Landsburg

COC No(s): \_\_\_\_\_ (NA)

Delivered by: Fed-Ex UPS Courier Hand Delivered Other: \_\_\_\_\_

Assigned ARI Job No: ZL38 | 2L39

Tracking No: \_\_\_\_\_ (NA)

## Preliminary Examination Phase:

Were intact, properly signed and dated custody seals attached to the outside of to cooler? \_\_\_\_\_

YES ☐ NO ☒

Were custody papers included with the cooler? \_\_\_\_\_

YES ☒ NO ☐

Were custody papers properly filled out (ink, signed, etc.) \_\_\_\_\_

YES ☒ NO ☐

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry)

Time: 1447

3.3 3.6 1.3 3.8 2.1 4.7 0.4

If cooler temperature is out of compliance fill out form 00070F

Temp Gun ID#: \_\_\_\_\_

Cooler Accepted by: TS

Date: 11/19/14

Time: 1447

**Complete custody forms and attach all shipping documents**

## Log-In Phase:

Was a temperature blank included in the cooler? \_\_\_\_\_

YES ☐ NO ☒

What kind of packing material was used? ... Bubble Wrap ☒ Wet Ice ☒ Gel Packs ☒ Baggies ☒ Foam Block Paper Other: \_\_\_\_\_

Was sufficient ice used (if appropriate)? \_\_\_\_\_

NA ☒ YES ☐ NO ☐

Were all bottles sealed in individual plastic bags? \_\_\_\_\_

YES ☒ NO ☐

Did all bottles arrive in good condition (unbroken)? \_\_\_\_\_

YES ☒ NO ☐

Were all bottle labels complete and legible? \_\_\_\_\_

YES ☒ NO ☐

Did the number of containers listed on COC match with the number of containers received? \_\_\_\_\_

YES ☒ NO ☐

Did all bottle labels and tags agree with custody papers? \_\_\_\_\_

YES ☒ NO ☐

Were all bottles used correct for the requested analyses? \_\_\_\_\_

YES ☒ NO ☐

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)...

NA ☒ YES ☐ NO ☐

Were all VOC vials free of air bubbles? \_\_\_\_\_

NA ☐ YES ☒ NO ☒

Was sufficient amount of sample sent in each bottle? \_\_\_\_\_

YES ☒ NO ☐

Date VOC Trip Blank was made at ARI: \_\_\_\_\_

NA 11/14/14

Was Sample Split by ARI: ☒ YES

Date/Time: \_\_\_\_\_

Equipment: \_\_\_\_\_

Split by: \_\_\_\_\_

Samples Logged by: AV

Date: 11/20/14

Time: 1047

**\*\* Notify Project Manager of discrepancies or concerns \*\***

| Sample ID on Bottle | Sample ID on COC | Sample ID on Bottle | Sample ID on COC |
|---------------------|------------------|---------------------|------------------|
|                     |                  |                     |                  |
|                     |                  |                     |                  |
|                     |                  |                     |                  |
|                     |                  |                     |                  |

### Additional Notes, Discrepancies, & Resolutions:

LMW-1 = 4sm LMW-2 = 2sm LMW-4 = 2pd LMW-5 = 4cg TB-1119 = 1sm  
3 TB for 1119 are not marked with client ID.

By: AV

Date: 11/20/14

|                                       |                                 |                                        |
|---------------------------------------|---------------------------------|----------------------------------------|
| <b>Small Air Bubbles</b><br>- 2mm<br> | <b>Peabubbles</b><br>2-4 mm<br> | <b>LARGE Air Bubbles</b><br>> 4 mm<br> |
|---------------------------------------|---------------------------------|----------------------------------------|

Small → "sm" (< 2 mm)

Peabubbles → "pb" (2 to < 4 mm)

Large → "lg" (4 to < 6 mm)

Headspace → "hs" (> 6 mm)

# Sample ID Cross Reference Report



ARI Job No: ZL38  
Client: Golder Associates  
Project Event: 923-1000-002.R273  
Project Name: Landsburg

| Sample ID             | ARI<br>Lab ID | ARI<br>LIMS ID | Matrix | Sample Date/Time | VTSR           |
|-----------------------|---------------|----------------|--------|------------------|----------------|
| 1. LMW-10-1114        | ZL38A         | 14-25246       | Water  | 11/18/14 09:45   | 11/19/14 14:42 |
| 2. LMW-2-1114         | ZL38B         | 14-25247       | Water  | 11/18/14 11:40   | 11/19/14 14:42 |
| 3. LMW-4-1114         | ZL38C         | 14-25248       | Water  | 11/18/14 13:10   | 11/19/14 14:42 |
| 4. LMW-3-1114         | ZL38D         | 14-25249       | Water  | 11/19/14 09:20   | 11/19/14 14:42 |
| 5. EB-1114            | ZL38E         | 14-25250       | Water  | 11/19/14 09:00   | 11/19/14 14:42 |
| 6. LMW-8-1114         | ZL38F         | 14-25251       | Water  | 11/19/14 10:05   | 11/19/14 14:42 |
| 7. LMW-5-1114         | ZL38G         | 14-25252       | Water  | 11/19/14 11:30   | 11/19/14 14:42 |
| 8. LMW-9-1114         | ZL38H         | 14-25253       | Water  | 11/19/14 13:15   | 11/19/14 14:42 |
| 9. TRIP BLANK-111814  | ZL38I         | 14-25254       | Water  | 11/18/14         | 11/19/14 14:42 |
| 10. TRIP BLANK-111914 | ZL38J         | 14-25255       | Water  | 11/19/14         | 11/19/14 14:42 |

# Sample ID Cross Reference Report



ARI Job No: ZL39  
Client: Golder Associates  
Project Event: 923-1000-002.R273  
Project Name: Landsburg

| Sample ID      | ARI<br>Lab ID | ARI<br>LIMS ID | Matrix | Sample Date/Time | VTSR           |
|----------------|---------------|----------------|--------|------------------|----------------|
| 1. LMW-10-1114 | ZL39A         | 14-25256       | Water  | 11/18/14 09:45   | 11/19/14 14:42 |
| 2. LMW-2-1114  | ZL39B         | 14-25257       | Water  | 11/18/14 11:40   | 11/19/14 14:42 |
| 3. LMW-4-1114  | ZL39C         | 14-25258       | Water  | 11/18/14 13:10   | 11/19/14 14:42 |
| 4. LMW-3-1114  | ZL39D         | 14-25259       | Water  | 11/19/14 09:20   | 11/19/14 14:42 |
| 5. EB-1114     | ZL39E         | 14-25260       | Water  | 11/19/14 09:00   | 11/19/14 14:42 |
| 6. LMW-8-1114  | ZL39F         | 14-25261       | Water  | 11/19/14 10:05   | 11/19/14 14:42 |
| 7. LMW-5-1114  | ZL39G         | 14-25262       | Water  | 11/19/14 11:30   | 11/19/14 14:42 |
| 8. LMW-9-1114  | ZL39H         | 14-25263       | Water  | 11/19/14 13:15   | 11/19/14 14:42 |

**PRESERVATION VERIFICATION 11/20/14**

Page 1 of 1



ARI Job No: **ZL38**

PC: Kelly  
VTSR: 11/19/14

Inquiry Number: NONE  
Analysis Requested: 11/20/14  
Contact: Morell, Douglas  
Client: Golder Associates  
Logged by: AV  
Sample Set Used: Yes-119  
Validatable Package: Lv4  
Deliverables:

Project #: 923-1000-002.R273  
Project: Landsburg  
Sample Site:  
SDG No:  
Analytical Protocol: In-house

| LOGNUM<br>ARI ID         | CLIENT ID   | CN<br>>12 | WAD<br>>12 | NH3<br><2 | COD<br><2 | FOG<br><2 | MET<br><2 | PHEN<br><2 | PHOS<br><2 | TKN<br><2 | NO23<br><2 | TOC<br><2 | S2<br>>9 | TPHD<br><2 | Fe2+<br><2 | DMET<br>FLT | DOC<br>FLT | PARAMETER | ADJUSTED<br>TO | LOT<br>NUMBER | AMOUNT<br>ADDED | DATE/BY |
|--------------------------|-------------|-----------|------------|-----------|-----------|-----------|-----------|------------|------------|-----------|------------|-----------|----------|------------|------------|-------------|------------|-----------|----------------|---------------|-----------------|---------|
| 14-25246<br><b>ZL38A</b> | LMW-10-1114 |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25247<br><b>ZL38B</b> | LMW-2-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25248<br><b>ZL38C</b> | LMW-4-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25249<br><b>ZL38D</b> | LMW-3-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25250<br><b>ZL38E</b> | EB-1114     |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25251<br><b>ZL38F</b> | LMW-8-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25252<br><b>ZL38G</b> | LMW-5-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25253<br><b>ZL38H</b> | LMW-9-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |

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Checked By AV Date 11/20/14

900000:8512

**PRESERVATION VERIFICATION 11/20/14**

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ARI Job No: ZL39

PC: Kelly  
VTSR: 11/19/14

Inquiry Number: NONE  
Analysis Requested: 11/20/14  
Contact: Morell, Douglas  
Client: Golder Associates  
Logged by: AV  
Sample Set Used: Yes-119  
Validatable Package: Lv4  
Deliverables:

Project #: 923-1000-002.R273  
Project: Landsburg  
Sample Site:  
SDG No:  
Analytical Protocol: In-house

| LOGNUM<br>ARI ID  | CLIENT ID   | CN<br>>12 | WAD<br>>12 | NH3<br><2 | COD<br><2 | FOG<br><2 | MET<br><2 | PHEN<br><2 | PHOS<br><2 | TKN<br><2 | NO23<br><2 | TOC<br><2 | S2<br>>9 | TPHD<br><2 | Fe2+<br><2 | DMET<br>FLT | DOC<br>FLT | PARAMETER | ADJUSTED<br>TO | LOT<br>NUMBER | AMOUNT<br>ADDED | DATE/BY |
|-------------------|-------------|-----------|------------|-----------|-----------|-----------|-----------|------------|------------|-----------|------------|-----------|----------|------------|------------|-------------|------------|-----------|----------------|---------------|-----------------|---------|
| 14-25256<br>ZL39A | LMW-10-1114 |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25257<br>ZL39B | LMW-2-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25258<br>ZL39C | LMW-4-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25259<br>ZL39D | LMW-3-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25260<br>ZL39E | EB-1114     |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25261<br>ZL39F | LMW-8-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25262<br>ZL39G | LMW-5-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25263<br>ZL39H | LMW-9-1114  |           |            |           |           |           | TOT       |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |

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Checked By AV Date 11/20/14

ZL39: 00007

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-10-1114

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SAMPLE

Lab Sample ID: ZL38A

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25246

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *mmw*

Date Sampled: 11/18/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 17:40

Purge Volume: 10.0 mL

| CAS Number     | Analyte                               | DL          | LOQ         | Result        |
|----------------|---------------------------------------|-------------|-------------|---------------|
| 74-87-3        | Chloromethane                         | 0.09        | 0.50        | < 0.50 U      |
| 74-83-9        | Bromomethane                          | 0.25        | 1.0         | < 1.0 U       |
| 75-01-4        | Vinyl Chloride                        | 0.06        | 0.10        | < 0.10 U      |
| 75-00-3        | Chloroethane                          | 0.09        | 0.20        | < 0.20 U      |
| 75-09-2        | Methylene Chloride                    | 0.48        | 1.0         | < 1.0 U       |
| 67-64-1        | Acetone                               | 2.1         | 5.0         | < 5.0 U       |
| <b>75-15-0</b> | <b>Carbon Disulfide</b>               | <b>0.04</b> | <b>0.20</b> | <b>0.14 J</b> |
| 75-35-4        | 1,1-Dichloroethene                    | 0.05        | 0.20        | < 0.20 U      |
| 75-34-3        | 1,1-Dichloroethane                    | 0.05        | 0.20        | < 0.20 U      |
| 156-60-5       | trans-1,2-Dichloroethene              | 0.05        | 0.20        | < 0.20 U      |
| 156-59-2       | cis-1,2-Dichloroethene                | 0.04        | 0.20        | < 0.20 U      |
| 67-66-3        | Chloroform                            | 0.03        | 0.20        | < 0.20 U      |
| 107-06-2       | 1,2-Dichloroethane                    | 0.07        | 0.20        | < 0.20 U      |
| 78-93-3        | 2-Butanone                            | 0.81        | 5.0         | < 5.0 U       |
| 71-55-6        | 1,1,1-Trichloroethane                 | 0.04        | 0.20        | < 0.20 U      |
| 56-23-5        | Carbon Tetrachloride                  | 0.04        | 0.20        | < 0.20 U      |
| 108-05-4       | Vinyl Acetate                         | 0.07        | 0.20        | < 0.20 U      |
| 75-27-4        | Bromodichloromethane                  | 0.05        | 0.20        | < 0.20 U      |
| 78-87-5        | 1,2-Dichloropropane                   | 0.04        | 0.20        | < 0.20 U      |
| 10061-01-5     | cis-1,3-Dichloropropene               | 0.06        | 0.20        | < 0.20 U      |
| 79-01-6        | Trichloroethene                       | 0.05        | 0.20        | < 0.20 U      |
| 124-48-1       | Dibromochloromethane                  | 0.05        | 0.20        | < 0.20 U      |
| 79-00-5        | 1,1,2-Trichloroethane                 | 0.13        | 0.20        | < 0.20 U      |
| 71-43-2        | Benzene                               | 0.03        | 0.20        | < 0.20 U      |
| 10061-02-6     | trans-1,3-Dichloropropene             | 0.08        | 0.20        | < 0.20 U      |
| 110-75-8       | 2-Chloroethylvinylether               | 0.25        | 0.50        | < 0.50 U      |
| 75-25-2        | Bromoform                             | 0.06        | 0.20        | < 0.20 U      |
| 108-10-1       | 4-Methyl-2-Pentanone (MIBK)           | 0.97        | 2.5         | < 2.5 U       |
| 591-78-6       | 2-Hexanone                            | 0.90        | 5.0         | < 5.0 U       |
| 127-18-4       | Tetrachloroethene                     | 0.05        | 0.20        | < 0.20 U      |
| 79-34-5        | 1,1,2,2-Tetrachloroethane             | 0.06        | 0.10        | < 0.10 U      |
| 108-88-3       | Toluene                               | 0.04        | 0.20        | < 0.20 U      |
| 108-90-7       | Chlorobenzene                         | 0.02        | 0.20        | < 0.20 U      |
| 100-41-4       | Ethylbenzene                          | 0.04        | 0.20        | < 0.20 U      |
| 100-42-5       | Styrene                               | 0.05        | 0.20        | < 0.20 U      |
| 75-69-4        | Trichlorofluoromethane                | 0.04        | 0.20        | < 0.20 U      |
| 76-13-1        | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04        | 0.20        | < 0.20 U      |
| 179601-23-1    | m,p-Xylene                            | 0.05        | 0.40        | < 0.40 U      |
| 95-47-6        | o-Xylene                              | 0.03        | 0.20        | < 0.20 U      |
| 95-50-1        | 1,2-Dichlorobenzene                   | 0.04        | 0.20        | < 0.20 U      |
| 541-73-1       | 1,3-Dichlorobenzene                   | 0.04        | 0.20        | < 0.20 U      |
| 106-46-7       | 1,4-Dichlorobenzene                   | 0.04        | 0.20        | < 0.20 U      |
| 107-02-8       | Acrolein                              | 2.5         | 2.5         | < 2.5 U       |
| 74-88-4        | Iodomethane                           | 0.23        | 0.50        | < 0.50 U      |
| 107-13-1       | Acrylonitrile                         | 0.60        | 1.0         | < 1.0 U       |
| 563-58-6       | 1,1-Dichloropropene                   | 0.03        | 0.10        | < 0.10 U      |
| 74-95-3        | Dibromomethane                        | 0.14        | 0.20        | < 0.20 U      |
| 630-20-6       | 1,1,1,2-Tetrachloroethane             | 0.04        | 0.20        | < 0.20 U      |
| 96-12-8        | 1,2-Dibromo-3-chloropropane           | 0.04        | 0.50        | < 0.50 U      |
| 96-18-4        | 1,2,3-Trichloropropane                | 0.13        | 0.20        | < 0.20 U      |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-10-1114  
SAMPLE

Lab Sample ID: ZL38A

LIMS ID: 14-25246

Matrix: Water

Date Analyzed: 11/28/14 17:40

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | 0.33 B   |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 102%  |
| d8-Toluene             | 100%  |
| Bromofluorobenzene     | 96.5% |
| d4-1,2-Dichlorobenzene | 101%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-2-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38B

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25247

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MMW*

Date Sampled: 11/18/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 18:07

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-2-1114  
SAMPLE

Lab Sample ID: ZL38B

LIMS ID: 14-25247

Matrix: Water

Date Analyzed: 11/28/14 18:07

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 105%  |
| d8-Toluene             | 98.0% |
| Bromofluorobenzene     | 96.5% |
| d4-1,2-Dichlorobenzene | 103%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-4-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38C

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25248

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *[Signature]*

Date Sampled: 11/18/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 18:34

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-4-1114

Page 2 of 2

SAMPLE

Lab Sample ID: ZL38C

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25248

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 18:34

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 106%  |
| d8-Toluene             | 97.8% |
| Bromofluorobenzene     | 98.1% |
| d4-1,2-Dichlorobenzene | 106%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-3-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38D

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25249

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *[Signature]*

Date Sampled: 11/19/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 19:00

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-3-1114  
SAMPLE

Lab Sample ID: ZL38D

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25249

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 19:00

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 103%  |
| d8-Toluene             | 102%  |
| Bromofluorobenzene     | 95.1% |
| d4-1,2-Dichlorobenzene | 105%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 2Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E

LIMS ID: 14-25250

Matrix: Water

Data Release Authorized: *mm*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Date Analyzed: 11/28/14 19:27

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | 1.4 B    |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: EB-1114

Page 2 of 2

SAMPLE

Lab Sample ID: ZL38E

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25250

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 19:27

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 104%  |
| d8-Toluene             | 98.6% |
| Bromofluorobenzene     | 95.7% |
| d4-1,2-Dichlorobenzene | 104%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-8-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38F

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25251

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *mmw*

Date Sampled: 11/19/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 19:53

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25251

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 19:53

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 104%  |
| d8-Toluene             | 102%  |
| Bromofluorobenzene     | 94.7% |
| d4-1,2-Dichlorobenzene | 103%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 2Sample ID: LMW-5-1114  
SAMPLE

Lab Sample ID: ZL38G

LIMS ID: 14-25252

Matrix: Water

Data Release Authorized: *MMW*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Date Analyzed: 11/28/14 20:20

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-5-1114

Page 2 of 2

SAMPLE

Lab Sample ID: ZL38G

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25252

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 20:20

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 108%  |
| d8-Toluene             | 99.6% |
| Bromofluorobenzene     | 95.8% |
| d4-1,2-Dichlorobenzene | 104%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-9-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38H

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25253

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/19/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 20:47

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-9-1114  
SAMPLE

Lab Sample ID: ZL38H

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25253

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 20:47

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 107%  |
| d8-Toluene             | 100%  |
| Bromofluorobenzene     | 94.9% |
| d4-1,2-Dichlorobenzene | 105%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: TRIP BLANK-111814

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38I

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25254

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *mw*

Date Sampled: 11/18/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 21:14

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | 2.3 B    |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: TRIP BLANK-111814

Page 2 of 2

SAMPLE

Lab Sample ID: ZL38I

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25254

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 21:14

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 108%  |
| d8-Toluene             | 100%  |
| Bromofluorobenzene     | 94.2% |
| d4-1,2-Dichlorobenzene | 105%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: MB-112814A

Page 1 of 2

METHOD BLANK

Lab Sample ID: MB-112814A

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25246

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: 

Date Sampled: NA

Reported: 12/03/14

Date Received: NA

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 17:13

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | 0.92 J   |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Page 2 of 2

Sample ID: MB-112814A

METHOD BLANK

Lab Sample ID: MB-112814A

LIMS ID: 14-25246

Matrix: Water

Date Analyzed: 11/28/14 17:13

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | 0.74     |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | 0.11 J   |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | 0.21 J   |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | 0.26     |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 103%  |
| d8-Toluene             | 100%  |
| Bromofluorobenzene     | 95.1% |
| d4-1,2-Dichlorobenzene | 102%  |

## VOA SURROGATE RECOVERY SUMMARY

Matrix: Water

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| ARI ID       | Client ID         | PV | DCE   | TOL   | BFB   | DCB   | TOT OUT |
|--------------|-------------------|----|-------|-------|-------|-------|---------|
| MB-112814A   | Method Blank      | 10 | 103%  | 100%  | 95.1% | 102%  | 0       |
| LCS-112814A  | Lab Control       | 10 | 97.8% | 99.1% | 104%  | 99.4% | 0       |
| LCSD-112814A | Lab Control Dup   | 10 | 98.7% | 99.6% | 102%  | 100%  | 0       |
| ZL38A        | LMW-10-1114       | 10 | 102%  | 100%  | 96.5% | 101%  | 0       |
| ZL38B        | LMW-2-1114        | 10 | 105%  | 98.0% | 96.5% | 103%  | 0       |
| ZL38C        | LMW-4-1114        | 10 | 106%  | 97.8% | 98.1% | 106%  | 0       |
| ZL38D        | LMW-3-1114        | 10 | 103%  | 102%  | 95.1% | 105%  | 0       |
| ZL38E        | EB-1114           | 10 | 104%  | 98.6% | 95.7% | 104%  | 0       |
| ZL38F        | LMW-8-1114        | 10 | 104%  | 102%  | 94.7% | 103%  | 0       |
| ZL38G        | LMW-5-1114        | 10 | 108%  | 99.6% | 95.8% | 104%  | 0       |
| ZL38H        | LMW-9-1114        | 10 | 107%  | 100%  | 94.9% | 105%  | 0       |
| ZL38I        | TRIP BLANK-111814 | 10 | 108%  | 100%  | 94.2% | 105%  | 0       |

## LCS/MB LIMITS

## QC LIMITS

## SW8260C

|                                |          |          |
|--------------------------------|----------|----------|
| (DCE) = d4-1,2-Dichloroethane  | (80-120) | (80-120) |
| (TOL) = d8-Toluene             | (80-120) | (80-120) |
| (BFB) = Bromofluorobenzene     | (80-120) | (80-120) |
| (DCB) = d4-1,2-Dichlorobenzene | (80-120) | (80-120) |

Prep Method: SW5030B  
Log Number Range: 14-25246 to 14-25254

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Page 1 of 2

Sample ID: LCS-112814A

LAB CONTROL SAMPLE

Lab Sample ID: LCS-112814A

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: *mw*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: NA

Date Received: NA

Instrument/Analyst LCS: NT2/LH

LCSD: NT2/LH

Date Analyzed LCS: 11/28/14 15:42

LCSD: 11/28/14 16:09

Sample Amount LCS: 10.0 mL

LCSD: 10.0 mL

Purge Volume LCS: 10.0 mL

LCSD: 10.0 mL

| Analyte                             | LCS    | Spike<br>Added-LCS | LCS<br>Recovery | LCSD   | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD  |
|-------------------------------------|--------|--------------------|-----------------|--------|---------------------|------------------|------|
| Chloromethane                       | 10.2   | 10.0               | 102%            | 10.0   | 10.0                | 100%             | 2.0% |
| Bromomethane                        | 10.6   | 10.0               | 106%            | 10.2   | 10.0                | 102%             | 3.8% |
| Vinyl Chloride                      | 10.9   | 10.0               | 109%            | 10.6   | 10.0                | 106%             | 2.8% |
| Chloroethane                        | 11.5   | 10.0               | 115%            | 11.2   | 10.0                | 112%             | 2.6% |
| Methylene Chloride                  | 10.9 B | 10.0               | 109%            | 10.8 B | 10.0                | 108%             | 0.9% |
| Acetone                             | 48.4   | 50.0               | 96.8%           | 47.9   | 50.0                | 95.8%            | 1.0% |
| Carbon Disulfide                    | 10.4   | 10.0               | 104%            | 9.99   | 10.0                | 99.9%            | 4.0% |
| 1,1-Dichloroethene                  | 10.4   | 10.0               | 104%            | 10.0   | 10.0                | 100%             | 3.9% |
| 1,1-Dichloroethane                  | 10.6   | 10.0               | 106%            | 10.4   | 10.0                | 104%             | 1.9% |
| trans-1,2-Dichloroethene            | 10.8   | 10.0               | 108%            | 10.4   | 10.0                | 104%             | 3.8% |
| cis-1,2-Dichloroethene              | 10.4   | 10.0               | 104%            | 10.2   | 10.0                | 102%             | 1.9% |
| Chloroform                          | 10.5   | 10.0               | 105%            | 10.3   | 10.0                | 103%             | 1.9% |
| 1,2-Dichloroethane                  | 10.4   | 10.0               | 104%            | 10.4   | 10.0                | 104%             | 0.0% |
| 2-Butanone                          | 51.3   | 50.0               | 103%            | 53.7   | 50.0                | 107%             | 4.6% |
| 1,1,1-Trichloroethane               | 10.7   | 10.0               | 107%            | 10.4   | 10.0                | 104%             | 2.8% |
| Carbon Tetrachloride                | 10.9   | 10.0               | 109%            | 10.5   | 10.0                | 105%             | 3.7% |
| Vinyl Acetate                       | 8.92   | 10.0               | 89.2%           | 9.17   | 10.0                | 91.7%            | 2.8% |
| Bromodichloromethane                | 10.3   | 10.0               | 103%            | 10.9   | 10.0                | 109%             | 5.7% |
| 1,2-Dichloropropane                 | 10.7   | 10.0               | 107%            | 10.7   | 10.0                | 107%             | 0.0% |
| cis-1,3-Dichloropropene             | 9.82   | 10.0               | 98.2%           | 10.1   | 10.0                | 101%             | 2.8% |
| Trichloroethene                     | 11.0   | 10.0               | 110%            | 10.8   | 10.0                | 108%             | 1.8% |
| Dibromochloromethane                | 8.80   | 10.0               | 88.0%           | 8.94   | 10.0                | 89.4%            | 1.6% |
| 1,1,2-Trichloroethane               | 10.0   | 10.0               | 100%            | 10.4   | 10.0                | 104%             | 3.9% |
| Benzene                             | 10.9   | 10.0               | 109%            | 10.8   | 10.0                | 108%             | 0.9% |
| trans-1,3-Dichloropropene           | 11.2   | 10.0               | 112%            | 11.6   | 10.0                | 116%             | 3.5% |
| 2-Chloroethylvinylether             | 8.66   | 10.0               | 86.6%           | 9.53   | 10.0                | 95.3%            | 9.6% |
| Bromoform                           | 10.8   | 10.0               | 108%            | 11.0   | 10.0                | 110%             | 1.8% |
| 4-Methyl-2-Pentanone (MIBK)         | 50.6   | 50.0               | 101%            | 52.0   | 50.0                | 104%             | 2.7% |
| 2-Hexanone                          | 49.4   | 50.0               | 98.8%           | 48.7   | 50.0                | 97.4%            | 1.4% |
| Tetrachloroethene                   | 10.9   | 10.0               | 109%            | 10.6   | 10.0                | 106%             | 2.8% |
| 1,1,2,2-Tetrachloroethane           | 10.6   | 10.0               | 106%            | 10.7   | 10.0                | 107%             | 0.9% |
| Toluene                             | 10.5   | 10.0               | 105%            | 10.4   | 10.0                | 104%             | 1.0% |
| Chlorobenzene                       | 10.8   | 10.0               | 108%            | 10.6   | 10.0                | 106%             | 1.9% |
| Ethylbenzene                        | 11.1   | 10.0               | 111%            | 10.9   | 10.0                | 109%             | 1.8% |
| Styrene                             | 10.3   | 10.0               | 103%            | 10.4   | 10.0                | 104%             | 1.0% |
| Trichlorofluoromethane              | 11.9   | 10.0               | 119%            | 11.1   | 10.0                | 111%             | 7.0% |
| 1,1,2-Trichloro-1,2,2-trifluoroetha | 10.5   | 10.0               | 105%            | 10.4   | 10.0                | 104%             | 1.0% |
| m,p-Xylene                          | 23.0   | 20.0               | 115%            | 22.5   | 20.0                | 112%             | 2.2% |

**ORGANICS ANALYSIS DATA SHEET**

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2

Sample ID: LCS-112814A

LAB CONTROL SAMPLE

Lab Sample ID: LCS-112814A  
LIMS ID: 14-25246  
Matrix: Water

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| Analyte                     | LCS    | Spike<br>Added-LCS | LCS<br>Recovery | LCSD   | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|-----------------------------|--------|--------------------|-----------------|--------|---------------------|------------------|-------|
| o-Xylene                    | 10.1   | 10.0               | 101%            | 9.79   | 10.0                | 97.9%            | 3.1%  |
| 1,2-Dichlorobenzene         | 10.5   | 10.0               | 105%            | 10.5   | 10.0                | 105%             | 0.0%  |
| 1,3-Dichlorobenzene         | 11.1   | 10.0               | 111%            | 11.0   | 10.0                | 110%             | 0.9%  |
| 1,4-Dichlorobenzene         | 10.7   | 10.0               | 107%            | 10.7   | 10.0                | 107%             | 0.0%  |
| Acrolein                    | 53.9   | 50.0               | 108%            | 54.1   | 50.0                | 108%             | 0.4%  |
| Iodomethane                 | 10.4   | 10.0               | 104%            | 10.0   | 10.0                | 100%             | 3.9%  |
| Acrylonitrile               | 9.12   | 10.0               | 91.2%           | 10.3   | 10.0                | 103%             | 12.2% |
| 1,1-Dichloropropene         | 10.4   | 10.0               | 104%            | 10.4   | 10.0                | 104%             | 0.0%  |
| Dibromomethane              | 10.3   | 10.0               | 103%            | 10.4   | 10.0                | 104%             | 1.0%  |
| 1,1,1,2-Tetrachloroethane   | 10.8   | 10.0               | 108%            | 10.5   | 10.0                | 105%             | 2.8%  |
| 1,2-Dibromo-3-chloropropane | 9.26   | 10.0               | 92.6%           | 9.21   | 10.0                | 92.1%            | 0.5%  |
| 1,2,3-Trichloropropane      | 9.97   | 10.0               | 99.7%           | 10.3   | 10.0                | 103%             | 3.3%  |
| trans-1,4-Dichloro-2-butene | 10.9   | 10.0               | 109%            | 11.2   | 10.0                | 112%             | 2.7%  |
| 1,3,5-Trimethylbenzene      | 11.8   | 10.0               | 118%            | 11.8   | 10.0                | 118%             | 0.0%  |
| 1,2,4-Trimethylbenzene      | 11.7   | 10.0               | 117%            | 11.6   | 10.0                | 116%             | 0.9%  |
| Hexachlorobutadiene         | 11.6 B | 10.0               | 116%            | 11.3 B | 10.0                | 113%             | 2.6%  |
| 1,2-Dibromoethane           | 11.1   | 10.0               | 111%            | 11.5   | 10.0                | 115%             | 3.5%  |
| Bromochloromethane          | 10.1   | 10.0               | 101%            | 10.1   | 10.0                | 101%             | 0.0%  |
| 2,2-Dichloropropane         | 11.8   | 10.0               | 118%            | 11.3   | 10.0                | 113%             | 4.3%  |
| 1,3-Dichloropropane         | 10.8   | 10.0               | 108%            | 10.9   | 10.0                | 109%             | 0.9%  |
| Isopropylbenzene            | 9.38   | 10.0               | 93.8%           | 9.27   | 10.0                | 92.7%            | 1.2%  |
| n-Propylbenzene             | 11.7   | 10.0               | 117%            | 11.7   | 10.0                | 117%             | 0.0%  |
| Bromobenzene                | 10.8   | 10.0               | 108%            | 10.6   | 10.0                | 106%             | 1.9%  |
| 2-Chlorotoluene             | 11.8   | 10.0               | 118%            | 11.8   | 10.0                | 118%             | 0.0%  |
| 4-Chlorotoluene             | 11.7   | 10.0               | 117%            | 11.6   | 10.0                | 116%             | 0.9%  |
| tert-Butylbenzene           | 10.9   | 10.0               | 109%            | 10.8   | 10.0                | 108%             | 0.9%  |
| sec-Butylbenzene            | 9.30   | 10.0               | 93.0%           | 9.18   | 10.0                | 91.8%            | 1.3%  |
| 4-Isopropyltoluene          | 11.6   | 10.0               | 116%            | 11.5   | 10.0                | 115%             | 0.9%  |
| n-Butylbenzene              | 11.8 B | 10.0               | 118%            | 11.6 B | 10.0                | 116%             | 1.7%  |
| 1,2,4-Trichlorobenzene      | 11.3   | 10.0               | 113%            | 11.2   | 10.0                | 112%             | 0.9%  |
| Naphthalene                 | 9.01 B | 10.0               | 90.1%           | 9.48 B | 10.0                | 94.8%            | 5.1%  |
| 1,2,3-Trichlorobenzene      | 10.9 B | 10.0               | 109%            | 11.4 B | 10.0                | 114%             | 4.5%  |

Reported in µg/L (ppb)

RPD calculated using sample concentrations per SW846.

**Volatile Surrogate Recovery**

|                        | LCS   | LCSD  |
|------------------------|-------|-------|
| d4-1,2-Dichloroethane  | 97.8% | 98.7% |
| d8-Toluene             | 99.1% | 99.6% |
| Bromofluorobenzene     | 104%  | 102%  |
| d4-1,2-Dichlorobenzene | 99.4% | 100%  |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-10-1114  
SAMPLE

Lab Sample ID: ZL38A

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: *MB*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/24/14

Date Analyzed: 11/28/14 17:17

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-10-1114  
SAMPLE

Lab Sample ID: ZL38A  
LIMS ID: 14-25246  
Matrix: Water  
Date Analyzed: 11/28/14 17:17

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 78.4% | 2-Fluorobiphenyl       | 83.2% |
| d14-p-Terphenyl      | 86.0% | d4-1,2-Dichlorobenzene | 65.2% |
| d5-Phenol            | 83.7% | 2-Fluorophenol         | 69.6% |
| 2,4,6-Tribromophenol | 86.1% | d4-2-Chlorophenol      | 76.8% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-2-1114  
SAMPLE

Lab Sample ID: ZL38B

LIMS ID: 14-25247

Matrix: Water

Data Release Authorized: *AB*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/24/14

Date Analyzed: 12/02/14 19:49

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-2-1114  
SAMPLE

Lab Sample ID: ZL38B  
LIMS ID: 14-25247  
Matrix: Water  
Date Analyzed: 12/02/14 19:49

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 48.4% | 2-Fluorobiphenyl       | 48.4% |
| d14-p-Terphenyl      | 53.6% | d4-1,2-Dichlorobenzene | 47.2% |
| d5-Phenol            | 58.7% | 2-Fluorophenol         | 51.2% |
| 2,4,6-Tribromophenol | 49.9% | d4-2-Chlorophenol      | 54.4% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-4-1114  
SAMPLE

Lab Sample ID: ZL38C  
LIMS ID: 14-25248  
Matrix: Water  
Data Release Authorized: *A*  
Reported: 12/03/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/18/14  
Date Received: 11/19/14

Date Extracted: 11/24/14  
Date Analyzed: 11/28/14 18:25  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-4-1114  
SAMPLE

Lab Sample ID: ZL38C  
LIMS ID: 14-25248  
Matrix: Water  
Date Analyzed: 11/28/14 18:25

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | 12      |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 72.8% | 2-Fluorobiphenyl       | 68.8% |
| d14-p-Terphenyl      | 72.4% | d4-1,2-Dichlorobenzene | 63.2% |
| d5-Phenol            | 77.3% | 2-Fluorophenol         | 69.3% |
| 2,4,6-Tribromophenol | 69.6% | d4-2-Chlorophenol      | 71.5% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-3-1114  
SAMPLE

Lab Sample ID: ZL38D  
LIMS ID: 14-25249  
Matrix: Water  
Data Release Authorized: *RB*  
Reported: 12/03/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/24/14  
Date Analyzed: 11/28/14 18:59  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-3-1114  
SAMPLE

Lab Sample ID: ZL38D  
LIMS ID: 14-25249  
Matrix: Water  
Date Analyzed: 11/28/14 18:59

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 66.8% | 2-Fluorobiphenyl       | 64.4% |
| d14-p-Terphenyl      | 70.0% | d4-1,2-Dichlorobenzene | 60.8% |
| d5-Phenol            | 69.9% | 2-Fluorophenol         | 61.1% |
| 2,4,6-Tribromophenol | 63.7% | d4-2-Chlorophenol      | 65.1% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E  
LIMS ID: 14-25250  
Matrix: Water  
Data Release Authorized:   
Reported: 12/03/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/24/14  
Date Analyzed: 11/28/14 19:33  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E  
LIMS ID: 14-25250  
Matrix: Water  
Date Analyzed: 11/28/14 19:33

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 74.0% | 2-Fluorobiphenyl       | 71.2% |
| d14-p-Terphenyl      | 81.2% | d4-1,2-Dichlorobenzene | 66.0% |
| d5-Phenol            | 76.8% | 2-Fluorophenol         | 68.0% |
| 2,4,6-Tribromophenol | 69.3% | d4-2-Chlorophenol      | 72.8% |

**ORGANICS ANALYSIS DATA SHEET**  
**Semivolatiles by SW8270D GC/MS**  
**Extraction Method: SW3520C**  
 Page 1 of 2

**Sample ID: LMW-8-1114**  
**SAMPLE**

Lab Sample ID: ZL38F

LIMS ID: 14-25251

Matrix: Water

Data Release Authorized: *AB*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/24/14

Date Analyzed: 11/28/14 20:06

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F  
LIMS ID: 14-25251  
Matrix: Water  
Date Analyzed: 11/28/14 20:06

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 63.6% | 2-Fluorobiphenyl       | 60.8% |
| d14-p-Terphenyl      | 68.8% | d4-1,2-Dichlorobenzene | 56.0% |
| d5-Phenol            | 67.5% | 2-Fluorophenol         | 59.2% |
| 2,4,6-Tribromophenol | 62.1% | d4-2-Chlorophenol      | 61.9% |

Sample ID: LMW-5-1114  
SAMPLE



Lab Sample ID: ZL38G  
LIMS ID: 14-25252  
Matrix: Water  
Data Release Authorized:   
Reported: 02/04/15

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/24/14  
Date Analyzed: 11/28/14 20:40  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

Sample ID: LMW-5-1114  
SAMPLE

Lab Sample ID: ZL38G  
LIMS ID: 14-25252  
Matrix: Water  
Date Analyzed: 11/28/14 20:40

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 54.4% | 2-Fluorobiphenyl       | 54.8% |
| d14-p-Terphenyl      | 55.6% | d4-1,2-Dichlorobenzene | 46.8% |
| d5-Phenol            | 57.6% | 2-Fluorophenol         | 50.9% |
| 2,4,6-Tribromophenol | 53.6% | d4-2-Chlorophenol      | 53.6% |

Lab Sample ID: ZL38H

LIMS ID: 14-25253

Matrix: Water

Data Release Authorized:

Reported: 02/04/15

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/24/14

Date Analyzed: 11/28/14 21:14

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-9-1114  
SAMPLE



Lab Sample ID: ZL38H  
LIMS ID: 14-25253  
Matrix: Water  
Date Analyzed: 11/28/14 21:14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 70.8% | 2-Fluorobiphenyl       | 63.2% |
| d14-p-Terphenyl      | 65.2% | d4-1,2-Dichlorobenzene | 57.6% |
| d5-Phenol            | 64.0% | 2-Fluorophenol         | 56.5% |
| 2,4,6-Tribromophenol | 65.9% | d4-2-Chlorophenol      | 58.9% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: MB-112414  
METHOD BLANK

Lab Sample ID: MB-112414  
LIMS ID: 14-25246  
Matrix: Water  
Data Release Authorized: *AB*  
Reported: 12/03/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: NA  
Date Received: NA

Date Extracted: 11/24/14  
Date Analyzed: 11/28/14 11:04  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: MB-112414  
METHOD BLANK

Lab Sample ID: MB-112414  
LIMS ID: 14-25246  
Matrix: Water  
Date Analyzed: 11/28/14 11:04

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 78.0% | 2-Fluorobiphenyl       | 73.2% |
| d14-p-Terphenyl      | 78.0% | d4-1,2-Dichlorobenzene | 68.8% |
| d5-Phenol            | 82.9% | 2-Fluorophenol         | 72.8% |
| 2,4,6-Tribromophenol | 71.5% | d4-2-Chlorophenol      | 76.3% |

**SW8270 SEMIVOLATILES WATER SURROGATE RECOVERY SUMMARY**

Matrix: Water

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

| Client ID   | NBZ   | FBP   | TPH   | DCB   | PHL   | 2FP   | TBP    | 2CP   | TOT | OUT |
|-------------|-------|-------|-------|-------|-------|-------|--------|-------|-----|-----|
| MB-112414   | 78.0% | 73.2% | 78.0% | 68.8% | 82.9% | 72.8% | 71.5%  | 76.3% |     | 0   |
| LCS-112414  | 74.8% | 69.6% | 76.4% | 63.6% | 79.7% | 66.1% | 81.9%  | 72.3% |     | 0   |
| LCSD-112414 | 78.0% | 71.6% | 76.0% | 59.2% | 74.1% | 72.8% | 90.7%  | 69.1% |     | 0   |
| LMW-10-1114 | 78.4% | 83.2% | 86.0% | 65.2% | 83.7% | 69.6% | 86.1%  | 76.8% |     | 0   |
| LMW-2-1114  | 48.4% | 48.4% | 53.6% | 47.2% | 58.7% | 51.2% | 49.9%* | 54.4% |     | 1   |
| LMW-4-1114  | 72.8% | 68.8% | 72.4% | 63.2% | 77.3% | 69.3% | 69.6%  | 71.5% |     | 0   |
| LMW-3-1114  | 66.8% | 64.4% | 70.0% | 60.8% | 69.9% | 61.1% | 63.7%  | 65.1% |     | 0   |
| EB-1114     | 74.0% | 71.2% | 81.2% | 66.0% | 76.8% | 68.0% | 69.3%  | 72.8% |     | 0   |
| LMW-8-1114  | 63.6% | 60.8% | 68.8% | 56.0% | 67.5% | 59.2% | 62.1%  | 61.9% |     | 0   |
| LMW-5-1114  | 54.4% | 54.8% | 55.6% | 46.8% | 57.6% | 50.9% | 53.6%  | 53.6% |     | 0   |
| LMW-9-1114  | 70.8% | 63.2% | 65.2% | 57.6% | 64.0% | 56.5% | 65.9%  | 58.9% |     | 0   |

|                                | LCS/MB LIMITS | QC LIMITS |
|--------------------------------|---------------|-----------|
| (NBZ) = d5-Nitrobenzene        | (27-120)      | (27-120)  |
| (FBP) = 2-Fluorobiphenyl       | (33-120)      | (33-120)  |
| (TPH) = d14-p-Terphenyl        | (28-130)      | (28-130)  |
| (DCB) = d4-1,2-Dichlorobenzene | (20-120)      | (20-120)  |
| (PHL) = d5-Phenol              | (38-120)      | (38-120)  |
| (2FP) = 2-Fluorophenol         | (33-120)      | (33-120)  |
| (TBP) = 2,4,6-Tribromophenol   | (52-131)      | (52-131)  |
| (2CP) = d4-2-Chlorophenol      | (41-120)      | (41-120)  |

Prep Method: SW3520C  
Log Number Range: 14-25246 to 14-25253

**ORGANICS ANALYSIS DATA SHEET**  
Semivolatiles by SW8270D GC/MS  
Page 1 of 3

Sample ID: LCS-112414  
LCS/LCSD

Lab Sample ID: LCS-112414

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: 

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted LCS/LCSD: 11/24/14

Sample Amount LCS: 500 mL

LCSD: 500 mL

Date Analyzed LCS: 11/28/14 11:37

Final Extract Volume LCS: 0.50 mL

LCSD: 11/28/14 12:11

LCSD: 0.50 mL

Instrument/Analyst LCS: NT6/JZ

Dilution Factor LCS: 1.00

LCSD: NT6/JZ

LCSD: 1.00

GPC Cleanup: NO

| Analyte                      | LCS  | Spike<br>Added-LCS | LCS<br>Recovery | LCSD | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|------------------------------|------|--------------------|-----------------|------|---------------------|------------------|-------|
| Phenol                       | 19.2 | 25.0               | 76.8%           | 19.4 | 25.0                | 77.6%            | 1.0%  |
| Bis-(2-Chloroethyl) Ether    | 17.6 | 25.0               | 70.4%           | 18.6 | 25.0                | 74.4%            | 5.5%  |
| 2-Chlorophenol               | 18.1 | 25.0               | 72.4%           | 17.2 | 25.0                | 68.8%            | 5.1%  |
| 1,3-Dichlorobenzene          | 13.8 | 25.0               | 55.2%           | 13.4 | 25.0                | 53.6%            | 2.9%  |
| 1,4-Dichlorobenzene          | 14.4 | 25.0               | 57.6%           | 14.4 | 25.0                | 57.6%            | 0.0%  |
| Benzyl Alcohol               | 22.1 | 25.0               | 88.4%           | 20.9 | 25.0                | 83.6%            | 5.6%  |
| 1,2-Dichlorobenzene          | 15.5 | 25.0               | 62.0%           | 14.4 | 25.0                | 57.6%            | 7.4%  |
| 2-Methylphenol               | 18.1 | 25.0               | 72.4%           | 18.6 | 25.0                | 74.4%            | 2.7%  |
| 2,2'-Oxybis(1-Chloropropane) | 17.7 | 25.0               | 70.8%           | 18.6 | 25.0                | 74.4%            | 5.0%  |
| 4-Methylphenol               | 19.2 | 25.0               | 76.8%           | 18.3 | 25.0                | 73.2%            | 4.8%  |
| N-Nitroso-Di-N-Propylamine   | 19.3 | 25.0               | 77.2%           | 19.6 | 25.0                | 78.4%            | 1.5%  |
| Hexachloroethane             | 13.8 | 25.0               | 55.2%           | 13.4 | 25.0                | 53.6%            | 2.9%  |
| Nitrobenzene                 | 18.9 | 25.0               | 75.6%           | 20.7 | 25.0                | 82.8%            | 9.1%  |
| Isophorone                   | 20.0 | 25.0               | 80.0%           | 21.9 | 25.0                | 87.6%            | 9.1%  |
| 2-Nitrophenol                | 20.2 | 25.0               | 80.8%           | 21.8 | 25.0                | 87.2%            | 7.6%  |
| 2,4-Dimethylphenol           | 63.0 | 75.0               | 84.0%           | 65.7 | 75.0                | 87.6%            | 4.2%  |
| Benzoic Acid                 | 134  | 138                | 97.1%           | 144  | 138                 | 104%             | 7.2%  |
| bis(2-Chloroethoxy) Methane  | 18.5 | 25.0               | 74.0%           | 19.4 | 25.0                | 77.6%            | 4.7%  |
| 2,4-Dichlorophenol           | 61.8 | 75.0               | 82.4%           | 68.3 | 75.0                | 91.1%            | 10.0% |
| 1,2,4-Trichlorobenzene       | 15.5 | 25.0               | 62.0%           | 16.6 | 25.0                | 66.4%            | 6.9%  |
| Naphthalene                  | 17.0 | 25.0               | 68.0%           | 18.8 | 25.0                | 75.2%            | 10.1% |
| 4-Chloroaniline              | 56.6 | 75.0               | 75.5%           | 68.2 | 75.0                | 90.9%            | 18.6% |
| Hexachlorobutadiene          | 13.6 | 25.0               | 54.4%           | 15.7 | 25.0                | 62.8%            | 14.3% |
| 4-Chloro-3-methylphenol      | 73.3 | 75.0               | 97.7%           | 77.0 | 75.0                | 103%             | 4.9%  |
| 2-Methylnaphthalene          | 19.6 | 25.0               | 78.4%           | 22.1 | 25.0                | 88.4%            | 12.0% |
| Hexachlorocyclopentadiene    | 47.1 | 75.0               | 62.8%           | 52.7 | 75.0                | 70.3%            | 11.2% |
| 2,4,6-Trichlorophenol        | 65.2 | 75.0               | 86.9%           | 68.1 | 75.0                | 90.8%            | 4.4%  |
| 2,4,5-Trichlorophenol        | 64.9 | 75.0               | 86.5%           | 70.6 | 75.0                | 94.1%            | 8.4%  |
| 2-Chloronaphthalene          | 19.1 | 25.0               | 76.4%           | 17.8 | 25.0                | 71.2%            | 7.0%  |
| 2-Nitroaniline               | 75.4 | 75.0               | 101%            | 74.2 | 75.0                | 98.9%            | 1.6%  |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Page 2 of 3

Sample ID: LCS-112414  
LCS/LCSD

Lab Sample ID: LCS-112414  
LIMS ID: 14-25246  
Matrix: Water  
Date Analyzed LCS: 11/28/14 11:37  
LCSD: 11/28/14 12:11

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| Analyte                    | LCS  | Spike<br>Added-LCS | LCS<br>Recovery | LCSD | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|----------------------------|------|--------------------|-----------------|------|---------------------|------------------|-------|
| Dimethylphthalate          | 21.0 | 25.0               | 84.0%           | 21.5 | 25.0                | 86.0%            | 2.4%  |
| Acenaphthylene             | 20.5 | 25.0               | 82.0%           | 20.6 | 25.0                | 82.4%            | 0.5%  |
| 3-Nitroaniline             | 76.6 | 75.0               | 102%            | 82.8 | 75.0                | 110%             | 7.8%  |
| Acenaphthene               | 20.3 | 25.0               | 81.2%           | 20.4 | 25.0                | 81.6%            | 0.5%  |
| 2,4-Dinitrophenol          | 121  | 138                | 87.7%           | 150  | 138                 | 109%             | 21.4% |
| 4-Nitrophenol              | 65.9 | 75.0               | 87.9%           | 68.3 | 75.0                | 91.1%            | 3.6%  |
| Dibenzofuran               | 21.3 | 25.0               | 85.2%           | 21.2 | 25.0                | 84.8%            | 0.5%  |
| 2,6-Dinitrotoluene         | 63.7 | 75.0               | 84.9%           | 67.5 | 75.0                | 90.0%            | 5.8%  |
| 2,4-Dinitrotoluene         | 62.0 | 75.0               | 82.7%           | 69.3 | 75.0                | 92.4%            | 11.1% |
| Diethylphthalate           | 21.5 | 25.0               | 86.0%           | 21.1 | 25.0                | 84.4%            | 1.9%  |
| 4-Chlorophenyl-phenylether | 20.8 | 25.0               | 83.2%           | 19.8 | 25.0                | 79.2%            | 4.9%  |
| Fluorene                   | 22.2 | 25.0               | 88.8%           | 21.4 | 25.0                | 85.6%            | 3.7%  |
| 4-Nitroaniline             | 79.4 | 75.0               | 106%            | 80.1 | 75.0                | 107%             | 0.9%  |
| 4,6-Dinitro-2-Methylphenol | 143  | 138                | 104%            | 136  | 138                 | 98.6%            | 5.0%  |
| N-Nitrosodiphenylamine     | 19.2 | 25.0               | 76.8%           | 19.1 | 25.0                | 76.4%            | 0.5%  |
| 4-Bromophenyl-phenylether  | 21.6 | 25.0               | 86.4%           | 20.6 | 25.0                | 82.4%            | 4.7%  |
| Hexachlorobenzene          | 18.7 | 25.0               | 74.8%           | 19.6 | 25.0                | 78.4%            | 4.7%  |
| Pentachlorophenol          | 62.7 | 75.0               | 83.6%           | 62.1 | 75.0                | 82.8%            | 1.0%  |
| Phenanthrene               | 22.4 | 25.0               | 89.6%           | 22.2 | 25.0                | 88.8%            | 0.9%  |
| Carbazole                  | 21.2 | 25.0               | 84.8%           | 19.8 | 25.0                | 79.2%            | 6.8%  |
| Anthracene                 | 21.8 | 25.0               | 87.2%           | 22.7 | 25.0                | 90.8%            | 4.0%  |
| Di-n-Butylphthalate        | 23.6 | 25.0               | 94.4%           | 23.3 | 25.0                | 93.2%            | 1.3%  |
| Fluoranthene               | 22.7 | 25.0               | 90.8%           | 22.4 | 25.0                | 89.6%            | 1.3%  |
| Pyrene                     | 20.7 | 25.0               | 82.8%           | 21.7 | 25.0                | 86.8%            | 4.7%  |
| Butylbenzylphthalate       | 20.4 | 25.0               | 81.6%           | 23.3 | 25.0                | 93.2%            | 13.3% |
| 3,3'-Dichlorobenzidine     | 60.2 | 75.0               | 80.3%           | 65.6 | 75.0                | 87.5%            | 8.6%  |
| Benzo(a)anthracene         | 21.5 | 25.0               | 86.0%           | 23.7 | 25.0                | 94.8%            | 9.7%  |
| bis(2-Ethylhexyl)phthalate | 23.0 | 25.0               | 92.0%           | 20.3 | 25.0                | 81.2%            | 12.5% |
| Chrysene                   | 19.9 | 25.0               | 79.6%           | 20.2 | 25.0                | 80.8%            | 1.5%  |
| Di-n-Octyl phthalate       | 21.3 | 25.0               | 85.2%           | 21.7 | 25.0                | 86.8%            | 1.9%  |
| Benzo(b)fluoranthene       | 24.6 | 25.0               | 98.4%           | 24.2 | 25.0                | 96.8%            | 1.6%  |
| Benzo(k)fluoranthene       | 21.1 | 25.0               | 84.4%           | 23.4 | 25.0                | 93.6%            | 10.3% |
| Benzo(a)pyrene             | 23.5 | 25.0               | 94.0%           | 23.5 | 25.0                | 94.0%            | 0.0%  |
| Indeno(1,2,3-cd)pyrene     | 23.0 | 25.0               | 92.0%           | 22.0 | 25.0                | 88.0%            | 4.4%  |
| Dibenz(a,h)anthracene      | 22.9 | 25.0               | 91.6%           | 21.8 | 25.0                | 87.2%            | 4.9%  |
| Benzo(g,h,i)perylene       | 22.8 | 25.0               | 91.2%           | 25.8 | 25.0                | 103%             | 12.3% |
| 3&4-Methylphenol           | 19.2 | 25.0               | 76.8%           | 18.3 | 25.0                | 73.2%            | 4.8%  |
| 1-Methylnaphthalene        | 19.9 | 25.0               | 79.6%           | 22.5 | 25.0                | 90.0%            | 12.3% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Page 3 of 3

Sample ID: LCS-112414  
LCS/LCSD

Lab Sample ID: LCS-112414  
LIMS ID: 14-25246  
Matrix: Water  
Date Analyzed LCS: 11/28/14 11:37  
LCSD: 11/28/14 12:11

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| Analyte                  | LCS  | Spike<br>Added-LCS | LCS<br>Recovery | LCSD | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD  |
|--------------------------|------|--------------------|-----------------|------|---------------------|------------------|------|
| Total Benzofluoranthenes | 44.9 | 50.0               | 89.8%           | 46.7 | 50.0                | 93.4%            | 3.9% |

Semivolatile Surrogate Recovery

|                        | LCS   | LCSD  |
|------------------------|-------|-------|
| d5-Nitrobenzene        | 74.8% | 78.0% |
| 2-Fluorobiphenyl       | 69.6% | 71.6% |
| d14-p-Terphenyl        | 76.4% | 76.0% |
| d4-1,2-Dichlorobenzene | 63.6% | 59.2% |
| d5-Phenol              | 79.7% | 74.1% |
| 2-Fluorophenol         | 66.1% | 72.8% |
| 2,4,6-Tribromophenol   | 81.9% | 90.7% |
| d4-2-Chlorophenol      | 72.3% | 69.1% |

Results reported in µg/L  
RPD calculated using sample concentrations per SW846.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-10-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38A

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25246

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/18/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 15:01

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 58.8% |
| Tetrachlorometaxylene | 75.5% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-2-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38B

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25247

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/18/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 15:19

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 64.8% |
| Tetrachlorometaxylene | 84.8% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Extraction Method: SW3510C

Page 1 of 1

Sample ID: LMW-4-1114

SAMPLE

Lab Sample ID: ZL38C

LIMS ID: 14-25248

Matrix: Water

Data Release Authorized: *mm*

Reported: 12/17/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/16/14 15:36

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 69.0% |
| Tetrachlorometaxylene | 116%  |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-3-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38D

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25249

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/19/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 15:53

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 68.2% |
| Tetrachlorometaxylene | 83.8% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: EB-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38E

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25250

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: 

Date Sampled: 11/19/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 16:10

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 65.2% |
| Tetrachlorometaxylene | 82.8% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-8-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38F

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25251

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/19/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 16:27

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 52.2% |
| Tetrachlorometaxylene | 75.5% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Extraction Method: SW3510C

Page 1 of 1

Sample ID: LMW-5-1114

SAMPLE

Lab Sample ID: ZL38G

LIMS ID: 14-25252

Matrix: Water

Data Release Authorized: *MM*

Reported: 12/17/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/16/14 16:45

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 77.5% |
| Tetrachlorometaxylene | 109%  |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Extraction Method: SW3510C

Page 1 of 1

Sample ID: LMW-9-1114

SAMPLE

Lab Sample ID: ZL38H

LIMS ID: 14-25253

Matrix: Water

Data Release Authorized: *MMW*

Reported: 12/17/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/16/14 17:02

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 83.8% |
| Tetrachlorometaxylene | 89.2% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: MB-112514

Extraction Method: SW3510C

METHOD BLANK

Page 1 of 1

Lab Sample ID: MB-112514

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: *MW*

Reported: 12/17/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: NA

Date Received: NA

Date Extracted: 11/25/14

Date Analyzed: 12/16/14 14:10

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 45.2% |
| Tetrachlorometaxylene | 63.2% |

## SW8081/PESTICIDE WATER SURROGATE RECOVERY SUMMARY

Matrix: Water

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| Client ID   | DCBP  | TCMX  | TOT OUT |
|-------------|-------|-------|---------|
| MB-112514   | 45.2% | 63.2% | 0       |
| LCS-112514  | 60.8% | 79.8% | 0       |
| LCSD-112514 | 64.2% | 85.0% | 0       |
| LMW-10-1114 | 58.8% | 75.5% | 0       |
| LMW-2-1114  | 64.8% | 84.8% | 0       |
| LMW-4-1114  | 69.0% | 116%  | 0       |
| LMW-3-1114  | 68.2% | 83.8% | 0       |
| EB-1114     | 65.2% | 82.8% | 0       |
| LMW-8-1114  | 52.2% | 75.5% | 0       |
| LMW-5-1114  | 77.5% | 109%  | 0       |
| LMW-9-1114  | 83.8% | 89.2% | 0       |

|                                | LCS/MB LIMITS | QC LIMITS |
|--------------------------------|---------------|-----------|
| (DCBP) = Decachlorobiphenyl    | (11-144)      | (11-144)  |
| (TCMX) = Tetrachlorometaxylene | (30-120)      | (30-120)  |

Prep Method: SW3510C  
Log Number Range: 14-25246 to 14-25253

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Page 1 of 1

Sample ID: LCS-112514

LCS/LCSD

Lab Sample ID: LCS-112514

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: *W*

Reported: 12/17/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted LCS/LCSD: 11/25/14

Sample Amount LCS: 500 mL

LCSD: 500 mL

Date Analyzed LCS: 12/16/14 14:27

Final Extract Volume LCS: 5.0 mL

LCSD: 12/16/14 14:44

LCSD: 5.0 mL

Instrument/Analyst LCS: ECD6/YZ

Dilution Factor LCS: 1.00

LCSD: ECD6/YZ

LCSD: 1.00

GPC Cleanup: No

Sulfur Cleanup: No

Florisol Cleanup: No

Silica Gel: No

| Analyte             | LCS   | Spike<br>Added-LCS | LCS<br>Recovery | LCSD  | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|---------------------|-------|--------------------|-----------------|-------|---------------------|------------------|-------|
| alpha-BHC           | 0.204 | 0.200              | 102%            | 0.222 | 0.200               | 111%             | 8.5%  |
| beta-BHC            | 0.189 | 0.200              | 94.5%           | 0.211 | 0.200               | 106%             | 11.0% |
| delta-BHC           | 0.158 | 0.200              | 79.0%           | 0.174 | 0.200               | 87.0%            | 9.6%  |
| gamma-BHC (Lindane) | 0.213 | 0.200              | 106%            | 0.231 | 0.200               | 116%             | 8.1%  |
| Heptachlor          | 0.184 | 0.200              | 92.0%           | 0.217 | 0.200               | 108%             | 16.5% |
| Aldrin              | 0.145 | 0.200              | 72.5%           | 0.163 | 0.200               | 81.5%            | 11.7% |
| Heptachlor Epoxide  | 0.152 | 0.200              | 76.0%           | 0.161 | 0.200               | 80.5%            | 5.8%  |
| Endosulfan I        | 0.156 | 0.200              | 78.0%           | 0.164 | 0.200               | 82.0%            | 5.0%  |
| Dieldrin            | 0.330 | 0.400              | 82.5%           | 0.350 | 0.400               | 87.5%            | 5.9%  |
| 4,4'-DDE            | 0.335 | 0.400              | 83.8%           | 0.351 | 0.400               | 87.8%            | 4.7%  |
| Endrin              | 0.199 | 0.400              | 49.8%           | 0.396 | 0.400               | 99.0%            | 66.2% |
| Endosulfan II       | 0.345 | 0.400              | 86.2%           | 0.365 | 0.400               | 91.2%            | 5.6%  |
| 4,4'-DDD            | 0.364 | 0.400              | 91.0%           | 0.384 | 0.400               | 96.0%            | 5.3%  |
| Endosulfan Sulfate  | 0.336 | 0.400              | 84.0%           | 0.378 | 0.400               | 94.5%            | 11.8% |
| 4,4'-DDT            | 0.404 | 0.400              | 101%            | 0.416 | 0.400               | 104%             | 2.9%  |
| Methoxychlor        | 1.37  | 2.00               | 68.5%           | 1.49  | 2.00                | 74.5%            | 8.4%  |
| Endrin Ketone       | 0.385 | 0.400              | 96.2%           | 0.338 | 0.400               | 84.5%            | 13.0% |
| Endrin Aldehyde     | 0.371 | 0.400              | 92.8%           | 0.350 | 0.400               | 87.5%            | 5.8%  |
| trans-Chlordane     | 0.164 | 0.200              | 82.0%           | 0.175 | 0.200               | 87.5%            | 6.5%  |
| cis-Chlordane       | 0.162 | 0.200              | 81.0%           | 0.172 | 0.200               | 86.0%            | 6.0%  |

## Pest/PCB Surrogate Recovery

|                       | LCS   | LCSD  |
|-----------------------|-------|-------|
| Decachlorobiphenyl    | 60.8% | 64.2% |
| Tetrachlorometaxylene | 79.8% | 85.0% |

Results reported in µg/L (ppb)

RPD calculated using sample concentrations per SW846.

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-10-1114  
SAMPLE

Lab Sample ID: ZL38A  
LIMS ID: 14-25246  
Matrix: Water  
Data Release Authorized:   
Reported: 12/10/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/18/14  
Date Received: 11/19/14

Date Extracted: 11/25/14  
Date Analyzed: 12/08/14 18:15  
Instrument/Analyst: ECD5/JGR  
GPC Cleanup: No  
Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00  
Silica Gel: Yes  
Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 66.8% |
| Tetrachlorometaxylene | 46.0% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-2-1114  
SAMPLE

Lab Sample ID: ZL38B

LIMS ID: 14-25247

Matrix: Water

Data Release Authorized:

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 18:34

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 74.2% |
| Tetrachlorometaxylene | 56.5% |

**ORGANICS ANALYSIS DATA SHEET**  
**PCB by GC/ECD Method SW8082A**  
**Extraction Method: SW3510C**  
 Page 1 of 1

**Sample ID: LMW-4-1114**  
**SAMPLE**

Lab Sample ID: ZL38C

LIMS ID: 14-25248

Matrix: Water

Data Release Authorized: *[Signature]*

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 18:54

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 76.8% |
| Tetrachlorometaxylene | 55.5% |

**ORGANICS ANALYSIS DATA SHEET**  
**PCB by GC/ECD Method SW8082A**  
**Extraction Method: SW3510C**  
 Page 1 of 1

**Sample ID: LMW-3-1114**  
**SAMPLE**

Lab Sample ID: ZL38D  
 LIMS ID: 14-25249  
 Matrix: Water  
 Data Release Authorized: *[Signature]*  
 Reported: 12/10/14

QC Report No: ZL38-Golder Associates  
 Project: Landsburg  
 923-1000-002.R273  
 Date Sampled: 11/19/14  
 Date Received: 11/19/14

Date Extracted: 11/25/14  
 Date Analyzed: 12/08/14 19:13  
 Instrument/Analyst: ECD5/JGR  
 GPC Cleanup: No  
 Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1.00  
 Silica Gel: No  
 Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 70.5% |
| Tetrachlorometaxylene | 55.2% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E  
LIMS ID: 14-25250  
Matrix: Water  
Data Release Authorized:   
Reported: 12/10/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/25/14  
Date Analyzed: 12/08/14 19:33  
Instrument/Analyst: ECD5/JGR  
GPC Cleanup: No  
Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00  
Silica Gel: No  
Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 50.0% |
| Tetrachlorometaxylene | 54.8% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F

LIMS ID: 14-25251

Matrix: Water

Data Release Authorized:

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 19:53

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 66.0% |
| Tetrachlorometaxylene | 57.5% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-5-1114  
SAMPLE

Lab Sample ID: ZL38G  
LIMS ID: 14-25252  
Matrix: Water  
Data Release Authorized:   
Reported: 12/10/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/25/14  
Date Analyzed: 12/08/14 20:51  
Instrument/Analyst: ECD5/JGR  
GPC Cleanup: No  
Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00  
Silica Gel: No  
Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 67.8% |
| Tetrachlorometaxylene | 56.2% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-9-1114  
SAMPLE

Lab Sample ID: ZL38H  
LIMS ID: 14-25253  
Matrix: Water  
Data Release Authorized:   
Reported: 12/10/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/25/14  
Date Analyzed: 12/08/14 21:11  
Instrument/Analyst: ECD5/JGR  
GPC Cleanup: No  
Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00  
Silica Gel: No  
Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 74.8% |
| Tetrachlorometaxylene | 53.8% |

**ORGANICS ANALYSIS DATA SHEET**  
**PCB by GC/ECD Method SW8082A**  
**Extraction Method: SW3510C**  
 Page 1 of 1

**Sample ID: MB-112514**  
**METHOD BLANK**

Lab Sample ID: MB-112514  
 LIMS ID: 14-25246  
 Matrix: Water  
 Data Release Authorized: *AL*  
 Reported: 12/10/14

QC Report No: ZL38-Golder Associates  
 Project: Landsburg  
 923-1000-002.R273  
 Date Sampled: NA  
 Date Received: NA

Date Extracted: 11/25/14  
 Date Analyzed: 12/08/14 17:16  
 Instrument/Analyst: ECD5/JGR  
 GPC Cleanup: No  
 Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1.00  
 Silica Gel: Yes  
 Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 62.2% |
| Tetrachlorometaxylene | 50.5% |

**SW8082/PCB WATER SURROGATE RECOVERY SUMMARY**

Matrix: Water

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| Client ID   | DCBP<br>% REC | DCBP<br>LCL-UCL | TCMX<br>% REC | TCMX<br>LCL-UCL | TOT OUT |
|-------------|---------------|-----------------|---------------|-----------------|---------|
| MB-112514   | 62.2%         | 29-120          | 50.5%         | 32-120          | 0       |
| LCS-112514  | 79.5%         | 29-120          | 53.8%         | 32-120          | 0       |
| LCSD-112514 | 76.8%         | 29-120          | 52.8%         | 32-120          | 0       |
| LMW-10-1114 | 66.8%         | 29-120          | 46.0%         | 32-120          | 0       |
| LMW-2-1114  | 74.2%         | 29-120          | 56.5%         | 32-120          | 0       |
| LMW-4-1114  | 76.8%         | 29-120          | 55.5%         | 32-120          | 0       |
| LMW-3-1114  | 70.5%         | 29-120          | 55.2%         | 32-120          | 0       |
| EB-1114     | 50.0%         | 29-120          | 54.8%         | 32-120          | 0       |
| LMW-8-1114  | 66.0%         | 29-120          | 57.5%         | 32-120          | 0       |
| LMW-5-1114  | 67.8%         | 29-120          | 56.2%         | 32-120          | 0       |
| LMW-9-1114  | 74.8%         | 29-120          | 53.8%         | 32-120          | 0       |

Prep Method: SW3510C  
Log Number Range: 14-25246 to 14-25253

**ORGANICS ANALYSIS DATA SHEET**  
**PCB by GC/ECD Method SW8082A**  
 Page 1 of 1

**Sample ID: LCS-112514**  
**LCS/LCSD**

Lab Sample ID: LCS-112514  
 LIMS ID: 14-25246  
 Matrix: Water  
 Data Release Authorized: *AB*  
 Reported: 12/10/14

QC Report No: ZL38-Golder Associates  
 Project: Landsburg  
 923-1000-002.R273  
 Date Sampled: NA  
 Date Received: NA

Date Extracted LCS/LCSD: 11/25/14

Sample Amount LCS: 1000 mL

Date Analyzed LCS: 12/08/14 17:36

Final Extract Volume LCS: 0.50 mL

LCSD: 12/08/14 17:55

LCSD: 0.50 mL

Instrument/Analyst LCS: ECD5/JGR

Dilution Factor LCS: 1.00

LCSD: ECD5/JGR

LCSD: 1.00

GPC Cleanup: No

Silica Gel: Yes

Sulfur Cleanup: Yes

Acid Cleanup: Yes

| Analyte      | LCS   | Spike<br>Added-LCS | LCS<br>Recovery | LCSD  | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD  |
|--------------|-------|--------------------|-----------------|-------|---------------------|------------------|------|
| Aroclor 1016 | 0.042 | 0.050              | 84.0%           | 0.042 | 0.050               | 84.0%            | 0.0% |
| Aroclor 1260 | 0.045 | 0.050              | 90.0%           | 0.044 | 0.050               | 88.0%            | 2.2% |

**PCB Surrogate Recovery**

|                       | LCS   | LCSD  |
|-----------------------|-------|-------|
| Decachlorobiphenyl    | 79.5% | 76.8% |
| Tetrachlorometaxylene | 53.8% | 52.8% |

Results reported in µg/L

RPD calculated using sample concentrations per SW846.

## ORGANICS ANALYSIS DATA SHEET

NWTPH-HCID Method by GC/FID

Extraction Method: SW3510C

Page 1 of 2

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Matrix: Water

Data Release Authorized: *MW*

Reported: 11/25/14

| ARI ID                | Sample ID                 | Extraction Date | Analysis Date | DL  | Range       | Result   |
|-----------------------|---------------------------|-----------------|---------------|-----|-------------|----------|
| MB-112114<br>14-25246 | Method Blank              | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 79.4%    |
| ZL38A<br>14-25246     | LMW-10-1114<br>HC ID: --- | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 91.4%    |
| ZL38B<br>14-25247     | LMW-2-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 91.5%    |
| ZL38C<br>14-25248     | LMW-4-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 88.6%    |
| ZL38D<br>14-25249     | LMW-3-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 83.6%    |
| ZL38E<br>14-25250     | EB-1114<br>HC ID: ---     | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 90.2%    |
| ZL38F<br>14-25251     | LMW-8-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 86.0%    |
| ZL38G<br>14-25252     | LMW-5-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 92.3%    |
| ZL38H<br>14-25253     | LMW-9-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 78.2%    |

HCID SURROGATE RECOVERY SUMMARY

Matrix: Water

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| <u>Client ID</u> | <u>O-TER</u> | <u>TOT OUT</u> |
|------------------|--------------|----------------|
| MB-112114        | 79.4%        | 0              |
| LMW-10-1114      | 91.4%        | 0              |
| LMW-2-1114       | 91.5%        | 0              |
| LMW-4-1114       | 88.6%        | 0              |
| LMW-3-1114       | 83.6%        | 0              |
| EB-1114          | 90.2%        | 0              |
| LMW-8-1114       | 86.0%        | 0              |
| LMW-5-1114       | 92.3%        | 0              |
| LMW-9-1114       | 78.2%        | 0              |

|                       | <u>LCS/MB LIMITS</u> | <u>QC LIMITS</u> |
|-----------------------|----------------------|------------------|
| (O-TER) = o-Terphenyl | (50-150)             | (50-150)         |

Prep Method: SW3510C  
Log Number Range: 14-25246 to 14-25253

**ORGANICS ANALYSIS DATA SHEET**

NWTPH-HCID Method by GC/FID

Extraction Method: SW3510C

Page 2 of 2

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Matrix: Water

Data Release Authorized: *[Signature]*

Reported: 11/25/14

| ARI ID | Sample ID | Extraction<br>Date | Analysis<br>Date | DL | Range | Result |
|--------|-----------|--------------------|------------------|----|-------|--------|
|--------|-----------|--------------------|------------------|----|-------|--------|

Reported in mg/L (ppm)

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

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

Oil value based on the total peaks in the range from C24 to C38.

HC ID: DRO/RRO indicates results of organics or additional hydrocarbons in ranges are not identifiable.

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LMW-10-1114

**SAMPLE**

Lab Sample ID: ZL38A

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25246

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *XellH*

Date Sampled: 11/18/14

Reported: 12/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>7,840</b>  |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>300</b>    |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>2,950</b>  |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-96-5        | Manganese        | 0.3  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>1,330</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>81,300</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LMW-2-1114  
SAMPLE

Lab Sample ID: ZL38B

LIMS ID: 14-25247

Matrix: Water

Data Release Authorized: *Kelly*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result         | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|----------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000          | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3              | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>113,000</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000          | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>250</b>     |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>69,200</b>  |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>220</b>     |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>3,590</b>   |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>20,700</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20             | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**
**TOTAL METALS**

Page 1 of 1

Sample ID: LMW-4-1114

**SAMPLE**

Lab Sample ID: ZL38C

LIMS ID: 14-25248

Matrix: Water

Data Release Authorized: *Xe/14*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result         | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|----------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000          | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3              | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>110,000</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000          | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>1,090</b>   |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>66,300</b>  |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>160</b>     |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>3,730</b>   |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>28,400</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20             | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LMW-3-1114  
SAMPLE

Lab Sample ID: ZL38D

LIMS ID: 14-25249

Matrix: Water

Data Release Authorized: *hmk*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>38,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-89-6        | Iron             | 7.5  | 200   | 200           | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>16,000</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>60</b>     |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>1,720</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>10,200</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**
**TOTAL METALS**

Page 1 of 1

Sample ID: EB-1114

**SAMPLE**

Lab Sample ID: ZL38E

LIMS ID: 14-25250

Matrix: Water

Data Release Authorized: *12/14*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte   | DL   | LOQ   | Result | Q |
|-----------|-----------|-----------------|---------------|------------|-----------|------|-------|--------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5  | Aluminum  | 7.6  | 1,000 | 1,000  | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0  | Antimony  | 0.01 | 3     | 3      | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2  | Arsenic   | 0.05 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3  | Barium    | 1.3  | 500   | 500    | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7  | Beryllium | 0.16 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9  | Cadmium   | 0.18 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-70-2  | Calcium   | 11.3 | 500   | 500    | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3  | Chromium  | 1.2  | 1,000 | 1,000  | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4  | Cobalt    | 0.3  | 10    | 10     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8  | Copper    | 0.92 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-89-6  | Iron      | 7.5  | 200   | 200    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1  | Lead      | 0.0  | 10    | 10     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-95-4  | Magnesium | 9.6  | 1,000 | 1,000  | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-96-5  | Manganese | 0.3  | 20    | 20     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0  | Nickel    | 3.9  | 20    | 20     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-09-7  | Potassium | 65.7 | 500   | 500    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2  | Selenium  | 0.13 | 5     | 5      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4  | Silver    | 0.43 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-23-5  | Sodium    | 11.4 | 500   | 500    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0  | Thallium  | 0.00 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2  | Vanadium  | 0.27 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6  | Zinc      | 1.4  | 20    | 20     | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F

LIMS ID: 14-25251

Matrix: Water

Data Release Authorized: *[Signature]*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>63,200</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>12,400</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>33,700</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>510</b>    |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>2,120</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>11,500</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

**Sample ID: LMW-5-1114  
SAMPLE**

Lab Sample ID: ZL38G

LIMS ID: 14-25252

Matrix: Water

Data Release Authorized: *[Signature]*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>93,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-89-6        | Iron             | 7.5  | 200   | 200           | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>52,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>230</b>    |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>2,770</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>16,400</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LMW-9-1114  
SAMPLE

Lab Sample ID: ZL38H *helly*  
LIMS ID: 14-25253  
Matrix: Water  
Data Release Authorized:  
Reported: 12/19/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>81,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>1,510</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>44,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>160</b>    |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>2,490</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>15,500</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ  
LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**
**TOTAL METALS**

Page 1 of 1

**Sample ID: METHOD BLANK**

Lab Sample ID: ZL38MB

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25247

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *[Signature]*

Date Sampled: NA

Reported: 12/19/14

Date Received: NA

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte   | DL   | LOQ   | Result | Q |
|-----------|-----------|-----------------|---------------|------------|-----------|------|-------|--------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5  | Aluminum  | 7.6  | 1,000 | 1,000  | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0  | Antimony  | 0.01 | 3     | 3      | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2  | Arsenic   | 0.05 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3  | Barium    | 1.3  | 500   | 500    | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7  | Beryllium | 0.16 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9  | Cadmium   | 0.18 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-70-2  | Calcium   | 11.3 | 500   | 500    | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3  | Chromium  | 1.2  | 1,000 | 1,000  | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4  | Cobalt    | 0.3  | 10    | 10     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8  | Copper    | 0.92 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-89-6  | Iron      | 7.5  | 200   | 200    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1  | Lead      | 0.0  | 10    | 10     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-95-4  | Magnesium | 9.6  | 1,000 | 1,000  | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-96-5  | Manganese | 0.3  | 20    | 20     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0  | Nickel    | 3.9  | 20    | 20     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-09-7  | Potassium | 65.7 | 500   | 500    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2  | Selenium  | 0.13 | 5     | 5      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4  | Silver    | 0.43 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-23-5  | Sodium    | 11.4 | 500   | 500    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0  | Thallium  | 0.00 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2  | Vanadium  | 0.27 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6  | Zinc      | 1.4  | 20    | 20     | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LMW-10-1114

DUPLICATE

Lab Sample ID: ZL38A

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25246

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *lilly*

Date Sampled: 11/18/14

Reported: 12/19/14

Date Received: 11/19/14

## MATRIX DUPLICATE QUALITY CONTROL REPORT

| Analyte   | Analysis Method | Sample  | Duplicate | RPD  | Control Limit | Q |
|-----------|-----------------|---------|-----------|------|---------------|---|
| Aluminum  | 6010C           | 1,000 U | 1,000 U   | 0.0% | +/- 1,000     | L |
| Antimony  | 200.8           | 3 U     | 3 U       | 0.0% | +/- 3         | L |
| Arsenic   | 200.8           | 3 U     | 3 U       | 0.0% | +/- 3         | L |
| Barium    | 6010C           | 500 U   | 500 U     | 0.0% | +/- 500       | L |
| Beryllium | 6010C           | 2 U     | 2 U       | 0.0% | +/- 2         | L |
| Cadmium   | 6010C           | 2 U     | 2 U       | 0.0% | +/- 2         | L |
| Calcium   | 6010C           | 7,840   | 7,750     | 1.2% | +/- 20%       |   |
| Chromium  | 6010C           | 1,000 U | 1,000 U   | 0.0% | +/- 1,000     | L |
| Cobalt    | 6010C           | 10 U    | 10 U      | 0.0% | +/- 10        | L |
| Copper    | 6010C           | 3 U     | 3 U       | 0.0% | +/- 3         | L |
| Iron      | 6010C           | 300     | 300       | 0.0% | +/- 200       | L |
| Lead      | 200.8           | 10 U    | 10 U      | 0.0% | +/- 10        | L |
| Magnesium | 6010C           | 2,950   | 2,910     | 1.4% | +/- 1,000     | L |
| Manganese | 6010C           | 20 U    | 20 U      | 0.0% | +/- 20        | L |
| Nickel    | 6010C           | 20 U    | 20 U      | 0.0% | +/- 20        | L |
| Potassium | 6010C           | 1,330   | 1,330     | 0.0% | +/- 500       | L |
| Selenium  | 200.8           | 5 U     | 5 U       | 0.0% | +/- 5         | L |
| Silver    | 6010C           | 3 U     | 3 U       | 0.0% | +/- 3         | L |
| Sodium    | 6010C           | 81,300  | 80,400    | 1.1% | +/- 20%       |   |
| Thallium  | 200.8           | 2 U     | 2 U       | 0.0% | +/- 2         | L |
| Vanadium  | 6010C           | 3 U     | 3 U       | 0.0% | +/- 3         | L |
| Zinc      | 6010C           | 20 U    | 20 U      | 0.0% | +/- 20        | L |

Reported in µg/L

\*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LMW-10-1114

MATRIX SPIKE

Lab Sample ID: ZL38A

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: *12/14*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

## MATRIX SPIKE QUALITY CONTROL REPORT

| Analyte   | Analysis Method | Sample  | Spike   | Spike Added | % Recovery | Q |
|-----------|-----------------|---------|---------|-------------|------------|---|
| Aluminum  | 6010C           | 1,000 U | 2,520   | 2,000       | 126%       | N |
| Antimony  | 200.8           | 3 U     | 23      | 25          | 92.0%      |   |
| Arsenic   | 200.8           | 3 U     | 26      | 25          | 104%       |   |
| Barium    | 6010C           | 500 U   | 2,400   | 2,000       | 120%       |   |
| Beryllium | 6010C           | 2 U     | 555     | 500         | 111%       |   |
| Cadmium   | 6010C           | 2 U     | 575     | 500         | 115%       |   |
| Calcium   | 6010C           | 7,840   | 18,900  | 10,000      | 111%       |   |
| Chromium  | 6010C           | 1,000 U | 1,000 U | 500         | NR         | N |
| Cobalt    | 6010C           | 10 U    | 560     | 500         | 112%       |   |
| Copper    | 6010C           | 3 U     | 576     | 500         | 115%       |   |
| Iron      | 6010C           | 300     | 2,580   | 2,000       | 114%       |   |
| Lead      | 200.8           | 10 U    | 20      | 20          | 100%       |   |
| Magnesium | 6010C           | 2,950   | 14,700  | 10,000      | 118%       |   |
| Manganese | 6010C           | 20 U    | 580     | 500         | 116%       |   |
| Nickel    | 6010C           | 20 U    | 570     | 500         | 114%       |   |
| Potassium | 6010C           | 1,330   | 12,700  | 10,000      | 114%       |   |
| Selenium  | 200.8           | 5 U     | 80      | 80          | 100%       |   |
| Silver    | 6010C           | 3 U     | 592     | 500         | 118%       |   |
| Sodium    | 6010C           | 81,300  | 94,400  | 10,000      | 131%       | H |
| Thallium  | 200.8           | 2 U     | 25      | 25          | 100%       |   |
| Vanadium  | 6010C           | 3 U     | 570     | 500         | 114%       |   |
| Zinc      | 6010C           | 20 U    | 560     | 500         | 112%       |   |

Reported in µg/L

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

NR-Not Recovered

Percent Recovery Limits: 75-125%

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: ZL38LCS

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25247

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *7/2/14*

Date Sampled: NA

Reported: 12/19/14

Date Received: NA

**BLANK SPIKE QUALITY CONTROL REPORT**

| Analyte   | Analysis Method | Spike Found | Spike Added | % Recovery | Q |
|-----------|-----------------|-------------|-------------|------------|---|
| Aluminum  | 6010C           | 2280        | 2000        | 114%       |   |
| Antimony  | 200.8           | 23.8        | 25.0        | 95.2%      |   |
| Arsenic   | 200.8           | 25.4        | 25.0        | 102%       |   |
| Barium    | 6010C           | 2360        | 2000        | 118%       |   |
| Beryllium | 6010C           | 543         | 500         | 109%       |   |
| Cadmium   | 6010C           | 566         | 500         | 113%       |   |
| Calcium   | 6010C           | 11000       | 10000       | 110%       |   |
| Chromium  | 6010C           | 580         | 500         | 116%       |   |
| Cobalt    | 6010C           | 553         | 500         | 111%       |   |
| Copper    | 6010C           | 559         | 500         | 112%       |   |
| Iron      | 6010C           | 2300        | 2000        | 115%       |   |
| Lead      | 200.8           | 25.1        | 25.0        | 100%       |   |
| Magnesium | 6010C           | 11700       | 10000       | 117%       |   |
| Manganese | 6010C           | 533         | 500         | 107%       |   |
| Nickel    | 6010C           | 570         | 500         | 114%       |   |
| Potassium | 6010C           | 11000       | 10000       | 110%       |   |
| Selenium  | 200.8           | 78.9        | 80.0        | 98.6%      |   |
| Silver    | 6010C           | 587         | 500         | 117%       |   |
| Sodium    | 6010C           | 11000       | 10000       | 110%       |   |
| Thallium  | 200.8           | 25.2        | 25.0        | 101%       |   |
| Vanadium  | 6010C           | 567         | 500         | 113%       |   |
| Zinc      | 6010C           | 560         | 500         | 112%       |   |

Reported in µg/L

N-Control limit not met

Control Limits: 80-120%

INORGANICS ANALYSIS DATA SHEET  
Total Mercury by Method SW7470A



Data Release Authorized:   
Reported: 11/21/14  
Date Received: 11/19/14  
Page 1 of 1

QC Report No: ZL39-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| Client/<br>ARI ID             | Date<br>Sampled | Matrix | Prep Date<br>Anal Date | RL   | Result |
|-------------------------------|-----------------|--------|------------------------|------|--------|
| LMW-10-1114<br>ZL39A 14-25256 | 11/18/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-2-1114<br>ZL39B 14-25257  | 11/18/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-4-1114<br>ZL39C 14-25258  | 11/18/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-3-1114<br>ZL39D 14-25259  | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| EB-1114<br>ZL39E 14-25260     | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-8-1114<br>ZL39F 14-25261  | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-5-1114<br>ZL39G 14-25262  | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-9-1114<br>ZL39H 14-25263  | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| MB-112114<br>Method Blank     | NA              | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |

Reported in ng/L

RL-Analytical reporting limit  
U-Undetected at reported detection limit

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LMW-10-1114

DUPLICATE

Lab Sample ID: ZL39A

LIMS ID: 14-25256

Matrix: Water

Data Release Authorized: 

Reported: 11/21/14

QC Report No: ZL39-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

## MATRIX DUPLICATE QUALITY CONTROL REPORT

| Analyte | Analysis Method | Sample | Duplicate | RPD  | Control Limit | Q |
|---------|-----------------|--------|-----------|------|---------------|---|
| Mercury | 7470A           | 20.0 U | 20.0 U    | 0.0% | +/- 20.0      | L |

Reported in ng/L

\*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LMW-10-1114  
MATRIX SPIKE

Lab Sample ID: ZL39A

LIMS ID: 14-25256

Matrix: Water

Data Release Authorized 

Reported: 11/21/14

QC Report No: ZL39-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

## MATRIX SPIKE QUALITY CONTROL REPORT

| Analyte | Analysis Method | Sample | Spike | Spike Added | % Recovery | Q |
|---------|-----------------|--------|-------|-------------|------------|---|
| Mercury | 7470A           | 20.0 U | 109   | 100         | 109%       |   |

Reported in ng/L

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: ZL39LCS

LIMS ID: 14-25257

Matrix: Water

Data Release Authorized: 

Reported: 11/21/14

QC Report No: ZL39-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: NA

Date Received: NA

## BLANK SPIKE QUALITY CONTROL REPORT

| Analyte | Analysis Method | Spike Found | Spike Added | % Recovery | Q |
|---------|-----------------|-------------|-------------|------------|---|
| Mercury | 7470A           | 218         | 200         | 109%       |   |

Reported in ng/L

N-Control limit not met

Control Limits: 80-120%



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants

December 19, 2014

Doug Morell  
Golder Associates Inc.  
18300 NE Union Hill Road, Suite 200  
Redmond, WA 98052-3333

**Client Project Name: Landsburg Mine**  
**Client Project Number: 923-1000-002.R273**  
**ARI ID: ZL65 and ZL66**

Dear Mr. Morell:

Please find enclosed Chain-of-Custody (COC) record, sample receipt documentation, and the final results for the project referenced above. Analytical Resources, Inc. (ARI) accepted two water samples and a trip blank in good condition on November 20, 2014. There were no discrepancies between the COC and the sample containers' labels.

The samples were analyzed for VOCs, PCBs, HCID, Pesticides, SVOCs, Total Metals, as requested on the COC. Quality control analyses are included for your review.

The VOCs CCAL is out of control low for all associated FORM III "Q" flagged analytes. All associated samples that contain analyte have been flagged with a "Q" qualifier.

The VOCs method blank contained methylene chloride and hexachlorobutadiene. All associated samples that contain analyte have been flagged with a "B" qualifier.

The SVOCs CCAL is out of control high for all associated FORM III "Q" flagged analytes. All associated samples that contain analyte have been flagged with a "Q" qualifier.

The SVOCs LCSD is out of control low for 3,3'-Dichlorobenzidine with an RPD for 4-Chloroaniline outside of control limits.

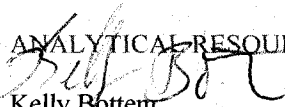
No other analytical complications were noted.

Per client request, the metals reporting limits were raised to meet client required limits.

An electronic copy of this report and all supporting raw data will remain on file at ARI. Please feel free to contact me if you have any questions or require any additional information.

Respectfully,

ANALYTICAL RESOURCES, INC.

  
Kelly Bottorri  
Client Services Manager  
(206) 695-6211  
[kellyb@arilabs.com](mailto:kellyb@arilabs.com)

|                              |                                    |
|------------------------------|------------------------------------|
| ARI Assigned Number:<br>7165 | Turn-around Requested:<br>Standard |
|------------------------------|------------------------------------|

ARI Client Company: *Golden* Phone: *475-883-0772*

Client Contact: D. Morell, J. Lamberts

Client Project Name: Landsburg

|                                       |                                |
|---------------------------------------|--------------------------------|
| Client Project #:<br>0231000002. R273 | Samplers:<br>Rudecki, Lamberts |
|---------------------------------------|--------------------------------|

Page: 1 of 1

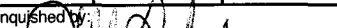
|                     |                   |
|---------------------|-------------------|
| Date:<br>11/20/2014 | Ice<br>Present? Y |
|---------------------|-------------------|

|                 |   |               |              |
|-----------------|---|---------------|--------------|
| No. of Coolers: | 3 | Cooler Temps: | 46.38<br>3.9 |
|-----------------|---|---------------|--------------|



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants  
4611 South 134th Place, Suite 100  
Tukwila, WA 98168  
206-695-6200 206-695-6201 (fax)

[illegible]

|                                                                                                                                                           |                                                                                                                                                                                                    |                                                                                                                                                                                                 |                                                                              |                                                                          |
|-----------------------------------------------------------------------------------------------------------------------------------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|------------------------------------------------------------------------------|--------------------------------------------------------------------------|
| Comments/Special Instructions<br>- Ecology EIM EDD<br>**CLIENTS SPECIFIC RLS<br>and ANALYTIC LIST**<br>Pls cc j.lamberts@golder.com<br>dmorell@golder.com | Relinquished by:<br>(Signature) <br>Printed Name: J. Lamberts<br>Company: Golder<br>Date & Time: 11/20/14 17:48 | Received by:<br>(Signature) <br>Printed Name: A. Volgardsen<br>Company: Arel<br>Date & Time: 11/20/14 17:48 | Relinquished by:<br>(Signature)<br>Printed Name:<br>Company:<br>Date & Time: | Received by:<br>(Signature)<br>Printed Name:<br>Company:<br>Date & Time: |
|-----------------------------------------------------------------------------------------------------------------------------------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|------------------------------------------------------------------------------|--------------------------------------------------------------------------|

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, notwithstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.



# Cooler Receipt Form

ARI Client: Golder

Project Name: Landsburg

COC No(s): \_\_\_\_\_ (NA)

Delivered by: Fed-Ex UPS Courier Hand Delivered Other: \_\_\_\_\_

Assigned ARI Job No: 2115

Tracking No: \_\_\_\_\_ (NA)

## Preliminary Examination Phase:

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES (NO)

Were custody papers included with the cooler? ..... YES (NO)

Were custody papers properly filled out (ink, signed, etc.) ..... YES (NO)

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry)

Time: 1748

4.6 3.8 3.9

If cooler temperature is out of compliance fill out form 00070F

Temp Gun ID#: 91677952

Cooler Accepted by: AV Date: 11/20/14 Time: 1748

**Complete custody forms and attach all shipping documents**

## Log-In Phase:

Was a temperature blank included in the cooler? ..... YES (NO)

What kind of packing material was used? ... Bubble Wrap (Wet Ice) Gel Packs (Baggies) Foam Block Paper Other: \_\_\_\_\_

Was sufficient ice used (if appropriate)? ..... NA (YES) (NO)

Were all bottles sealed in individual plastic bags? ..... (YES) (NO)

Did all bottles arrive in good condition (unbroken)? ..... (YES) (NO)

Were all bottle labels complete and legible? ..... (YES) (NO)

Did the number of containers listed on COC match with the number of containers received? ..... (YES) (NO)

Did all bottle labels and tags agree with custody papers? ..... (YES) (NO)

Were all bottles used correct for the requested analyses? ..... (YES) (NO)

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA (YES) (NO)

Were all VOC vials free of air bubbles? ..... NA (YES) (NO)

Was sufficient amount of sample sent in each bottle? ..... (YES) (NO)

Date VOC Trip Blank was made at ARI..... NA 11/14/14

Was Sample Split by ARI : (NA) YES Date/Time: \_\_\_\_\_ Equipment: \_\_\_\_\_ Split by: \_\_\_\_\_

Samples Logged by: AV Date: 11/21/14 Time: 1200

**\*\* Notify Project Manager of discrepancies or concerns \*\***

| Sample ID on Bottle | Sample ID on COC | Sample ID on Bottle | Sample ID on COC |
|---------------------|------------------|---------------------|------------------|
|                     |                  |                     |                  |
|                     |                  |                     |                  |
|                     |                  |                     |                  |
|                     |                  |                     |                  |

## Additional Notes, Discrepancies, & Resolutions:

TB=45m

By: AV

Date: 11/21/14



Small → "sm" (< 2 mm)

Peabubbles → "pb" (2 to < 4 mm)

Large → "lg" (4 to < 6 mm)

Headspace → "hs" (> 6 mm)

**PRESERVATION VERIFICATION 11/21/14**

Page 1 of 1



ARI Job No: **ZL65**

Inquiry Number: NONE  
 Analysis Requested: 11/21/14  
 Contact: Morell, Douglas  
 Client: Golder Associates  
 Logged by: AV  
 Sample Set Used: Yes-119  
 Validatable Package: Lv4  
 Deliverables:

PC: Kelly  
 VTSR: 11/20/14

Project #: 9231000002.R273  
 Project: Landsburg  
 Sample Site:  
 SDG No:  
 Analytical Protocol: In-house

| LOGNUM<br>ARI ID  | CLIENT ID   | CN<br>>12 | WAD<br>>12 | NH3<br><2 | COD<br><2 | FOG<br><2 | MET<br><2  | PHEN<br><2 | PHOS<br><2 | TKN<br><2 | NO23<br><2 | TOC<br><2 | S2<br>>9 | TPHD<br><2 | Fe2+<br><2 | DMET<br>FLT | DOC<br>FLT | PARAMETER | ADJUSTED<br>TO | LOT<br>NUMBER | AMOUNT<br>ADDED | DATE/BY |
|-------------------|-------------|-----------|------------|-----------|-----------|-----------|------------|------------|------------|-----------|------------|-----------|----------|------------|------------|-------------|------------|-----------|----------------|---------------|-----------------|---------|
| 14-25426<br>ZL65A | LMW-11-1114 |           |            |           |           |           | TOT<br>P/G |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |
| 14-25427<br>ZL65B | LMW-6-1114  |           |            |           |           |           | TOT<br>P/G |            |            |           |            |           |          |            |            |             |            |           |                |               |                 |         |

110000:5972

Checked By AV Date 11/21/14

**PRESERVATION VERIFICATION 11/21/14**

Page 1 of 1



ARI Job No: **ZL66**

PC: Kelly  
VTSR: 11/20/14

Inquiry Number: NONE  
Analysis Requested: 11/21/14  
Contact: Morell, Douglas  
Client: Golder Associates  
Logged by: AV  
Sample Set Used: Yes-119  
Validatable Package: Lv4  
Deliverables:

Project #: 9231000002.R273  
Project: Landsburg  
Sample Site:  
SDG No:  
Analytical Protocol: In-house

| LOGNUM<br>ARI ID         | CLIENT ID   | CN<br>>12 | WAD<br>>12 | NH3<br><2 | COD<br><2 | FOG<br><2 | MET<br><2 | PHEN<br><2 | PHOS<br><2 | TKN<br><2 | NO23<br><2 | TOC<br><2 | S2<br>>9 | TPHD<br><2 | Fe2+<br><2 | DMET DOC<br>FLT FLT | PARAMETER | ADJUSTED<br>TO | LOT<br>NUMBER | AMOUNT<br>ADDED | DATE/BY |
|--------------------------|-------------|-----------|------------|-----------|-----------|-----------|-----------|------------|------------|-----------|------------|-----------|----------|------------|------------|---------------------|-----------|----------------|---------------|-----------------|---------|
| 14-25429<br><b>ZL66A</b> | LMW-11-1114 |           |            |           |           |           | TOT<br>AS |            |            |           |            |           |          |            |            |                     |           |                |               |                 |         |
| 14-25430<br><b>ZL66B</b> | LMW-6-1114  |           |            |           |           |           | TOT<br>AS |            |            |           |            |           |          |            |            |                     |           |                |               |                 |         |

9231000002.R273

Checked By AV Date 11/21/14

# Sample ID Cross Reference Report



ARI Job No: ZL65  
Client: Golder Associates  
Project Event: 9231000002.R273  
Project Name: Landsburg

| Sample ID            | ARI<br>Lab ID | ARI<br>LIMS ID | Matrix | Sample Date/Time | VTSR           |
|----------------------|---------------|----------------|--------|------------------|----------------|
| 1. LMW-11-1114       | ZL65A         | 14-25426       | Water  | 11/20/14 10:45   | 11/20/14 17:48 |
| 2. LMW-6-1114        | ZL65B         | 14-25427       | Water  | 11/20/14 12:40   | 11/20/14 17:48 |
| 3. TRIP BLANK-112014 | ZL65C         | 14-25428       | Water  | 11/20/14         | 11/20/14 17:48 |

# Sample ID Cross Reference Report



ARI Job No: ZL66  
Client: Golder Associates  
Project Event: 9231000002.R273  
Project Name: Landsburg

| Sample ID      | ARI<br>Lab ID | ARI<br>LIMS ID | Matrix | Sample Date/Time | VTSR           |
|----------------|---------------|----------------|--------|------------------|----------------|
| 1. LMW-11-1114 | ZL66A         | 14-25429       | Water  | 11/20/14 10:45   | 11/20/14 17:48 |
| 2. LMW-6-1114  | ZL66B         | 14-25430       | Water  | 11/20/14 12:40   | 11/20/14 17:48 |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-11-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL65A

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25426

Project: Landsburg

Matrix: Water

9231000002.R273

Data Release Authorized: *W*

Date Sampled: 11/20/14

Reported: 12/03/14

Date Received: 11/20/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 12/01/14 14:09

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-11-1114

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SAMPLE

Lab Sample ID: ZL65A

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25426

Project: Landsburg

Matrix: Water

9231000002.R273

Date Analyzed: 12/01/14 14:09

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 102%  |
| d8-Toluene             | 104%  |
| Bromofluorobenzene     | 92.7% |
| d4-1,2-Dichlorobenzene | 103%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-6-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL65B

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25427

Project: Landsburg

Matrix: Water

9231000002.R273

Data Release Authorized: *W*

Date Sampled: 11/20/14

Reported: 12/03/14

Date Received: 11/20/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 12/01/14 14:36

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-6-1114

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SAMPLE

Lab Sample ID: ZL65B

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25427

Project: Landsburg

Matrix: Water

9231000002.R273

Date Analyzed: 12/01/14 14:36

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 106%  |
| d8-Toluene             | 101%  |
| Bromofluorobenzene     | 95.1% |
| d4-1,2-Dichlorobenzene | 106%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: TRIP BLANK-112014

Page 1 of 2

SAMPLE

Lab Sample ID: ZL65C

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25428

Project: Landsburg

Matrix: Water

9231000002.R273

Data Release Authorized: *W*

Date Sampled: 11/20/14

Reported: 12/03/14

Date Received: 11/20/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 12/01/14 15:03

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | 2.6 B    |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Page 2 of 2

Sample ID: TRIP BLANK-112014

SAMPLE

Lab Sample ID: ZL65C

LIMS ID: 14-25428

Matrix: Water

Date Analyzed: 12/01/14 15:03

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 108%  |
| d8-Toluene             | 101%  |
| Bromofluorobenzene     | 93.0% |
| d4-1,2-Dichlorobenzene | 105%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 2Sample ID: MB-120114A  
METHOD BLANK

Lab Sample ID: MB-120114A

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: *MW*

Reported: 12/03/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: NA

Date Received: NA

Instrument/Analyst: NT2/LH

Date Analyzed: 12/01/14 13:31

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | 0.92 J   |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: MB-120114A

Page 2 of 2

METHOD BLANK

Lab Sample ID: MB-120114A

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25426

Project: Landsburg

Matrix: Water

9231000002.R273

Date Analyzed: 12/01/14 13:31

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | 0.49     |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 103%  |
| d8-Toluene             | 98.2% |
| Bromofluorobenzene     | 96.9% |
| d4-1,2-Dichlorobenzene | 106%  |

## VOA SURROGATE RECOVERY SUMMARY

Matrix: Water

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| ARI ID       | Client ID         | PV | DCE   | TOL   | BFB   | DCB   | TOT OUT |
|--------------|-------------------|----|-------|-------|-------|-------|---------|
| MB-120114A   | Method Blank      | 10 | 103%  | 98.2% | 96.9% | 106%  | 0       |
| LCS-120114A  | Lab Control       | 10 | 98.3% | 102%  | 100%  | 97.3% | 0       |
| LCSD-120114A | Lab Control Dup   | 10 | 100%  | 101%  | 102%  | 98.7% | 0       |
| ZL65A        | LMW-11-1114       | 10 | 102%  | 104%  | 92.7% | 103%  | 0       |
| ZL65B        | LMW-6-1114        | 10 | 106%  | 101%  | 95.1% | 106%  | 0       |
| ZL65C        | TRIP BLANK-112014 | 10 | 108%  | 101%  | 93.0% | 105%  | 0       |

## LCS/MB LIMITS

## QC LIMITS

## SW8260C

(DCE) = d4-1,2-Dichloroethane  
(TOL) = d8-Toluene  
(BFB) = Bromofluorobenzene  
(DCB) = d4-1,2-Dichlorobenzene(80-120)  
(80-120)  
(80-120)  
(80-120)(80-120)  
(80-120)  
(80-120)  
(80-120)Prep Method: SW5030B  
Log Number Range: 14-25426 to 14-25428

**ORGANICS ANALYSIS DATA SHEET**

Volatiles by Purge & Trap GC/MS-Method SW8260C

Sample ID: LCS-120114A

Page 1 of 2

LAB CONTROL SAMPLE

Lab Sample ID: LCS-120114A

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25426

Project: Landsburg

Matrix: Water

9231000002.R273

Data Release Authorized: *[Signature]*

Date Sampled: NA

Reported: 12/03/14

Date Received: NA

Instrument/Analyst LCS: NT2/LH

Sample Amount LCS: 10.0 mL

LCSD: NT2/LH

LCSD: 10.0 mL

Date Analyzed LCS: 12/01/14 12:37

Purge Volume LCS: 10.0 mL

LCSD: 12/01/14 13:04

LCSD: 10.0 mL

| Analyte                             | LCS    | Spike<br>Added-LCS | LCS<br>Recovery | LCSD   | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD  |
|-------------------------------------|--------|--------------------|-----------------|--------|---------------------|------------------|------|
| Chloromethane                       | 10.8   | 10.0               | 108%            | 10.1   | 10.0                | 101%             | 6.7% |
| Bromomethane                        | 11.0   | 10.0               | 110%            | 10.5   | 10.0                | 105%             | 4.7% |
| Vinyl Chloride                      | 11.1   | 10.0               | 111%            | 10.7   | 10.0                | 107%             | 3.7% |
| Chloroethane                        | 12.5   | 10.0               | 125%            | 11.6   | 10.0                | 116%             | 7.5% |
| Methylene Chloride                  | 11.4 B | 10.0               | 114%            | 10.8 B | 10.0                | 108%             | 5.4% |
| Acetone                             | 47.4   | 50.0               | 94.8%           | 47.7   | 50.0                | 95.4%            | 0.6% |
| Carbon Disulfide                    | 11.2   | 10.0               | 112%            | 10.6   | 10.0                | 106%             | 5.5% |
| 1,1-Dichloroethene                  | 10.9   | 10.0               | 109%            | 10.4   | 10.0                | 104%             | 4.7% |
| 1,1-Dichloroethane                  | 11.0   | 10.0               | 110%            | 10.6   | 10.0                | 106%             | 3.7% |
| trans-1,2-Dichloroethene            | 11.1   | 10.0               | 111%            | 10.5   | 10.0                | 105%             | 5.6% |
| cis-1,2-Dichloroethene              | 10.8   | 10.0               | 108%            | 10.5   | 10.0                | 105%             | 2.8% |
| Chloroform                          | 11.2   | 10.0               | 112%            | 10.7   | 10.0                | 107%             | 4.6% |
| 1,2-Dichloroethane                  | 10.4   | 10.0               | 104%            | 10.7   | 10.0                | 107%             | 2.8% |
| 2-Butanone                          | 49.6   | 50.0               | 99.2%           | 52.7   | 50.0                | 105%             | 6.1% |
| 1,1,1-Trichloroethane               | 11.2   | 10.0               | 112%            | 10.9   | 10.0                | 109%             | 2.7% |
| Carbon Tetrachloride                | 11.4   | 10.0               | 114%            | 11.1   | 10.0                | 111%             | 2.7% |
| Vinyl Acetate                       | 8.05 Q | 10.0               | 80.5%           | 8.81 Q | 10.0                | 88.1%            | 9.0% |
| Bromodichloromethane                | 10.8   | 10.0               | 108%            | 10.6   | 10.0                | 106%             | 1.9% |
| 1,2-Dichloropropane                 | 10.9   | 10.0               | 109%            | 10.8   | 10.0                | 108%             | 0.9% |
| cis-1,3-Dichloropropene             | 10.0   | 10.0               | 100%            | 10.1   | 10.0                | 101%             | 1.0% |
| Trichloroethene                     | 11.1   | 10.0               | 111%            | 11.0   | 10.0                | 110%             | 0.9% |
| Dibromochloromethane                | 8.74   | 10.0               | 87.4%           | 8.98   | 10.0                | 89.8%            | 2.7% |
| 1,1,2-Trichloroethane               | 10.5   | 10.0               | 105%            | 10.5   | 10.0                | 105%             | 0.0% |
| Benzene                             | 11.2   | 10.0               | 112%            | 11.0   | 10.0                | 110%             | 1.8% |
| trans-1,3-Dichloropropene           | 11.6   | 10.0               | 116%            | 11.3   | 10.0                | 113%             | 2.6% |
| 2-Chloroethylvinylether             | 8.04 Q | 10.0               | 80.4%           | 8.22 Q | 10.0                | 82.2%            | 2.2% |
| Bromoform                           | 10.5   | 10.0               | 105%            | 10.6   | 10.0                | 106%             | 0.9% |
| 4-Methyl-2-Pentanone (MIBK)         | 49.9   | 50.0               | 99.8%           | 52.3   | 50.0                | 105%             | 4.7% |
| 2-Hexanone                          | 44.8   | 50.0               | 89.6%           | 49.2   | 50.0                | 98.4%            | 9.4% |
| Tetrachloroethene                   | 10.9   | 10.0               | 109%            | 11.2   | 10.0                | 112%             | 2.7% |
| 1,1,2,2-Tetrachloroethane           | 10.1   | 10.0               | 101%            | 10.5   | 10.0                | 105%             | 3.9% |
| Toluene                             | 11.1   | 10.0               | 111%            | 10.7   | 10.0                | 107%             | 3.7% |
| Chlorobenzene                       | 11.0   | 10.0               | 110%            | 10.8   | 10.0                | 108%             | 1.8% |
| Ethylbenzene                        | 11.2   | 10.0               | 112%            | 11.2   | 10.0                | 112%             | 0.0% |
| Styrene                             | 10.5   | 10.0               | 105%            | 10.5   | 10.0                | 105%             | 0.0% |
| Trichlorofluoromethane              | 13.5   | 10.0               | 135%            | 12.7   | 10.0                | 127%             | 6.1% |
| 1,1,2-Trichloro-1,2,2-trifluoroetha | 11.5   | 10.0               | 115%            | 11.0   | 10.0                | 110%             | 4.4% |
| m,p-Xylene                          | 23.3   | 20.0               | 116%            | 22.5   | 20.0                | 112%             | 3.5% |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LCS-120114A

Page 2 of 2

LAB CONTROL SAMPLE

Lab Sample ID: LCS-120114A

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25426

Project: Landsburg

Matrix: Water

9231000002.R273

| Analyte                     | LCS    | Spike<br>Added-LCS | LCS<br>Recovery | LCSD   | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|-----------------------------|--------|--------------------|-----------------|--------|---------------------|------------------|-------|
| o-Xylene                    | 9.96   | 10.0               | 99.6%           | 9.98   | 10.0                | 99.8%            | 0.2%  |
| 1,2-Dichlorobenzene         | 10.4   | 10.0               | 104%            | 10.5   | 10.0                | 105%             | 1.0%  |
| 1,3-Dichlorobenzene         | 11.1   | 10.0               | 111%            | 10.8   | 10.0                | 108%             | 2.7%  |
| 1,4-Dichlorobenzene         | 10.7   | 10.0               | 107%            | 10.5   | 10.0                | 105%             | 1.9%  |
| Acrolein                    | 48.5   | 50.0               | 97.0%           | 49.8   | 50.0                | 99.6%            | 2.6%  |
| Iodomethane                 | 11.1   | 10.0               | 111%            | 10.5   | 10.0                | 105%             | 5.6%  |
| Acrylonitrile               | 9.89   | 10.0               | 98.9%           | 9.81   | 10.0                | 98.1%            | 0.8%  |
| 1,1-Dichloropropene         | 10.7   | 10.0               | 107%            | 10.5   | 10.0                | 105%             | 1.9%  |
| Dibromomethane              | 10.4   | 10.0               | 104%            | 10.8   | 10.0                | 108%             | 3.8%  |
| 1,1,1,2-Tetrachloroethane   | 10.7   | 10.0               | 107%            | 11.0   | 10.0                | 110%             | 2.8%  |
| 1,2-Dibromo-3-chloropropane | 8.40   | 10.0               | 84.0%           | 9.51   | 10.0                | 95.1%            | 12.4% |
| 1,2,3-Trichloropropane      | 10.1   | 10.0               | 101%            | 10.2   | 10.0                | 102%             | 1.0%  |
| trans-1,4-Dichloro-2-butene | 9.88   | 10.0               | 98.8%           | 9.80   | 10.0                | 98.0%            | 0.8%  |
| 1,3,5-Trimethylbenzene      | 11.9   | 10.0               | 119%            | 11.5   | 10.0                | 115%             | 3.4%  |
| 1,2,4-Trimethylbenzene      | 11.6   | 10.0               | 116%            | 11.3   | 10.0                | 113%             | 2.6%  |
| Hexachlorobutadiene         | 10.6 B | 10.0               | 106%            | 10.7 B | 10.0                | 107%             | 0.9%  |
| 1,2-Dibromoethane           | 11.3   | 10.0               | 113%            | 11.2   | 10.0                | 112%             | 0.9%  |
| Bromochloromethane          | 10.7   | 10.0               | 107%            | 10.6   | 10.0                | 106%             | 0.9%  |
| 2,2-Dichloropropane         | 12.5   | 10.0               | 125%            | 11.8   | 10.0                | 118%             | 5.8%  |
| 1,3-Dichloropropane         | 10.5   | 10.0               | 105%            | 10.8   | 10.0                | 108%             | 2.8%  |
| Isopropylbenzene            | 9.19   | 10.0               | 91.9%           | 8.87   | 10.0                | 88.7%            | 3.5%  |
| n-Propylbenzene             | 11.6   | 10.0               | 116%            | 11.2   | 10.0                | 112%             | 3.5%  |
| Bromobenzene                | 10.4   | 10.0               | 104%            | 10.2   | 10.0                | 102%             | 1.9%  |
| 2-Chlorotoluene             | 11.7   | 10.0               | 117%            | 11.3   | 10.0                | 113%             | 3.5%  |
| 4-Chlorotoluene             | 11.5   | 10.0               | 115%            | 11.0   | 10.0                | 110%             | 4.4%  |
| tert-Butylbenzene           | 10.7   | 10.0               | 107%            | 10.6   | 10.0                | 106%             | 0.9%  |
| sec-Butylbenzene            | 9.23   | 10.0               | 92.3%           | 8.98   | 10.0                | 89.8%            | 2.7%  |
| 4-Isopropyltoluene          | 11.5   | 10.0               | 115%            | 11.2   | 10.0                | 112%             | 2.6%  |
| n-Butylbenzene              | 11.8   | 10.0               | 118%            | 11.5   | 10.0                | 115%             | 2.6%  |
| 1,2,4-Trichlorobenzene      | 11.0   | 10.0               | 110%            | 11.2   | 10.0                | 112%             | 1.8%  |
| Naphthalene                 | 8.43 Q | 10.0               | 84.3%           | 9.08 Q | 10.0                | 90.8%            | 7.4%  |
| 1,2,3-Trichlorobenzene      | 10.8 B | 10.0               | 108%            | 11.3 B | 10.0                | 113%             | 4.5%  |

Reported in µg/L (ppb)

RPD calculated using sample concentrations per SW846.

## Volatile Surrogate Recovery

|                        | LCS   | LCSD  |
|------------------------|-------|-------|
| d4-1,2-Dichloroethane  | 98.3% | 100%  |
| d8-Toluene             | 102%  | 101%  |
| Bromofluorobenzene     | 100%  | 102%  |
| d4-1,2-Dichlorobenzene | 97.3% | 98.7% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-11-1114  
SAMPLE

Lab Sample ID: ZL65A  
LIMS ID: 14-25426  
Matrix: Water  
Data Release Authorized: *MW*  
Reported: 12/04/14

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273  
Date Sampled: 11/20/14  
Date Received: 11/20/14

Date Extracted: 11/27/14  
Date Analyzed: 12/03/14 00:53  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

Lab Sample ID: ZL65A  
LIMS ID: 14-25426  
Matrix: Water  
Date Analyzed: 12/03/14 00:53

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 63.2% | 2-Fluorobiphenyl       | 65.6% |
| d14-p-Terphenyl      | 68.8% | d4-1,2-Dichlorobenzene | 64.8% |
| d5-Phenol            | 68.5% | 2-Fluorophenol         | 65.6% |
| 2,4,6-Tribromophenol | 59.7% | d4-2-Chlorophenol      | 66.7% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-6-1114  
SAMPLE

Lab Sample ID: ZL65B  
LIMS ID: 14-25427  
Matrix: Water  
Data Release Authorized: *mm*  
Reported: 12/04/14

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273  
Date Sampled: 11/20/14  
Date Received: 11/20/14

Date Extracted: 11/27/14  
Date Analyzed: 12/03/14 01:27  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-6-1114  
SAMPLE

Lab Sample ID: ZL65B  
LIMS ID: 14-25427  
Matrix: Water  
Date Analyzed: 12/03/14 01:27

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 60.8% | 2-Fluorobiphenyl       | 57.2% |
| d14-p-Terphenyl      | 64.4% | d4-1,2-Dichlorobenzene | 62.0% |
| d5-Phenol            | 71.5% | 2-Fluorophenol         | 68.5% |
| 2,4,6-Tribromophenol | 56.3% | d4-2-Chlorophenol      | 69.9% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: MB-112714  
METHOD BLANK

Lab Sample ID: MB-112714  
LIMS ID: 14-25426  
Matrix: Water  
Data Release Authorized: *mm*  
Reported: 12/04/14

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273  
Date Sampled: NA  
Date Received: NA

Date Extracted: 11/27/14  
Date Analyzed: 12/02/14 20:23  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: MB-112714  
METHOD BLANK

Lab Sample ID: MB-112714  
LIMS ID: 14-25426  
Matrix: Water  
Date Analyzed: 12/02/14 20:23

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 79.2% | 2-Fluorobiphenyl       | 72.4% |
| d14-p-Terphenyl      | 86.0% | d4-1,2-Dichlorobenzene | 62.4% |
| d5-Phenol            | 74.1% | 2-Fluorophenol         | 67.5% |
| 2,4,6-Tribromophenol | 68.0% | d4-2-Chlorophenol      | 69.3% |

## SW8270 SEMIVOLATILES WATER SURROGATE RECOVERY SUMMARY

Matrix: Water

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| Client ID   | NBZ   | FBP   | TPH   | DCB   | PHL   | 2FP   | TBP   | 2CP   | TOT | OUT |
|-------------|-------|-------|-------|-------|-------|-------|-------|-------|-----|-----|
| MB-112714   | 79.2% | 72.4% | 86.0% | 62.4% | 74.1% | 67.5% | 68.0% | 69.3% | 0   |     |
| LCS-112714  | 76.0% | 69.2% | 70.0% | 68.0% | 82.7% | 83.7% | 86.1% | 77.6% | 0   |     |
| LCSD-112714 | 70.8% | 71.6% | 82.4% | 60.0% | 75.2% | 71.2% | 79.2% | 70.7% | 0   |     |
| LMW-11-1114 | 63.2% | 65.6% | 68.8% | 64.8% | 68.5% | 65.6% | 59.7% | 66.7% | 0   |     |
| LMW-6-1114  | 60.8% | 57.2% | 64.4% | 62.0% | 71.5% | 68.5% | 56.3% | 69.9% | 0   |     |

|                                | LCS/MB LIMITS | QC LIMITS |
|--------------------------------|---------------|-----------|
| (NBZ) = d5-Nitrobenzene        | (27-120)      | (27-120)  |
| (FBP) = 2-Fluorobiphenyl       | (33-120)      | (33-120)  |
| (TPH) = d14-p-Terphenyl        | (28-130)      | (28-130)  |
| (DCB) = d4-1,2-Dichlorobenzene | (20-120)      | (20-120)  |
| (PHL) = d5-Phenol              | (38-120)      | (38-120)  |
| (2FP) = 2-Fluorophenol         | (33-120)      | (33-120)  |
| (TBP) = 2,4,6-Tribromophenol   | (52-131)      | (52-131)  |
| (2CP) = d4-2-Chlorophenol      | (41-120)      | (41-120)  |

Prep Method: SW3520C  
Log Number Range: 14-25426 to 14-25427

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Page 1 of 3Sample ID: LCS-112714  
LCS/LCSDLab Sample ID: LCS-112714  
LIMS ID: 14-25426  
Matrix: Water  
Data Release Authorized: *mw*  
Reported: 12/04/14QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273  
Date Sampled: 11/20/14  
Date Received: 11/20/14

Date Extracted LCS/LCSD: 11/27/14

Sample Amount LCS: 500 mL

Date Analyzed LCS: 12/02/14 20:56

Final Extract Volume LCS: 0.50 mL

LCSD: 12/02/14 21:30

LCSD: 0.50 mL

Instrument/Analyst LCS: NT6/JZ

Dilution Factor LCS: 1.00

LCSD: NT6/JZ

LCSD: 1.00

GPC Cleanup: NO

| Analyte                      | LCS  | Spike<br>Added-LCS | LCS<br>Recovery | LCSD | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|------------------------------|------|--------------------|-----------------|------|---------------------|------------------|-------|
| Phenol                       | 19.8 | 25.0               | 79.2%           | 18.3 | 25.0                | 73.2%            | 7.9%  |
| Bis-(2-Chloroethyl) Ether    | 19.3 | 25.0               | 77.2%           | 18.2 | 25.0                | 72.8%            | 5.9%  |
| 2-Chlorophenol               | 19.8 | 25.0               | 79.2%           | 17.2 | 25.0                | 68.8%            | 14.1% |
| 1,3-Dichlorobenzene          | 14.6 | 25.0               | 58.4%           | 15.2 | 25.0                | 60.8%            | 4.0%  |
| 1,4-Dichlorobenzene          | 15.2 | 25.0               | 60.8%           | 15.4 | 25.0                | 61.6%            | 1.3%  |
| Benzyl Alcohol               | 20.6 | 25.0               | 82.4%           | 18.5 | 25.0                | 74.0%            | 10.7% |
| 1,2-Dichlorobenzene          | 15.3 | 25.0               | 61.2%           | 14.6 | 25.0                | 58.4%            | 4.7%  |
| 2-Methylphenol               | 18.4 | 25.0               | 73.6%           | 17.6 | 25.0                | 70.4%            | 4.4%  |
| 2,2'-Oxybis(1-Chloropropane) | 17.5 | 25.0               | 70.0%           | 16.8 | 25.0                | 67.2%            | 4.1%  |
| 4-Methylphenol               | 18.1 | 25.0               | 72.4%           | 17.5 | 25.0                | 70.0%            | 3.4%  |
| N-Nitroso-Di-N-Propylamine   | 18.5 | 25.0               | 74.0%           | 18.3 | 25.0                | 73.2%            | 1.1%  |
| Hexachloroethane             | 14.6 | 25.0               | 58.4%           | 15.0 | 25.0                | 60.0%            | 2.7%  |
| Nitrobenzene                 | 18.1 | 25.0               | 72.4%           | 18.0 | 25.0                | 72.0%            | 0.6%  |
| Isophorone                   | 19.1 | 25.0               | 76.4%           | 18.6 | 25.0                | 74.4%            | 2.7%  |
| 2-Nitrophenol                | 18.9 | 25.0               | 75.6%           | 18.1 | 25.0                | 72.4%            | 4.3%  |
| 2,4-Dimethylphenol           | 53.8 | 75.0               | 71.7%           | 50.4 | 75.0                | 67.2%            | 6.5%  |
| Benzoic Acid                 | 122  | 138                | 88.4%           | 124  | 138                 | 89.9%            | 1.6%  |
| bis(2-Chloroethoxy) Methane  | 17.2 | 25.0               | 68.8%           | 16.0 | 25.0                | 64.0%            | 7.2%  |
| 2,4-Dichlorophenol           | 61.0 | 75.0               | 81.3%           | 59.0 | 75.0                | 78.7%            | 3.3%  |
| 1,2,4-Trichlorobenzene       | 15.5 | 25.0               | 62.0%           | 16.2 | 25.0                | 64.8%            | 4.4%  |
| Naphthalene                  | 17.0 | 25.0               | 68.0%           | 17.1 | 25.0                | 68.4%            | 0.6%  |
| 4-Chloroaniline              | 67.4 | 75.0               | 89.9%           | 36.2 | 75.0                | 48.3%            | 60.2% |
| Hexachlorobutadiene          | 15.2 | 25.0               | 60.8%           | 15.4 | 25.0                | 61.6%            | 1.3%  |
| 4-Chloro-3-methylphenol      | 65.9 | 75.0               | 87.9%           | 61.3 | 75.0                | 81.7%            | 7.2%  |
| 2-Methylnaphthalene          | 17.8 | 25.0               | 71.2%           | 16.9 | 25.0                | 67.6%            | 5.2%  |
| Hexachlorocyclopentadiene    | 38.9 | 75.0               | 51.9%           | 42.0 | 75.0                | 56.0%            | 7.7%  |
| 2,4,6-Trichlorophenol        | 59.6 | 75.0               | 79.5%           | 65.4 | 75.0                | 87.2%            | 9.3%  |
| 2,4,5-Trichlorophenol        | 63.9 | 75.0               | 85.2%           | 67.4 | 75.0                | 89.9%            | 5.3%  |
| 2-Chloronaphthalene          | 17.6 | 25.0               | 70.4%           | 17.8 | 25.0                | 71.2%            | 1.1%  |
| 2-Nitroaniline               | 63.1 | 75.0               | 84.1%           | 66.2 | 75.0                | 88.3%            | 4.8%  |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Page 2 of 3

Sample ID: LCS-112714  
LCS/LCSD

Lab Sample ID: LCS-112714  
LIMS ID: 14-25426  
Matrix: Water  
Date Analyzed LCS: 12/02/14 20:56  
LCSD: 12/02/14 21:30

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| Analyte                    | LCS    | Spike<br>Added-LCS | LCS<br>Recovery | LCSD    | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|----------------------------|--------|--------------------|-----------------|---------|---------------------|------------------|-------|
| Dimethylphthalate          | 18.9   | 25.0               | 75.6%           | 19.9    | 25.0                | 79.6%            | 5.2%  |
| Acenaphthylene             | 17.7   | 25.0               | 70.8%           | 18.0    | 25.0                | 72.0%            | 1.7%  |
| 3-Nitroaniline             | 68.2   | 75.0               | 90.9%           | 61.7    | 75.0                | 82.3%            | 10.0% |
| Acenaphthene               | 17.6   | 25.0               | 70.4%           | 18.4    | 25.0                | 73.6%            | 4.4%  |
| 2,4-Dinitrophenol          | 109    | 138                | 79.0%           | 114     | 138                 | 82.6%            | 4.5%  |
| 4-Nitrophenol              | 62.0   | 75.0               | 82.7%           | 67.5    | 75.0                | 90.0%            | 8.5%  |
| Dibenzofuran               | 18.7   | 25.0               | 74.8%           | 19.2    | 25.0                | 76.8%            | 2.6%  |
| 2,6-Dinitrotoluene         | 59.0   | 75.0               | 78.7%           | 63.0    | 75.0                | 84.0%            | 6.6%  |
| 2,4-Dinitrotoluene         | 57.4   | 75.0               | 76.5%           | 58.6    | 75.0                | 78.1%            | 2.1%  |
| Diethylphthalate           | 18.3   | 25.0               | 73.2%           | 18.3    | 25.0                | 73.2%            | 0.0%  |
| 4-Chlorophenyl-phenylether | 17.2   | 25.0               | 68.8%           | 18.5    | 25.0                | 74.0%            | 7.3%  |
| Fluorene                   | 18.5   | 25.0               | 74.0%           | 18.5    | 25.0                | 74.0%            | 0.0%  |
| 4-Nitroaniline             | 69.5 Q | 75.0               | 92.7%           | 62.9 Q  | 75.0                | 83.9%            | 10.0% |
| 4,6-Dinitro-2-Methylphenol | 114    | 138                | 82.6%           | 117     | 138                 | 84.8%            | 2.6%  |
| N-Nitrosodiphenylamine     | 17.4   | 25.0               | 69.6%           | 14.6    | 25.0                | 58.4%            | 17.5% |
| 4-Bromophenyl-phenylether  | 18.0   | 25.0               | 72.0%           | 21.7    | 25.0                | 86.8%            | 18.6% |
| Hexachlorobenzene          | 18.3   | 25.0               | 73.2%           | 20.7    | 25.0                | 82.8%            | 12.3% |
| Pentachlorophenol          | 57.2   | 75.0               | 76.3%           | 65.3    | 75.0                | 87.1%            | 13.2% |
| Phenanthrene               | 19.3   | 25.0               | 77.2%           | 19.9    | 25.0                | 79.6%            | 3.1%  |
| Carbazole                  | 16.6   | 25.0               | 66.4%           | 17.3    | 25.0                | 69.2%            | 4.1%  |
| Anthracene                 | 19.2   | 25.0               | 76.8%           | 21.5    | 25.0                | 86.0%            | 11.3% |
| Di-n-Butylphthalate        | 19.8   | 25.0               | 79.2%           | 21.8    | 25.0                | 87.2%            | 9.6%  |
| Fluoranthene               | 18.9   | 25.0               | 75.6%           | 21.0    | 25.0                | 84.0%            | 10.5% |
| Pyrene                     | 19.1   | 25.0               | 76.4%           | 24.3    | 25.0                | 97.2%            | 24.0% |
| Butylbenzylphthalate       | 19.2   | 25.0               | 76.8%           | 21.1    | 25.0                | 84.4%            | 9.4%  |
| 3,3'-Dichlorobenzidine     | 62.2   | 75.0               | 82.9%           | < 5.0 U | 75.0                | NA%              | NA    |
| Benzo(a)anthracene         | 19.3   | 25.0               | 77.2%           | 21.4    | 25.0                | 85.6%            | 10.3% |
| bis(2-Ethylhexyl)phthalate | 20.2   | 25.0               | 80.8%           | 21.6    | 25.0                | 86.4%            | 6.7%  |
| Chrysene                   | 18.6   | 25.0               | 74.4%           | 18.6    | 25.0                | 74.4%            | 0.0%  |
| Di-n-Octyl phthalate       | 18.4   | 25.0               | 73.6%           | 19.5    | 25.0                | 78.0%            | 5.8%  |
| Benzo(b)fluoranthene       | 19.1   | 25.0               | 76.4%           | 23.5    | 25.0                | 94.0%            | 20.7% |
| Benzo(k)fluoranthene       | 19.7   | 25.0               | 78.8%           | 24.6    | 25.0                | 98.4%            | 22.1% |
| Benzo(a)pyrene             | 19.8   | 25.0               | 79.2%           | 21.2    | 25.0                | 84.8%            | 6.8%  |
| Indeno(1,2,3-cd)pyrene     | 19.9   | 25.0               | 79.6%           | 24.4    | 25.0                | 97.6%            | 20.3% |
| Dibenz(a,h)anthracene      | 19.2   | 25.0               | 76.8%           | 23.7    | 25.0                | 94.8%            | 21.0% |
| Benzo(g,h,i)perylene       | 18.5   | 25.0               | 74.0%           | 23.8    | 25.0                | 95.2%            | 25.1% |
| 3&4-Methylphenol           | 18.1   | 25.0               | 72.4%           | 17.5    | 25.0                | 70.0%            | 3.4%  |
| 1-Methylnaphthalene        | 18.7   | 25.0               | 74.8%           | 18.1    | 25.0                | 72.4%            | 3.3%  |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Page 3 of 3

Sample ID: LCS-112714  
LCS/LCSD

Lab Sample ID: LCS-112714  
LIMS ID: 14-25426  
Matrix: Water  
Date Analyzed LCS: 12/02/14 20:56  
LCSD: 12/02/14 21:30

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| Analyte                  | LCS  | Spike<br>Added-LCS | LCS<br>Recovery | LCSD | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|--------------------------|------|--------------------|-----------------|------|---------------------|------------------|-------|
| Total Benzofluoranthenes | 39.2 | 50.0               | 78.4%           | 48.8 | 50.0                | 97.6%            | 21.8% |

Semivolatile Surrogate Recovery

|                        | LCS   | LCSD  |
|------------------------|-------|-------|
| d5-Nitrobenzene        | 76.0% | 70.8% |
| 2-Fluorobiphenyl       | 69.2% | 71.6% |
| d14-p-Terphenyl        | 70.0% | 82.4% |
| d4-1,2-Dichlorobenzene | 68.0% | 60.0% |
| d5-Phenol              | 82.7% | 75.2% |
| 2-Fluorophenol         | 83.7% | 71.2% |
| 2,4,6-Tribromophenol   | 86.1% | 79.2% |
| d4-2-Chlorophenol      | 77.6% | 70.7% |

Results reported in µg/L  
RPD calculated using sample concentrations per SW846.

**ORGANICS ANALYSIS DATA SHEET**

**Pesticides/PCB by GC/ECD Method SW8081B**

**Extraction Method: SW3510C**

Page 1 of 1

**Sample ID: LMW-11-1114**

**SAMPLE**

Lab Sample ID: ZL65A

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: 

Reported: 12/17/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

Date Extracted: 11/27/14

Date Analyzed: 12/16/14 19:54

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

**Pest/PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 80.5% |
| Tetrachlorometaxylene | 88.0% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Extraction Method: SW3510C

Page 1 of 1

Sample ID: LMW-6-1114

SAMPLE

Lab Sample ID: ZL65B

LIMS ID: 14-25427

Matrix: Water

Data Release Authorized: *AB*

Reported: 12/17/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

Date Extracted: 11/27/14

Date Analyzed: 12/16/14 20:11

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 86.0% |
| Tetrachlorometaxylene | 93.5% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

**SW8081/PESTICIDE WATER SURROGATE RECOVERY SUMMARY**

Matrix: Water

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| <u>Client ID</u> | <u>DCBP</u> | <u>TCMX</u> | <u>TOT OUT</u> |
|------------------|-------------|-------------|----------------|
| MB-112714        | 80.5%       | 88.0%       | 0              |
| LCS-112714       | 83.8%       | 83.0%       | 0              |
| LCSD-112714      | 81.0%       | 87.2%       | 0              |
| LMW-11-1114      | 80.5%       | 88.0%       | 0              |
| LMW-6-1114       | 86.0%       | 93.5%       | 0              |

|                                | <u>LCS/MB LIMITS</u> | <u>QC LIMITS</u> |
|--------------------------------|----------------------|------------------|
| (DCBP) = Decachlorobiphenyl    | (11-144)             | (11-144)         |
| (TCMX) = Tetrachlorometaxylene | (30-120)             | (30-120)         |

Prep Method: SW3510C  
Log Number Range: 14-25426 to 14-25427

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Extraction Method: SW3510C

Page 1 of 1

Sample ID: MB-112714

METHOD BLANK

Lab Sample ID: MB-112714

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: 

Reported: 12/17/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: NA

Date Received: NA

Date Extracted: 11/27/14

Date Analyzed: 12/16/14 18:11

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 80.5% |
| Tetrachlorometaxylene | 88.0% |

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Page 1 of 1

Sample ID: LCS-112714

LCS/LCSD

Lab Sample ID: LCS-112714

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: 

Reported: 12/17/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

Date Extracted LCS/LCSD: 11/27/14

Sample Amount LCS: 500 mL

LCSD: 500 mL

Date Analyzed LCS: 12/16/14 18:28

Final Extract Volume LCS: 5.0 mL

LCSD: 12/16/14 18:45

LCSD: 5.0 mL

Instrument/Analyst LCS: ECD6/YZ

Dilution Factor LCS: 1.00

LCSD: ECD6/YZ

LCSD: 1.00

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Silica Gel: No

| Analyte             | Spike   |           |          | Spike |            |          | RPD   |
|---------------------|---------|-----------|----------|-------|------------|----------|-------|
|                     | LCS     | Added-LCS | Recovery | LCSD  | Added-LCSD | Recovery |       |
| alpha-BHC           | 0.227   | 0.200     | 114%     | 0.231 | 0.200      | 116%     | 1.7%  |
| beta-BHC            | 0.220   | 0.200     | 110%     | 0.218 | 0.200      | 109%     | 0.9%  |
| delta-BHC           | 0.217   | 0.200     | 108%     | 0.211 | 0.200      | 106%     | 2.8%  |
| gamma-BHC (Lindane) | 0.242   | 0.200     | 121%     | 0.243 | 0.200      | 122%     | 0.4%  |
| Heptachlor          | 0.230   | 0.200     | 115%     | 0.230 | 0.200      | 115%     | 0.0%  |
| Aldrin              | 0.204   | 0.200     | 102%     | 0.207 | 0.200      | 104%     | 1.5%  |
| Heptachlor Epoxide  | 0.231   | 0.200     | 116%     | 0.227 | 0.200      | 114%     | 1.7%  |
| Endosulfan I        | 0.236   | 0.200     | 118%     | 0.234 | 0.200      | 117%     | 0.9%  |
| Dieldrin            | 0.467   | 0.400     | 117%     | 0.461 | 0.400      | 115%     | 1.3%  |
| 4,4'-DDE            | 0.464   | 0.400     | 116%     | 0.459 | 0.400      | 115%     | 1.1%  |
| Endrin              | 0.529   | 0.400     | 132%     | 0.462 | 0.400      | 116%     | 13.5% |
| Endosulfan II       | 0.498 P | 0.400     | 124%     | 0.461 | 0.400      | 115%     | 7.7%  |
| 4,4'-DDD            | 0.521   | 0.400     | 130%     | 0.479 | 0.400      | 120%     | 8.4%  |
| Endosulfan Sulfate  | 0.529   | 0.400     | 132%     | 0.446 | 0.400      | 112%     | 17.0% |
| 4,4'-DDT            | 0.581   | 0.400     | 145%     | 0.538 | 0.400      | 134%     | 7.7%  |
| Methoxychlor        | 1.95    | 2.00      | 97.5%    | 1.80  | 2.00       | 90.0%    | 8.0%  |
| Endrin Ketone       | 0.489   | 0.400     | 122%     | 0.465 | 0.400      | 116%     | 5.0%  |
| Endrin Aldehyde     | 0.474   | 0.400     | 118%     | 0.450 | 0.400      | 112%     | 5.2%  |
| trans-Chlordane     | 0.229   | 0.200     | 114%     | 0.223 | 0.200      | 112%     | 2.7%  |
| cis-Chlordane       | 0.236   | 0.200     | 118%     | 0.233 | 0.200      | 116%     | 1.3%  |

## Pest/PCB Surrogate Recovery

|                       | LCS   | LCSD  |
|-----------------------|-------|-------|
| Decachlorobiphenyl    | 83.8% | 81.0% |
| Tetrachlorometaxylene | 83.0% | 87.2% |

Results reported in µg/L (ppb)

RPD calculated using sample concentrations per SW846.

**ORGANICS ANALYSIS DATA SHEET**

NWTPH-HCID Method by GC/FID

Extraction Method: SW3510C

Page 1 of 1

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Matrix: Water

Data Release Authorized: *MMW*

Reported: 11/25/14

| ARI ID                | Sample ID                 | Extraction Date | Analysis Date | DL  | Range                               | Result                                    |
|-----------------------|---------------------------|-----------------|---------------|-----|-------------------------------------|-------------------------------------------|
| MB-112214<br>14-25426 | Method Blank              | 11/22/14        | 11/24/14      | 1.0 | Gas<br>Diesel<br>Oil<br>o-Terphenyl | < 0.25 U<br>< 0.50 U<br>< 0.50 U<br>92.2% |
| ZL65A<br>14-25426     | LMW-11-1114<br>HC ID: --- | 11/22/14        | 11/24/14      | 1.0 | Gas<br>Diesel<br>Oil<br>o-Terphenyl | < 0.25 U<br>< 0.50 U<br>< 0.50 U<br>94.0% |
| ZL65B<br>14-25427     | LMW-6-1114<br>HC ID: ---  | 11/22/14        | 11/24/14      | 1.0 | Gas<br>Diesel<br>Oil<br>o-Terphenyl | < 0.25 U<br>< 0.50 U<br>< 0.50 U<br>97.0% |

Reported in mg/L (ppm)

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

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

Oil value based on the total peaks in the range from C24 to C38.

HC ID: DRO/RRO indicates results of organics or additional hydrocarbons in ranges are not identifiable.

HCID SURROGATE RECOVERY SUMMARY

Matrix: Water

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| <u>Client ID</u> | <u>O-TER</u> | <u>TOT OUT</u> |
|------------------|--------------|----------------|
| MB-112214        | 92.2%        | 0              |
| LMW-11-1114      | 94.0%        | 0              |
| LMW-6-1114       | 97.0%        | 0              |

|                       | <u>LCS/MB LIMITS</u> | <u>QC LIMITS</u> |
|-----------------------|----------------------|------------------|
| (O-TER) = o-Terphenyl | (50-150)             | (50-150)         |

Prep Method: SW3510C  
Log Number Range: 14-25426 to 14-25427

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-11-1114  
SAMPLE

Lab Sample ID: ZL65A  
LIMS ID: 14-25426  
Matrix: Water  
Data Release Authorized: *[Signature]*  
Reported: 12/10/14

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273  
Date Sampled: 11/20/14  
Date Received: 11/20/14

Date Extracted: 11/27/14  
Date Analyzed: 12/09/14 01:45  
Instrument/Analyst: ECD5/JGR  
GPC Cleanup: No  
Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00  
Silica Gel: Yes  
Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 72.8% |
| Tetrachlorometaxylene | 52.8% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-6-1114  
SAMPLE

Lab Sample ID: ZL65B

LIMS ID: 14-25427

Matrix: Water

Data Release Authorized:

Reported: 12/10/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

Date Extracted: 11/27/14

Date Analyzed: 12/09/14 02:05

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 58.8% |
| Tetrachlorometaxylene | 48.2% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: MB-112714  
METHOD BLANK

Lab Sample ID: MB-112714  
LIMS ID: 14-25426  
Matrix: Water  
Data Release Authorized:   
Reported: 12/10/14

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273  
Date Sampled: NA  
Date Received: NA

Date Extracted: 11/27/14  
Date Analyzed: 12/08/14 23:08  
Instrument/Analyst: ECD5/JGR  
GPC Cleanup: No  
Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00  
Silica Gel: Yes  
Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 62.0% |
| Tetrachlorometaxylene | 53.5% |

**SW8082/PCB WATER SURROGATE RECOVERY SUMMARY**

Matrix: Water

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| <u>Client ID</u> | <u>DCBP<br/>% REC</u> | <u>DCBP<br/>LCL-UCL</u> | <u>TCMX<br/>% REC</u> | <u>TCMX<br/>LCL-UCL</u> | <u>TOT OUT</u> |
|------------------|-----------------------|-------------------------|-----------------------|-------------------------|----------------|
| MB-112714        | 62.0%                 | 29-120                  | 53.5%                 | 32-120                  | 0              |
| LCS-112714       | 59.2%                 | 29-120                  | 49.2%                 | 32-120                  | 0              |
| LCSD-112714      | 59.5%                 | 29-120                  | 49.5%                 | 32-120                  | 0              |
| LMW-11-1114      | 72.8%                 | 29-120                  | 52.8%                 | 32-120                  | 0              |
| LMW-6-1114       | 58.8%                 | 29-120                  | 48.2%                 | 32-120                  | 0              |

Prep Method: SW3510C  
Log Number Range: 14-25426 to 14-25427

**ORGANICS ANALYSIS DATA SHEET**  
**PCB by GC/ECD Method SW8082A**  
 Page 1 of 1

Sample ID: LCS-112714  
 LCS/LCSD

Lab Sample ID: LCS-112714  
 LIMS ID: 14-25426  
 Matrix: Water  
 Data Release Authorized: *AB*  
 Reported: 12/10/14

QC Report No: ZL65-Golder Associates  
 Project: Landsburg  
 9231000002.R273  
 Date Sampled: NA  
 Date Received: NA

Date Extracted LCS/LCSD: 11/27/14

Sample Amount LCS: 1000 mL

Date Analyzed LCS: 12/08/14 23:28

Final Extract Volume LCS: 0.50 mL

LCSD: 12/09/14 00:27

LCSD: 0.50 mL

Instrument/Analyst LCS: ECD5/JGR

Dilution Factor LCS: 1.00

LCSD: ECD5/JGR

LCSD: 1.00

GPC Cleanup: No

Silica Gel: Yes

Sulfur Cleanup: Yes

Acid Cleanup: Yes

| Analyte      | LCS   | Spike<br>Added-LCS | LCS<br>Recovery | LCSD  | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD   |
|--------------|-------|--------------------|-----------------|-------|---------------------|------------------|-------|
| Aroclor 1016 | 0.041 | 0.050              | 82.0%           | 0.047 | 0.050               | 94.0%            | 13.6% |
| Aroclor 1260 | 0.039 | 0.050              | 78.0%           | 0.038 | 0.050               | 76.0%            | 2.6%  |

**PCB Surrogate Recovery**

|                       | LCS   | LCSD  |
|-----------------------|-------|-------|
| Decachlorobiphenyl    | 59.2% | 59.5% |
| Tetrachlorometaxylene | 49.2% | 49.5% |

Results reported in µg/L  
 RPD calculated using sample concentrations per SW846.

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LMW-11-1114  
SAMPLE

Lab Sample ID: ZL65A

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: *[Signature]*

Reported: 12/19/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | <b>7440-38-2</b> | <b>Arsenic</b>   | 0.05 | 3     | <b>7</b>      |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>58,200</b> |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>1,410</b>  |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>26,300</b> |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>170</b>    |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>2,100</b>  |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>28,800</b> |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**
**TOTAL METALS**

Page 1 of 1

**Sample ID: LMW-6-1114  
SAMPLE**

Lab Sample ID: ZL65B

LIMS ID: 14-25427

Matrix: Water

Data Release Authorized: *8/2/14*

Reported: 12/19/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>27,300</b> |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>2,410</b>  |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>14,000</b> |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>30</b>     |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>720</b>    |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>7,070</b>  |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**
**TOTAL METALS**

Page 1 of 1

**Sample ID: METHOD BLANK**

Lab Sample ID: ZL65MB

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: *[Signature]*

Reported: 12/19/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: NA

Date Received: NA

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte   | DL   | LOQ   | Result | Q |
|-----------|-----------|-----------------|---------------|------------|-----------|------|-------|--------|---|
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7429-90-5  | Aluminum  | 7.6  | 1,000 | 1,000  | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-36-0  | Antimony  | 0.01 | 3     | 3      | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-38-2  | Arsenic   | 0.05 | 3     | 3      | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-39-3  | Barium    | 1.3  | 500   | 500    | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-41-7  | Beryllium | 0.16 | 2     | 2      | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-43-9  | Cadmium   | 0.18 | 2     | 2      | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-70-2  | Calcium   | 11.3 | 500   | 500    | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-47-3  | Chromium  | 1.2  | 1,000 | 1,000  | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-48-4  | Cobalt    | 0.3  | 10    | 10     | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-50-8  | Copper    | 0.92 | 3     | 3      | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7439-89-6  | Iron      | 7.5  | 200   | 200    | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7439-92-1  | Lead      | 0.0  | 10    | 10     | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7439-95-4  | Magnesium | 9.6  | 1,000 | 1,000  | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7439-96-5  | Manganese | 0.3  | 20    | 20     | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-02-0  | Nickel    | 3.9  | 20    | 20     | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-09-7  | Potassium | 65.7 | 500   | 500    | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7782-49-2  | Selenium  | 0.13 | 5     | 5      | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-22-4  | Silver    | 0.43 | 3     | 3      | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-23-5  | Sodium    | 11.4 | 500   | 500    | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-28-0  | Thallium  | 0.00 | 2     | 2      | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-62-2  | Vanadium  | 0.27 | 3     | 3      | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-66-6  | Zinc      | 1.4  | 20    | 20     | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**
**TOTAL METALS**

Page 1 of 1

**Sample ID: LAB CONTROL**

Lab Sample ID: ZL65LCS

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25426

Project: Landsburg

Matrix: Water

9231000002.R273

Data Release Authorized:

Date Sampled: NA

Reported: 12/19/14

Date Received: NA

**BLANK SPIKE QUALITY CONTROL REPORT**

| Analyte   | Analysis Method | Spike Found | Spike Added | % Recovery | Q |
|-----------|-----------------|-------------|-------------|------------|---|
| Aluminum  | 6010C           | 2080        | 2000        | 104%       |   |
| Antimony  | 200.8           | 23.8        | 25.0        | 95.2%      |   |
| Arsenic   | 200.8           | 26.0        | 25.0        | 104%       |   |
| Barium    | 6010C           | 2080        | 2000        | 104%       |   |
| Beryllium | 6010C           | 522         | 500         | 104%       |   |
| Cadmium   | 6010C           | 535         | 500         | 107%       |   |
| Calcium   | 6010C           | 10200       | 10000       | 102%       |   |
| Chromium  | 6010C           | 529         | 500         | 106%       |   |
| Cobalt    | 6010C           | 515         | 500         | 103%       |   |
| Copper    | 6010C           | 530         | 500         | 106%       |   |
| Iron      | 6010C           | 2170        | 2000        | 108%       |   |
| Lead      | 200.8           | 25.9        | 25.0        | 104%       |   |
| Magnesium | 6010C           | 10500       | 10000       | 105%       |   |
| Manganese | 6010C           | 511         | 500         | 102%       |   |
| Nickel    | 6010C           | 510         | 500         | 102%       |   |
| Potassium | 6010C           | 10600       | 10000       | 106%       |   |
| Selenium  | 200.8           | 81.5        | 80.0        | 102%       |   |
| Silver    | 6010C           | 539         | 500         | 108%       |   |
| Sodium    | 6010C           | 10600       | 10000       | 106%       |   |
| Thallium  | 200.8           | 25.7        | 25.0        | 103%       |   |
| Vanadium  | 6010C           | 523         | 500         | 105%       |   |
| Zinc      | 6010C           | 500         | 500         | 100%       |   |

Reported in µg/L

N-Control limit not met

Control Limits: 80-120%

INORGANICS ANALYSIS DATA SHEET  
Total Mercury by Method SW7470A



Data Release Authorized: *gr*  
Reported: 12/01/14  
Date Received: 11/20/14  
Page 1 of 1

QC Report No: ZL66-Golder Associates  
Project: Landsburg  
9231000002.R273

| Client/<br>ARI ID             | Date<br>Sampled | Matrix | Prep Date<br>Anal Date | RL   | Result |
|-------------------------------|-----------------|--------|------------------------|------|--------|
| LMW-11-1114<br>ZL66A 14-25429 | 11/20/14        | Water  | 11/24/14<br>11/28/14   | 20.0 | 20.0 U |
| LMW-6-1114<br>ZL66B 14-25430  | 11/20/14        | Water  | 11/24/14<br>11/28/14   | 20.0 | 20.0 U |
| MB-112414<br>Method Blank     | NA              | Water  | 11/24/14<br>11/28/14   | 20.0 | 20.0 U |

Reported in ng/L

RL-Analytical reporting limit  
U-Undetected at reported detection limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: ZL66LCS

LIMS ID: 14-25429

Matrix: Water

Data Release Authorized:

Reported: 12/01/14

QC Report No: ZL66-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: NA

Date Received: NA

**BLANK SPIKE QUALITY CONTROL REPORT**

| Analyte | Analysis Method | Spike Found | Spike Added | % Recovery | Q |
|---------|-----------------|-------------|-------------|------------|---|
| Mercury | 7470A           | 228         | 200         | 114%       |   |

Reported in ng/L

N-Control limit not met

Control Limits: 80-120%



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants

January 23, 2015

Doug Morell  
Golder Associates Inc.  
18300 NE Union Hill Road, Suite 200  
Redmond, WA 98052-3333

**Client Project Name: Landsburg Mine**  
**Client Project Number: 923-1000-002.R273**  
**ARI ID: ZS69**

Dear Mr. Morell:

Please find enclosed Chain-of-Custody (COC) record, sample receipt documentation, and the final results for the project referenced above. Analytical Resources, Inc. (ARI) accepted eight water samples and two trip blanks in good condition on December 4, 2014. There were no discrepancies between the COC and the sample containers' labels.

The samples were originally analyzed for VOCs, PCBs, HCID, Pesticides, SVOCs, Total Metals, as requested on the COC and reported under ARI SDGs ZL38 and ZL39. Quality control analyses are included for your review.

At the request of the client, sample LMW-4-1114 was re-extracted for SVOCs outside of the method recommended holding time.

The SVOCs CCAL is out of control low for all associated FORM III "Q" flagged analytes with the exception of 3-Nitroaniline which is out of control high.

No other analytical complications were noted.

Per client request, the metals reporting limits were raised to meet client required limits.

An electronic copy of this report and all supporting raw data will remain on file at ARI. Please feel free to contact me if you have any questions or require any additional information.

Respectfully,

  
ANALYTICAL RESOURCES, INC.  
Kelly Bottom  
Client Services Manager  
(206) 695-6211  
[kellyb@arilabs.com](mailto:kellyb@arilabs.com)

# Chain of Custody Record & Laboratory Analysis Request

|                                                      |                                           |                                                           |
|------------------------------------------------------|-------------------------------------------|-----------------------------------------------------------|
| ARI Assigned Number:<br><b>2138-7565</b>             | Turn-around Requested:<br><b>Standard</b> | Page: <b>1</b> of <b>1</b>                                |
| ARI Client Company:<br><b>Goldner Associates</b>     | Phone:<br><b>425 883-0777</b>             | Date:<br><b>11/18/2014</b>                                |
| Client Contact:<br><b>Doug Morell, Jill Lamberts</b> | No. of Coolers:<br><b>7</b>               | Ice Present? <b>y</b><br>Cooler Temps: <b>0.41 - 4.17</b> |
| Client Project Name:<br><b>Landsburg</b>             |                                           |                                                           |
| Client Project #:<br><b>923-1000-002-R273</b>        | Samplers:<br><b>Rydecki, Lamberts</b>     |                                                           |



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants  
4611 South 134th Place, Suite 100  
Tukwila, WA 98168  
206-695-6200 206-695-6201 (fax)  
www.arilabs.com

| Sample ID         | Date     | Time | Matrix | No. Containers | Analysis Requested  |                               |            |                           |          |                        |                      |  | Notes/Comments |  |
|-------------------|----------|------|--------|----------------|---------------------|-------------------------------|------------|---------------------------|----------|------------------------|----------------------|--|----------------|--|
|                   |          |      |        |                | VOCs<br>Client list | PCBs (CLL)                    | Pesticides | SVOCS 8270<br>Client list | TPH-HCID | TAM L<br>-Total Metals | TAM L<br>Diss Metals |  |                |  |
| Trip Blank-111814 | 11/18/14 | -    | W      | 9-6            | X                   |                               |            |                           |          |                        |                      |  |                |  |
| LMW-10-1114       |          | 945  | W      | 17             | X                   | X                             | X          | X                         | X        | X                      | HOLD                 |  |                |  |
| LMW-2-1114        |          | 1140 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| LMW-4-1114        |          | 1310 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| Trip Blank-111914 | 11/19/14 | -    | W      | 6              | X                   | HOLD unless there are detects |            |                           |          |                        |                      |  |                |  |
| LMW-3-1114        |          | 0920 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      | HOLD                 |  |                |  |
| EB-1114           |          | 0900 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| LMW-8-1114        |          | 1005 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| LMW-5-1114        |          | 1130 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| LMW-9-1114        |          | 1315 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |

|                                                                                                                                                                            |                                                    |                                                |                                 |                             |
|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----------------------------------------------------|------------------------------------------------|---------------------------------|-----------------------------|
| Comments/Special Instructions<br><b>-Ecology ETM EDD</b><br><br><b>** CLIENT SPECIFIC RLS and ANALYTE LIST **</b><br><br>PLS cc jlamberts@golder.com<br>dmorell@golder.com | Relinquished by:<br>(Signature) <i>[Signature]</i> | Received by:<br>(Signature) <i>[Signature]</i> | Relinquished by:<br>(Signature) | Received by:<br>(Signature) |
|                                                                                                                                                                            | Printed Name:<br><b>J. Lamberts</b>                | Printed Name:<br><b>Taylor Street</b>          | Printed Name:                   | Printed Name:               |
|                                                                                                                                                                            | Company:<br><b>Goldner</b>                         | Company:<br><b>ARL</b>                         | Company:                        | Company:                    |
|                                                                                                                                                                            | Date & Time:<br><b>11/19/14 1442</b>               | Date & Time:<br><b>11/19/14 1442</b>           | Date & Time:                    | Date & Time:                |

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

2569-00002

# Sample ID Cross Reference Report



ARI Job No: ZS69  
Client: Golder Associates  
Project Event: 923-1000-002.R273  
Project Name: Landsburg

| Sample ID     | ARI<br>Lab ID | ARI<br>LIMS ID | Matrix | Sample Date/Time | VTSR           |
|---------------|---------------|----------------|--------|------------------|----------------|
| 1. LMW-4-1114 | ZS69A         | 15-518         | Water  | 11/18/14 13:10   | 11/19/14 14:42 |

**ORGANICS ANALYSIS DATA SHEET**  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-4-1114  
SAMPLE

Lab Sample ID: ZS69A  
LIMS ID: 15-518  
Matrix: Water  
Data Release Authorized:  
Reported: 01/22/15

QC Report No: ZS69-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/18/14  
Date Received: 11/19/14

Date Extracted: 01/15/15  
Date Analyzed: 01/21/15 17:13  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

Sample ID: LMW-4-1114  
SAMPLE

Lab Sample ID: ZS69A  
LIMS ID: 15-518  
Matrix: Water  
Date Analyzed: 01/21/15 17:13

QC Report No: ZS69-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 62.0% | 2-Fluorobiphenyl       | 57.2% |
| d14-p-Terphenyl      | 67.6% | d4-1,2-Dichlorobenzene | 56.8% |
| d5-Phenol            | 83.5% | 2-Fluorophenol         | 77.9% |
| 2,4,6-Tribromophenol | 85.1% | d4-2-Chlorophenol      | 82.7% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: MB-011515  
METHOD BLANK

Lab Sample ID: MB-011515  
LIMS ID: 15-518  
Matrix: Water  
Data Release Authorized: *AB*  
Reported: 01/22/15

QC Report No: ZS69-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: NA  
Date Received: NA

Date Extracted: 01/15/15  
Date Analyzed: 01/21/15 15:31  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

Sample ID: MB-011515  
METHOD BLANK

Lab Sample ID: MB-011515  
LIMS ID: 15-518  
Matrix: Water  
Date Analyzed: 01/21/15 15:31

QC Report No: ZS69-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 60.4% | 2-Fluorobiphenyl       | 56.8% |
| d14-p-Terphenyl      | 69.6% | d4-1,2-Dichlorobenzene | 52.0% |
| d5-Phenol            | 85.6% | 2-Fluorophenol         | 75.5% |
| 2,4,6-Tribromophenol | 82.9% | d4-2-Chlorophenol      | 80.8% |

**SW8270 SEMIVOLATILES WATER SURROGATE RECOVERY SUMMARY**

Matrix: Water

QC Report No: ZS69-Golder Associates

Project: Landsburg

923-1000-002.R273

| Client ID   | NBZ   | FBP   | TPH   | DCB   | PHL   | 2FP   | TBP   | 2CP   | TOT | OUT |
|-------------|-------|-------|-------|-------|-------|-------|-------|-------|-----|-----|
| MB-011515   | 60.4% | 56.8% | 69.6% | 52.0% | 85.6% | 75.5% | 82.9% | 80.8% | 0   |     |
| LCS-011515  | 61.6% | 61.2% | 68.4% | 53.2% | 91.2% | 72.5% | 104%  | 83.2% | 0   |     |
| LCSD-011515 | 64.8% | 64.0% | 64.8% | 55.2% | 96.5% | 79.7% | 105%  | 88.3% | 0   |     |
| LMW-4-1114  | 62.0% | 57.2% | 67.6% | 56.8% | 83.5% | 77.9% | 85.1% | 82.7% | 0   |     |

|                                | LCS/MB LIMITS | QC LIMITS |
|--------------------------------|---------------|-----------|
| (NBZ) = d5-Nitrobenzene        | (27-120)      | (27-120)  |
| (FBP) = 2-Fluorobiphenyl       | (33-120)      | (33-120)  |
| (TPH) = d14-p-Terphenyl        | (28-130)      | (28-130)  |
| (DCB) = d4-1,2-Dichlorobenzene | (20-120)      | (20-120)  |
| (PHL) = d5-Phenol              | (38-120)      | (38-120)  |
| (2FP) = 2-Fluorophenol         | (33-120)      | (33-120)  |
| (TBP) = 2,4,6-Tribromophenol   | (52-131)      | (52-131)  |
| (2CP) = d4-2-Chlorophenol      | (41-120)      | (41-120)  |

Prep Method: SW3520C  
Log Number Range: 15-518 to 15-518

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Page 1 of 3



Sample ID: LCS-011515  
LCS/LCSD

Lab Sample ID: LCS-011515  
LIMS ID: 15-518  
Matrix: Water  
Data Release Authorized: *[Signature]*  
Reported: 01/22/15

QC Report No: ZS69-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/18/14  
Date Received: 11/19/14

Date Extracted LCS/LCSD: 01/15/15

Sample Amount LCS: 500 mL

Date Analyzed LCS: 01/21/15 16:05  
LCSD: 01/21/15 16:39

Final Extract Volume LCS: 0.50 mL  
LCSD: 0.50 mL

Instrument/Analyst LCS: NT6/JZ  
LCSD: NT6/JZ

Dilution Factor LCS: 1.00  
LCSD: 1.00

GPC Cleanup: NO

| Analyte                      | LCS  | Spike Added-LCS | LCS Recovery | LCS  | Spike Added-LCSD | LCSD Recovery | RPD   |
|------------------------------|------|-----------------|--------------|------|------------------|---------------|-------|
| Phenol                       | 15.4 | 25.0            | 61.6%        | 15.4 | 25.0             | 61.6%         | 7.5%  |
| Bis-(2-Chloroethyl) Ether    | 15.1 | 25.0            | 60.4%        | 15.4 | 25.0             | 61.6%         | 2.0%  |
| 2-Chlorophenol               | 14.4 | 25.0            | 57.6%        | 15.7 | 25.0             | 62.8%         | 8.6%  |
| 1,3-Dichlorobenzene          | 10.4 | 25.0            | 41.6%        | 11.4 | 25.0             | 45.6%         | 9.2%  |
| 1,4-Dichlorobenzene          | 10.8 | 25.0            | 43.2%        | 12.1 | 25.0             | 48.4%         | 11.4% |
| Benzyl Alcohol               | 13.2 | 25.0            | 52.8%        | 14.5 | 25.0             | 58.0%         | 9.4%  |
| 1,2-Dichlorobenzene          | 11.2 | 25.0            | 44.8%        | 12.2 | 25.0             | 48.8%         | 8.5%  |
| 2-Methylphenol               | 14.0 | 25.0            | 56.0%        | 15.0 | 25.0             | 60.0%         | 6.9%  |
| 2,2'-Oxybis(1-Chloropropane) | 13.0 | 25.0            | 52.0%        | 13.7 | 25.0             | 54.8%         | 5.2%  |
| 4-Methylphenol               | 15.0 | 25.0            | 60.0%        | 16.1 | 25.0             | 64.4%         | 7.1%  |
| N-Nitroso-Di-N-Propylamine   | 14.8 | 25.0            | 59.2%        | 15.9 | 25.0             | 63.6%         | 7.2%  |
| Hexachloroethane             | 10.5 | 25.0            | 42.0%        | 11.5 | 25.0             | 46.0%         | 9.1%  |
| Nitrobenzene                 | 14.4 | 25.0            | 57.6%        | 15.7 | 25.0             | 62.8%         | 8.6%  |
| Isophorone                   | 14.6 | 25.0            | 58.4%        | 15.6 | 25.0             | 62.4%         | 6.6%  |
| 2-Nitrophenol                | 16.4 | 25.0            | 65.6%        | 17.6 | 25.0             | 70.4%         | 7.1%  |
| 2,4-Dimethylphenol           | 48.4 | 75.0            | 64.5%        | 50.3 | 75.0             | 67.1%         | 3.9%  |
| Benzoic Acid                 | 117  | 138             | 84.8%        | 121  | 138              | 87.7%         | 3.4%  |
| bis(2-Chloroethoxy) Methane  | 14.7 | 25.0            | 58.8%        | 15.7 | 25.0             | 62.8%         | 6.6%  |
| 2,4-Dichlorophenol           | 51.4 | 75.0            | 68.5%        | 55.3 | 75.0             | 73.7%         | 7.3%  |
| 1,2,4-Trichlorobenzene       | 12.0 | 25.0            | 48.0%        | 13.2 | 25.0             | 52.8%         | 9.5%  |
| Naphthalene                  | 14.4 | 25.0            | 57.6%        | 15.7 | 25.0             | 62.8%         | 8.6%  |
| 4-Chloroaniline              | 65.5 | 75.0            | 87.3%        | 59.6 | 75.0             | 79.5%         | 9.4%  |
| Hexachlorobutadiene          | 9.7  | 25.0            | 38.8%        | 11.2 | 25.0             | 44.8%         | 14.4% |
| 4-Chloro-3-methylphenol      | 57.5 | 75.0            | 76.7%        | 61.2 | 75.0             | 81.6%         | 6.2%  |
| 2-Methylnaphthalene          | 15.0 | 25.0            | 60.0%        | 16.1 | 25.0             | 64.4%         | 7.1%  |
| Hexachlorocyclopentadiene    | 24.7 | 75.0            | 32.9%        | 27.6 | 75.0             | 36.8%         | 11.1% |
| 2,4,6-Trichlorophenol        | 52.0 | 75.0            | 69.3%        | 56.4 | 75.0             | 75.2%         | 8.1%  |
| 2,4,5-Trichlorophenol        | 53.0 | 75.0            | 70.7%        | 55.3 | 75.0             | 73.7%         | 4.2%  |
| 2-Chloronaphthalene          | 13.9 | 25.0            | 55.6%        | 15.2 | 25.0             | 60.8%         | 8.9%  |
| 2-Nitroaniline               | 64.2 | 75.0            | 85.6%        | 67.9 | 75.0             | 90.5%         | 5.6%  |

**ORGANICS ANALYSIS DATA SHEET**  
**Semivolatiles by SW8270D GC/MS**  
Page 2 of 3



Sample ID: LCS-011515  
LCS/LCSD

Lab Sample ID: LCS-011515  
LIMS ID: 15-518  
Matrix: Water  
Date Analyzed LCS: 01/21/15 16:05  
LCSD: 01/21/15 16:39

QC Report No: ZS69-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| Analyte                    | LCS    | Spike Added-LCS | LCS Recovery | LCSD   | Spike Added-LCSD | LCSD Recovery | RPD   |
|----------------------------|--------|-----------------|--------------|--------|------------------|---------------|-------|
| Dimethylphthalate          | 18.3   | 25.0            | 73.2%        | 18.5   | 25.0             | 74.0%         | 1.1%  |
| Acenaphthylene             | 16.3   | 25.0            | 65.2%        | 17.5   | 25.0             | 70.0%         | 7.1%  |
| 3-Nitroaniline             | 78.6 Q | 75.0            | 105%         | 74.9 Q | 75.0             | 99.9%         | 4.8%  |
| Acenaphthene               | 16.1   | 25.0            | 64.4%        | 17.1   | 25.0             | 68.4%         | 6.0%  |
| 2,4-Dinitrophenol          | 117    | 138             | 84.8%        | 128    | 138              | 92.8%         | 9.0%  |
| 4-Nitrophenol              | 67.8   | 75.0            | 90.4%        | 68.2   | 75.0             | 90.9%         | 0.6%  |
| Dibenzofuran               | 17.0   | 25.0            | 68.0%        | 18.3   | 25.0             | 73.2%         | 7.4%  |
| 2,6-Dinitrotoluene         | 56.2   | 75.0            | 74.9%        | 59.3   | 75.0             | 79.1%         | 5.4%  |
| 2,4-Dinitrotoluene         | 58.8   | 75.0            | 78.4%        | 62.6   | 75.0             | 83.5%         | 6.3%  |
| Diethylphthalate           | 18.9   | 25.0            | 75.6%        | 19.0   | 25.0             | 76.0%         | 0.5%  |
| 4-Chlorophenyl-phenylether | 15.5   | 25.0            | 62.0%        | 16.4   | 25.0             | 65.6%         | 5.6%  |
| Fluorene                   | 16.4   | 25.0            | 65.6%        | 17.5   | 25.0             | 70.0%         | 6.5%  |
| 4-Nitroaniline             | 76.5   | 75.0            | 102%         | 77.8   | 75.0             | 104%          | 1.7%  |
| 4,6-Dinitro-2-Methylphenol | 120    | 138             | 87.0%        | 125    | 138              | 90.6%         | 4.1%  |
| N-Nitrosodiphenylamine     | 15.8   | 25.0            | 63.2%        | 15.8   | 25.0             | 63.2%         | 0.0%  |
| 4-Bromophenyl-phenylether  | 16.4   | 25.0            | 65.6%        | 16.9   | 25.0             | 67.6%         | 3.0%  |
| Hexachlorobenzene          | 16.2   | 25.0            | 64.8%        | 16.7   | 25.0             | 66.8%         | 3.0%  |
| Pentachlorophenol          | 44.2 Q | 75.0            | 58.9%        | 49.0 Q | 75.0             | 65.3%         | 10.3% |
| Phenanthrene               | 17.4   | 25.0            | 69.6%        | 17.6   | 25.0             | 70.4%         | 1.1%  |
| Carbazole                  | 18.1   | 25.0            | 72.4%        | 18.0   | 25.0             | 72.0%         | 0.6%  |
| Anthracene                 | 17.9   | 25.0            | 71.6%        | 18.2   | 25.0             | 72.8%         | 1.7%  |
| Di-n-Butylphthalate        | 19.3   | 25.0            | 77.2%        | 18.6   | 25.0             | 74.4%         | 3.7%  |
| Fluoranthene               | 20.5   | 25.0            | 82.0%        | 20.4   | 25.0             | 81.6%         | 0.5%  |
| Pyrene                     | 15.8   | 25.0            | 63.2%        | 15.7   | 25.0             | 62.8%         | 0.6%  |
| Butylbenzylphthalate       | 17.8   | 25.0            | 71.2%        | 17.2   | 25.0             | 68.8%         | 3.4%  |
| 3,3'-Dichlorobenzidine     | 57.7   | 75.0            | 76.9%        | 52.4   | 75.0             | 69.9%         | 9.6%  |
| Benzo(a)anthracene         | 18.9   | 25.0            | 75.6%        | 18.2   | 25.0             | 72.8%         | 3.8%  |
| bis(2-Ethylhexyl)phthalate | 18.1   | 25.0            | 72.4%        | 17.3   | 25.0             | 69.2%         | 4.5%  |
| Chrysene                   | 19.3   | 25.0            | 77.2%        | 18.8   | 25.0             | 75.2%         | 2.6%  |
| Di-n-Octyl phthalate       | 19.1   | 25.0            | 76.4%        | 18.3   | 25.0             | 73.2%         | 4.3%  |
| Benzo(b)fluoranthene       | 19.0   | 25.0            | 76.0%        | 19.0   | 25.0             | 76.0%         | 0.0%  |
| Benzo(k)fluoranthene       | 21.8   | 25.0            | 87.2%        | 20.3   | 25.0             | 81.2%         | 7.1%  |
| Benzo(a)pyrene             | 20.1   | 25.0            | 80.4%        | 19.2   | 25.0             | 76.8%         | 4.6%  |
| Indeno(1,2,3-cd)pyrene     | 15.1   | 25.0            | 60.4%        | 14.5   | 25.0             | 58.0%         | 4.1%  |
| Dibenz(a,h)anthracene      | 15.5   | 25.0            | 62.0%        | 15.0   | 25.0             | 60.0%         | 3.3%  |
| Benzo(g,h,i)perylene       | 13.5 Q | 25.0            | 54.0%        | 13.0 Q | 25.0             | 52.0%         | 3.8%  |
| 3&4-Methylphenol           | 15.0   | 25.0            | 60.0%        | 16.1   | 25.0             | 64.4%         | 7.1%  |
| 1-Methylnaphthalene        | 15.4   | 25.0            | 61.6%        | 16.8   | 25.0             | 67.2%         | 8.7%  |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Page 3 of 3



Sample ID: LCS-011515  
LCS/LCSD

Lab Sample ID: LCS-011515

LIMS ID: 15-518

Matrix: Water

Date Analyzed LCS: 01/21/15 16:05

LCSD: 01/21/15 16:39

QC Report No: ZS69-Golder Associates

Project: Landsburg

923-1000-002.R273

| Analyte                  | LCS  | Spike<br>Added-LCS | LCS<br>Recovery | LCSD | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD  |
|--------------------------|------|--------------------|-----------------|------|---------------------|------------------|------|
| Total Benzofluoranthenes | 41.3 | 50.0               | 82.6%           | 39.5 | 50.0                | 79.0%            | 4.5% |

**Semivolatile Surrogate Recovery**

|                        | LCS   | LCSD  |
|------------------------|-------|-------|
| d5-Nitrobenzene        | 61.6% | 64.8% |
| 2-Fluorobiphenyl       | 61.2% | 64.0% |
| d14-p-Terphenyl        | 68.4% | 64.8% |
| d4-1,2-Dichlorobenzene | 53.2% | 55.2% |
| d5-Phenol              | 91.2% | 96.5% |
| 2-Fluorophenol         | 72.5% | 79.7% |
| 2,4,6-Tribromophenol   | 104%  | 105%  |
| d4-2-Chlorophenol      | 83.2% | 88.3% |

Results reported in µg/L

RPD calculated using sample concentrations per SW846.



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants

February 19, 2015

Doug Morell  
Golder Associates Inc.  
18300 NE Union Hill Road, Suite 200  
Redmond, WA 98052-3333

**Client Project Name: Landsburg Mine**  
**Client Project Number: 923-1000-002.R273**  
**ARI ID: ZW42**

Dear Mr. Morell:

Please find enclosed Chain-of-Custody (COC) record, sample receipt documentation, and the final results for the project referenced above. Analytical Resources, Inc. (ARI) accepted one water sample in good condition on February 12, 2015. There were no discrepancies between the COC and the sample containers' labels.

The sample was analyzed for SVOCs, as requested on the COC.

No analytical complications were noted.

An electronic copy of this report and all supporting raw data will remain on file at ARI. Please feel free to contact me if you have any questions or require any additional information.

Respectfully,

ANALYTICAL RESOURCES, INC.

  
Kelly Bottem  
Client Services Manager  
(206) 695-6211  
[kellyb@arilabs.com](mailto:kellyb@arilabs.com)

# Chain of Custody Record & Laboratory Analysis Request

|                                                                                                                                                                                   |                |                                           |          |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  |                |
|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----------------|-------------------------------------------|----------|---------------------------------------------------------------------------------------------------------------------|--|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|--|---------------------------------|--|-----------------------------|--|--------------------------------------------------|--|----------------|
| ARI Assigned Number:<br><b>2012</b>                                                                                                                                               |                | Turn-around Requested:<br><b>Standard</b> |          | Page: <b>1</b> of <b>1</b>                                                                                          |  |  <b>Analytical Resources, Incorporated</b><br>Analytical Chemists and Consultants<br>4611 South 134th Place, Suite 100<br>Tukwila, WA 98168<br>206-695-6200 206-695-6201 (fax) |  |                                 |  |                             |  |                                                  |  |                |
| ARI Client Company:<br><b>Golder Associates</b>                                                                                                                                   |                | Phone:<br><b>425-883-0777</b>             |          | Date:<br><b>2/12/15</b>                                                                                             |  |                                                                                                                                                                                                                                                                  |  | Ice Present? <b>Y</b>           |  |                             |  |                                                  |  |                |
| Client Contact:<br><b>Doug Morell, Jill Lamberts</b>                                                                                                                              |                |                                           |          | No. of Coolers: <b>1</b>                                                                                            |  |                                                                                                                                                                                                                                                                  |  | Cooler Temps: <b>5.7</b>        |  |                             |  |                                                  |  |                |
| Client Project Name:<br><b>Landsburg</b>                                                                                                                                          |                |                                           |          | Analysis Requested                                                                                                  |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  | Notes/Comments                                   |  |                |
| Client Project #:<br><b>923-1000-002-R273</b>                                                                                                                                     |                | Samplers:<br><b>Lamberts, Hunt</b>        |          | BEHP BY<br>8230 D                                                                                                   |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  | Pls analyze under existing MSA btwn Golder + ARI |  |                |
| Sample ID                                                                                                                                                                         | Date           | Time                                      | Matrix   |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  | No. Containers |
| <b>LMW-4-0215</b>                                                                                                                                                                 | <b>2/12/15</b> | <b>1450</b>                               | <b>W</b> |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  | <b>4</b>       |
|                                                                                                                                                                                   |                |                                           |          |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          |                                                                                                                     |  |                                                                                                                                                                                                                                                                  |  |                                 |  |                             |  |                                                  |  |                |
| Comments/Special Instructions<br><b>- Ecology Eim EDD</b><br><b>**Client specific RLS and analyze first**</b><br><b>Pls CC j.lamberts@golder.com</b><br><b>dmorell@golder.com</b> |                |                                           |          | Relinquished by:<br>(Signature)  |  | Received by:<br>(Signature)                                                                                                                                                  |  | Relinquished by:<br>(Signature) |  | Received by:<br>(Signature) |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          | Printed Name:<br><b>J. Lamberts</b>                                                                                 |  | Printed Name:<br><b>A. Volgardsen</b>                                                                                                                                                                                                                            |  | Printed Name:                   |  | Printed Name:               |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          | Company:<br><b>Golder</b>                                                                                           |  | Company:<br><b>ARI</b>                                                                                                                                                                                                                                           |  | Company:                        |  | Company:                    |  |                                                  |  |                |
|                                                                                                                                                                                   |                |                                           |          | Date & Time:<br><b>2/12/15 1607</b>                                                                                 |  | Date & Time:<br><b>2/12/15 1607</b>                                                                                                                                                                                                                              |  | Date & Time:                    |  | Date & Time:                |  |                                                  |  |                |

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

2012:000002

# Sample ID Cross Reference Report



ARI Job No: ZW42  
Client: Golder Associates  
Project Event: 923-1000-002-R273  
Project Name: Landsburg

| Sample ID     | ARI<br>Lab ID | ARI<br>LIMS ID | Matrix | Sample Date/Time | VTSR           |
|---------------|---------------|----------------|--------|------------------|----------------|
| 1. LMW-4-0215 | ZW42A         | 15-2811        | Water  | 02/12/15 14:50   | 02/12/15 16:07 |



# Cooler Receipt Form

ARI Client: Goldier

Project Name: Landsburg

COC No(s): \_\_\_\_\_ (NA)

Delivered by: Fed-Ex UPS Courier Hand Delivered Other: \_\_\_\_\_

Assigned ARI Job No. 2W42

Tracking No: \_\_\_\_\_ (NA)

## Preliminary Examination Phase:

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES (NO)

Were custody papers included with the cooler? ..... (YES) NO

Were custody papers properly filled out (ink, signed, etc) ..... (YES) NO

Temperature of Cooler(s) (°C) (recommended 2 0-6.0 °C for chemistry)

Time: 1607 5.7 Temp Gun ID#: 90877952

If cooler temperature is out of compliance fill out form 00070F

Cooler Accepted by A Date: 2/12/15 Time: 1607

**Complete custody forms and attach all shipping documents**

## Log-In Phase:

Was a temperature blank included in the cooler? ..... YES (NO)

What kind of packing material was used? ... Bubble Wrap Wet Ice Gel Packs (Baggies) Foam Block Paper Other: \_\_\_\_\_

Was sufficient ice used (if appropriate)? ..... NA (YES) NO

Were all bottles sealed in individual plastic bags? ..... (YES) NO

Did all bottles arrive in good condition (unbroken)? ..... (YES) NO

Were all bottle labels complete and legible? ..... (YES) NO

Did the number of containers listed on COC match with the number of containers received? ..... (YES) NO

Did all bottle labels and tags agree with custody papers? ..... (YES) NO

Were all bottles used correct for the requested analyses? ..... (YES) NO

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... (NA) YES NO

Were all VOC vials free of air bubbles? ..... (NA) YES NO

Was sufficient amount of sample sent in each bottle? .. (YES) NO

Date VOC Trip Blank was made at ARI ..... (NA)

Was Sample Split by ARI : (NA) YES Date/Time: \_\_\_\_\_ Equipment: \_\_\_\_\_ Split by: \_\_\_\_\_

Samples Logged by: AV Date: 2/12/15 Time: 1720

**\*\* Notify Project Manager of discrepancies or concerns \*\***

| Sample ID on Bottle | Sample ID on COC | Sample ID on Bottle | Sample ID on COC |
|---------------------|------------------|---------------------|------------------|
|                     |                  |                     |                  |
|                     |                  |                     |                  |
|                     |                  |                     |                  |
|                     |                  |                     |                  |

**Additional Notes, Discrepancies, & Resolutions:**

By: \_\_\_\_\_ Date: \_\_\_\_\_

|                                    |                              |                                        |                                                                                                                                                    |
|------------------------------------|------------------------------|----------------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------|
| <p>Small Air Bubbles<br/>- 2mm</p> | <p>Peabubbles<br/>2-4 mm</p> | <p>LARGE Air Bubbles<br/>&gt; 4 mm</p> | <p>Small → "sm" (&lt; 2 mm)</p> <p>Peabubbles → "pb" (2 to &lt; 4 mm)</p> <p>Large → "lg" (4 to &lt; 6 mm)</p> <p>Headspace → "hs" (&gt; 6 mm)</p> |
|------------------------------------|------------------------------|----------------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------|

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 1

Sample ID: LMW-4-0215  
SAMPLE



Lab Sample ID: ZW42A  
LIMS ID: 15-2811  
Matrix: Water  
Data Release Authorized: *[Signature]*  
Reported: 02/18/15

QC Report No: ZW42-Golder Associates  
Project: Landsburg  
923-1000-002-R273  
Date Sampled: 02/12/15  
Date Received: 02/12/15

Date Extracted: 02/16/15  
Date Analyzed: 02/18/15 12:04  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                    | DL  | LOQ | Result  |
|------------|----------------------------|-----|-----|---------|
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1 | 3.0 | < 3.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 64.0% | 2-Fluorobiphenyl       | 64.0% |
| d14-p-Terphenyl      | 80.0% | d4-1,2-Dichlorobenzene | 59.6% |
| d5-Phenol            | 60.8% | 2-Fluorophenol         | 61.1% |
| 2,4,6-Tribromophenol | 69.6% | d4-2-Chlorophenol      | 65.3% |

**ORGANICS ANALYSIS DATA SHEET**  
**Semivolatiles by SW8270D GC/MS**  
**Extraction Method: SW3520C**  
 Page 1 of 1

**Sample ID: MB-021615**  
**METHOD BLANK**

Lab Sample ID: MB-021615  
 LIMS ID: 15-2811  
 Matrix: Water  
 Data Release Authorized: *AB*  
 Reported: 02/18/15

QC Report No: ZW42-Golder Associates  
 Project: Landsburg  
 923-1000-002-R273  
 Date Sampled: NA  
 Date Received: NA

Date Extracted: 02/16/15  
 Date Analyzed: 02/18/15 09:51  
 Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1.00

| CAS Number | Analyte                    | DL  | LOQ | Result  |
|------------|----------------------------|-----|-----|---------|
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1 | 3.0 | < 3.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 60.0% | 2-Fluorobiphenyl       | 58.8% |
| d14-p-Terphenyl      | 88.0% | d4-1,2-Dichlorobenzene | 55.2% |
| d5-Phenol            | 56.0% | 2-Fluorophenol         | 54.1% |
| 2,4,6-Tribromophenol | 60.8% | d4-2-Chlorophenol      | 61.1% |

**SW8270 SEMIVOLATILES WATER SURROGATE RECOVERY SUMMARY**

Matrix: Water

QC Report No: ZW42-Golder Associates

Project: Landsburg

923-1000-002-R273

| Client ID   | NBZ   | FBP   | TPH   | DCB   | PHL   | 2FP   | TBP   | 2CP   | TOT | OUT |
|-------------|-------|-------|-------|-------|-------|-------|-------|-------|-----|-----|
| MB-021615   | 60.0% | 58.8% | 88.0% | 55.2% | 56.0% | 54.1% | 60.8% | 61.1% | 0   |     |
| LCS-021615  | 72.4% | 70.0% | 82.8% | 62.8% | 75.5% | 66.7% | 83.5% | 73.3% | 0   |     |
| LCSD-021615 | 56.8% | 55.6% | 76.8% | 52.0% | 57.3% | 52.5% | 69.9% | 58.4% | 0   |     |
| LMW-4-0215  | 64.0% | 64.0% | 80.0% | 59.6% | 60.8% | 61.1% | 69.6% | 65.3% | 0   |     |

|                                | LCS/MB LIMITS | QC LIMITS |
|--------------------------------|---------------|-----------|
| (NBZ) = d5-Nitrobenzene        | (27-120)      | (27-120)  |
| (FBP) = 2-Fluorobiphenyl       | (33-120)      | (33-120)  |
| (TPH) = d14-p-Terphenyl        | (28-130)      | (28-130)  |
| (DCB) = d4-1,2-Dichlorobenzene | (20-120)      | (20-120)  |
| (PHL) = d5-Phenol              | (38-120)      | (38-120)  |
| (2FP) = 2-Fluorophenol         | (33-120)      | (33-120)  |
| (TBP) = 2,4,6-Tribromophenol   | (52-131)      | (52-131)  |
| (2CP) = d4-2-Chlorophenol      | (41-120)      | (41-120)  |

Prep Method: SW3520C  
Log Number Range: 15-2811 to 15-2811

**ORGANICS ANALYSIS DATA SHEET**  
Semivolatiles by SW8270D GC/MS  
Page 1 of 1

Sample ID: LCS-021615  
LCS/LCSD

Lab Sample ID: LCS-021615

LIMS ID: 15-2811

Matrix: Water

Data Release Authorized: *AB*

Reported: 02/18/15

QC Report No: ZW42-Golder Associates

Project: Landsburg

923-1000-002-R273

Date Sampled: 02/12/15

Date Received: 02/12/15

Date Extracted LCS/LCSD: 02/16/15

Sample Amount LCS: 500 mL

LCSD: 500 mL

Date Analyzed LCS: 02/18/15 10:24

Final Extract Volume LCS: 0.50 mL

LCSD: 02/18/15 10:58

LCSD: 0.50 mL

Instrument/Analyst LCS: NT6/JZ

Dilution Factor LCS: 1.00

LCSD: NT6/JZ

LCSD: 1.00

GPC Cleanup: NO

| Analyte                    | LCS  | Spike<br>Added-LCS | LCS<br>Recovery | LCSD | Spike<br>Added-LCSD | LCSD<br>Recovery | RPD  |
|----------------------------|------|--------------------|-----------------|------|---------------------|------------------|------|
| bis(2-Ethylhexyl)phthalate | 21.2 | 25.0               | 84.8%           | 20.3 | 25.0                | 81.2%            | 4.3% |

**Semivolatile Surrogate Recovery**

|                        | LCS   | LCSD  |
|------------------------|-------|-------|
| d5-Nitrobenzene        | 72.4% | 56.8% |
| 2-Fluorobiphenyl       | 70.0% | 55.6% |
| d14-p-Terphenyl        | 82.8% | 76.8% |
| d4-1,2-Dichlorobenzene | 62.8% | 52.0% |
| d5-Phenol              | 75.5% | 57.3% |
| 2-Fluorophenol         | 66.7% | 52.5% |
| 2,4,6-Tribromophenol   | 83.5% | 69.9% |
| d4-2-Chlorophenol      | 73.3% | 58.4% |

Results reported in µg/L

RPD calculated using sample concentrations per SW846.

Analytical Resources, Inc.

Semivolatiles Report SW846 Method 8270D

Data file : /chem2/nt6.i/20150218.b/02181506.d  
Lab Smp Id: ZW42A Client Smp ID: LMW-4-0215  
Inj Date : 18-FEB-2015 12:04  
Operator : JZ Inst ID: nt6.i  
Smp Info : ZW42A  
Misc Info : 15-2811  
Comment : 1ul Injection  
Method : /chem2/nt6.i/20150218.b/SW846021115.m  
Meth Date : 18-Feb-2015 14:17 jianqing Quant Type: ISTD  
Cal Date : 11-FEB-2015 12:58 Cal File: 02111502.d  
Als bottle: 6  
Dil Factor: 1.00000  
Integrator: HP RTE Compound Sublist: LL.sub  
Target Version: 3.50

Concentration Formula: Amt \* DF \* Vt/Vo \* CpndVariable

*02/18/15*

| Name | Value     | Description                     |
|------|-----------|---------------------------------|
| DF   | 1.00000   | Dilution Factor                 |
| Vt   | 500.00000 | Volume of final extract (uL)    |
| Vo   | 500.00000 | Volume of sample extracted (mL) |

Cpnd Variable

Local Compound Variable

| Compounds                       | QUANT SIG | CONCENTRATIONS |                        |        |         |          |                      |                 |
|---------------------------------|-----------|----------------|------------------------|--------|---------|----------|----------------------|-----------------|
|                                 |           | MASS           | RT                     | EXP RT | REL RT  | RESPONSE | ON-COLUMN<br>(ug/mL) | FINAL<br>(ug/L) |
| =====                           | =====     | =====          | ==                     | =====  | =====   | =====    | =====                | =====           |
| \$ 1 2-Fluorophenol             | 112       |                | 6.487                  | 6.491  | (0.769) | 263067   | 22.9480              | 22.95           |
| \$ 2 Phenol-d5                  | 99        |                | 7.994                  | 7.998  | (0.948) | 315823   | 22.7951              | 22.80           |
| 3 Phenol                        | 94        |                | Compound Not Detected. |        |         |          |                      |                 |
| \$ 5 2-Chlorophenol-d4          | 132       |                | 8.133                  | 8.137  | (0.965) | 323374   | 24.5232              | 24.52           |
| 4 Bis(2-Chloroethyl) ether      | 93        |                | Compound Not Detected. |        |         |          |                      |                 |
| 6 2-Chlorophenol                | 128       |                | Compound Not Detected. |        |         |          |                      |                 |
| 7 1,3-Dichlorobenzene           | 146       |                | Compound Not Detected. |        |         |          |                      |                 |
| * 8 1,4-Dichlorobenzene-d4      | 152       |                | 8.432                  | 8.431  | (1.000) | 222362   | 20.0000              |                 |
| 9 1,4-Dichlorobenzene           | 146       |                | Compound Not Detected. |        |         |          |                      |                 |
| \$ 10 1,2-Dichlorobenzene-d4    | 152       |                | 8.731                  | 8.735  | (1.035) | 158523   | 14.8776              | 14.88           |
| 12 1,2-Dichlorobenzene          | 146       |                | Compound Not Detected. |        |         |          |                      |                 |
| 11 Benzyl alcohol               | 108       |                | Compound Not Detected. |        |         |          |                      |                 |
| 14 2,2'-oxybis(1-Chloropropane) | 45        |                | Compound Not Detected. |        |         |          |                      |                 |
| 13 2-Methylphenol               | 108       |                | Compound Not Detected. |        |         |          |                      |                 |
| 17 Hexachloroethane             | 117       |                | Compound Not Detected. |        |         |          |                      |                 |
| 16 N-Nitroso-di-n-propylamine   | 70        |                | Compound Not Detected. |        |         |          |                      |                 |

| Compounds                     | QUANT SIG | CONCENTRATIONS |        |         |                        |         |         |
|-------------------------------|-----------|----------------|--------|---------|------------------------|---------|---------|
|                               |           | ON-COLUMN      | FINAL  |         |                        |         |         |
|                               | MASS      | RT             | EXP RT | REL RT  | RESPONSE               | (ug/mL) | ( ug/L) |
| =====                         | =====     | ==             | =====  | =====   | =====                  | =====   | =====   |
| 15 4-Methylphenol             | 108       |                |        |         | Compound Not Detected. |         |         |
| \$ 18 Nitrobenzene-d5         | 82        | 9.356          | 9.360  | (0.893) | 223955                 | 16.0418 | 16.04   |
| 19 Nitrobenzene               | 77        |                |        |         | Compound Not Detected. |         |         |
| 20 Isophorone                 | 82        |                |        |         | Compound Not Detected. |         |         |
| 21 2-Nitrophenol              | 139       |                |        |         | Compound Not Detected. |         |         |
| 22 2,4-Dimethylphenol         | 107       |                |        |         | Compound Not Detected. |         |         |
| 23 Bis(2-Chloroethoxy)methane | 93        |                |        |         | Compound Not Detected. |         |         |
| 24 Benzoic acid               | 105       |                |        |         | Compound Not Detected. |         |         |
| 25 2,4-Dichlorophenol         | 162       |                |        |         | Compound Not Detected. |         |         |
| 26 1,2,4-Trichlorobenzene     | 180       |                |        |         | Compound Not Detected. |         |         |
| * 27 Naphthalene-d8           | 136       | 10.472         | 10.477 | (1.000) | 783563                 | 20.0000 |         |
| 28 Naphthalene                | 128       |                |        |         | Compound Not Detected. |         |         |
| 29 4-Chloroaniline            | 127       |                |        |         | Compound Not Detected. |         |         |
| 30 Hexachlorobutadiene        | 225       |                |        |         | Compound Not Detected. |         |         |
| 31 4-Chloro-3-methylphenol    | 107       |                |        |         | Compound Not Detected. |         |         |
| 32 2-Methylnaphthalene        | 141       |                |        |         | Compound Not Detected. |         |         |
| 33 Hexachlorocyclopentadiene  | 237       |                |        |         | Compound Not Detected. |         |         |
| 34 2,4,6-Trichlorophenol      | 196       |                |        |         | Compound Not Detected. |         |         |
| 35 2,4,5-Trichlorophenol      | 196       |                |        |         | Compound Not Detected. |         |         |
| \$ 36 2-Fluorobiphenyl        | 172       | 12.262         | 12.266 | (0.919) | 542232                 | 16.0443 | 16.04   |
| 37 2-Chloronaphthalene        | 162       |                |        |         | Compound Not Detected. |         |         |
| 38 2-Nitroaniline             | 65        |                |        |         | Compound Not Detected. |         |         |
| 39 Dimethylphthalate          | 163       |                |        |         | Compound Not Detected. |         |         |
| 40 Acenaphthylene             | 152       |                |        |         | Compound Not Detected. |         |         |
| 41 2,6-Dinitrotoluene         | 165       |                |        |         | Compound Not Detected. |         |         |
| * 42 Acenaphthene-d10         | 164       | 13.336         | 13.340 | (1.000) | 537749                 | 20.0000 |         |
| 43 3-Nitroaniline             | 138       |                |        |         | Compound Not Detected. |         |         |
| 44 Acenaphthene               | 153       |                |        |         | Compound Not Detected. |         |         |
| 45 2,4-Dinitrophenol          | 184       |                |        |         | Compound Not Detected. |         |         |
| 46 Dibenzofuran               | 168       |                |        |         | Compound Not Detected. |         |         |
| 47 4-Nitrophenol              | 109       |                |        |         | Compound Not Detected. |         |         |
| 48 2,4-Dinitrotoluene         | 165       |                |        |         | Compound Not Detected. |         |         |
| 50 Diethylphthalate           | 149       |                |        |         | Compound Not Detected. |         |         |
| 49 Fluorene                   | 166       |                |        |         | Compound Not Detected. |         |         |
| 51 4-Chlorophenyl-phenylether | 204       |                |        |         | Compound Not Detected. |         |         |
| 52 4-Nitroaniline             | 138       |                |        |         | Compound Not Detected. |         |         |
| 53 4,6-Dinitro-2-methylphenol | 198       |                |        |         | Compound Not Detected. |         |         |
| 54 N-Nitrosodiphenylamine     | 169       |                |        |         | Compound Not Detected. |         |         |
| \$ 55 2,4,6-Tribromophenol    | 330       | 14.629         | 14.638 | (1.097) | 162645                 | 26.1491 | 26.15   |
| 56 4-Bromophenyl-phenylether  | 248       |                |        |         | Compound Not Detected. |         |         |
| 57 Hexachlorobenzene          | 284       |                |        |         | Compound Not Detected. |         |         |
| 58 Pentachlorophenol          | 266       |                |        |         | Compound Not Detected. |         |         |
| * 59 Phenanthrene-d10         | 188       | 15.718         | 15.723 | (1.000) | 973135                 | 20.0000 |         |
| 60 Phenanthrene               | 178       |                |        |         | Compound Not Detected. |         |         |
| 61 Anthracene                 | 178       |                |        |         | Compound Not Detected. |         |         |
| 62 Carbazole                  | 167       |                |        |         | Compound Not Detected. |         |         |
| 63 Di-n-butylphthalate        | 149       |                |        |         | Compound Not Detected. |         |         |

| Compounds                         | QUANT SIG | RT     | EXP RT | REL RT  | RESPONSE               | CONCENTRATIONS       |                  |
|-----------------------------------|-----------|--------|--------|---------|------------------------|----------------------|------------------|
|                                   |           |        |        |         |                        | ON-COLUMN<br>(ug/mL) | FINAL<br>( ug/L) |
| =====                             | =====     | ==     | =====  | =====   | =====                  | =====                | =====            |
| 64 Fluoranthene                   | 202       |        |        |         | Compound Not Detected. |                      |                  |
| 65 Pyrene                         | 202       |        |        |         | Compound Not Detected. |                      |                  |
| \$ 66 Terphenyl-d14               | 244       | 18.352 | 18.356 | (0.916) | 625172                 | 20.0051              | 20.01            |
| 67 Butylbenzylphthalate           | 149       |        |        |         | Compound Not Detected. |                      |                  |
| 68 Benzo(a)anthracene             | 228       |        |        |         | Compound Not Detected. |                      |                  |
| * 69 Chrysene-d12                 | 240       | 20.040 | 20.050 | (1.000) | 1026859                | 20.0000              |                  |
| 70 3,3'-Dichlorobenzidine         | 252       |        |        |         | Compound Not Detected. |                      |                  |
| 71 Chrysene                       | 228       |        |        |         | Compound Not Detected. |                      |                  |
| 72 bis(2-Ethylhexyl)phthalate     | 149       |        |        |         | Compound Not Detected. |                      |                  |
| * 134 Di-n-octylphthalate-d4      | 153       | 21.141 | 21.150 | (1.000) | 1316049                | 20.0000              |                  |
| 73 Di-n-octylphthalate            | 149       |        |        |         | Compound Not Detected. |                      |                  |
| 74 Benzo(b)fluoranthene           | 252       |        |        |         | Compound Not Detected. |                      |                  |
| 75 Benzo(k)fluoranthene           | 252       |        |        |         | Compound Not Detected. |                      |                  |
| 76 Benzo(a)pyrene                 | 252       |        |        |         | Compound Not Detected. |                      |                  |
| * 77 Perylene-d12                 | 264       | 22.204 | 22.213 | (1.000) | 1110938                | 20.0000              |                  |
| 78 Indeno(1,2,3-cd)pyrene         | 276       |        |        |         | Compound Not Detected. |                      |                  |
| 79 Dibenzo(a,h)anthracene         | 278       |        |        |         | Compound Not Detected. |                      |                  |
| 80 Benzo(g,h,i)perylene           | 276       |        |        |         | Compound Not Detected. |                      |                  |
| 90 N-Nitrosodimethylamine         | 74        |        |        |         | Compound Not Detected. |                      |                  |
| 91 Aniline                        | 93        |        |        |         | Compound Not Detected. |                      |                  |
| 93 Benzidine                      | 184       |        |        |         | Compound Not Detected. |                      |                  |
| 103 Pyridine                      | 79        |        |        |         | Compound Not Detected. |                      |                  |
| 105 1-methylnaphthalene           | 141       |        |        |         | Compound Not Detected. |                      |                  |
| 111 Azobenzene (1,2-DP-Hydrazine) | 77        |        |        |         | Compound Not Detected. |                      |                  |
| 187 Total Benzofluoranthenes      | 252       |        |        |         | Compound Not Detected. |                      |                  |

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS  
 AREA AND RT SUMMARY

|                                                    |                               |
|----------------------------------------------------|-------------------------------|
| Instrument ID: nt6.i                               | Calibration Date: 18-FEB-2015 |
| Lab File ID: 02181506.d                            | Calibration Time: 09:17       |
| Lab Smp Id: ZW42A                                  | Client Smp ID: LMW-4-0215     |
| Analysis Type: SV                                  | Level: LOW                    |
| Quant Type: ISTD                                   | Sample Type: Water            |
| Operator: JZ                                       |                               |
| Method File: /chem2/nt6.i/20150218.b/SW846021115.m |                               |
| Misc Info: 15-2811                                 |                               |

Test Mode:  
 Use Initial Calibration Level 4.

| COMPOUND              | STANDARD | AREA LIMIT |         | SAMPLE  | %DIFF  |
|-----------------------|----------|------------|---------|---------|--------|
|                       |          | LOWER      | UPPER   |         |        |
| =====                 | =====    | =====      | =====   | =====   | =====  |
| 8 1,4-Dichlorobenze   | 274359   | 137180     | 548718  | 222362  | -18.95 |
| 27 Naphthalene-d8     | 982337   | 491168     | 1964674 | 783563  | -20.23 |
| 42 Acenaphthene-d10   | 661495   | 330748     | 1322990 | 537749  | -18.71 |
| 59 Phenanthrene-d10   | 1274103  | 637052     | 2548206 | 973135  | -23.62 |
| 69 Chrysene-d12       | 1469248  | 734624     | 2938496 | 1026859 | -30.11 |
| 134 Di-n-octylphthala | 1764428  | 882214     | 3528856 | 1316049 | -25.41 |
| 77 Perylene-d12       | 1586921  | 793460     | 3173842 | 1110938 | -29.99 |

| COMPOUND              | STANDARD | RT LIMIT |       | SAMPLE | %DIFF |
|-----------------------|----------|----------|-------|--------|-------|
|                       |          | LOWER    | UPPER |        |       |
| =====                 | =====    | =====    | ===== | =====  | ===== |
| 8 1,4-Dichlorobenze   | 8.43     | 7.93     | 8.93  | 8.43   | 0.01  |
| 27 Naphthalene-d8     | 10.48    | 9.98     | 10.98 | 10.47  | -0.04 |
| 42 Acenaphthene-d10   | 13.34    | 12.84    | 13.84 | 13.34  | -0.03 |
| 59 Phenanthrene-d10   | 15.72    | 15.22    | 16.22 | 15.72  | -0.03 |
| 69 Chrysene-d12       | 20.05    | 19.55    | 20.55 | 20.04  | -0.05 |
| 134 Di-n-octylphthala | 21.15    | 20.65    | 21.65 | 21.14  | -0.05 |
| 77 Perylene-d12       | 22.21    | 21.71    | 22.71 | 22.20  | -0.04 |

AREA UPPER LIMIT = +100% of internal standard area.  
 AREA LOWER LIMIT = - 50% of internal standard area.  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Analytical Resources, Inc.

RECOVERY REPORT

Client Name: Golder Associates  
Sample Matrix: LIQUID  
Lab Smp Id: ZW42A  
Level: LOW  
Data Type: MS DATA  
SpikeList File: LLLCS.spk  
Sublist File: LL.sub  
Method File: /chem2/nt6.i/20150218.b/SW846021115.m  
Misc Info: 15-2811

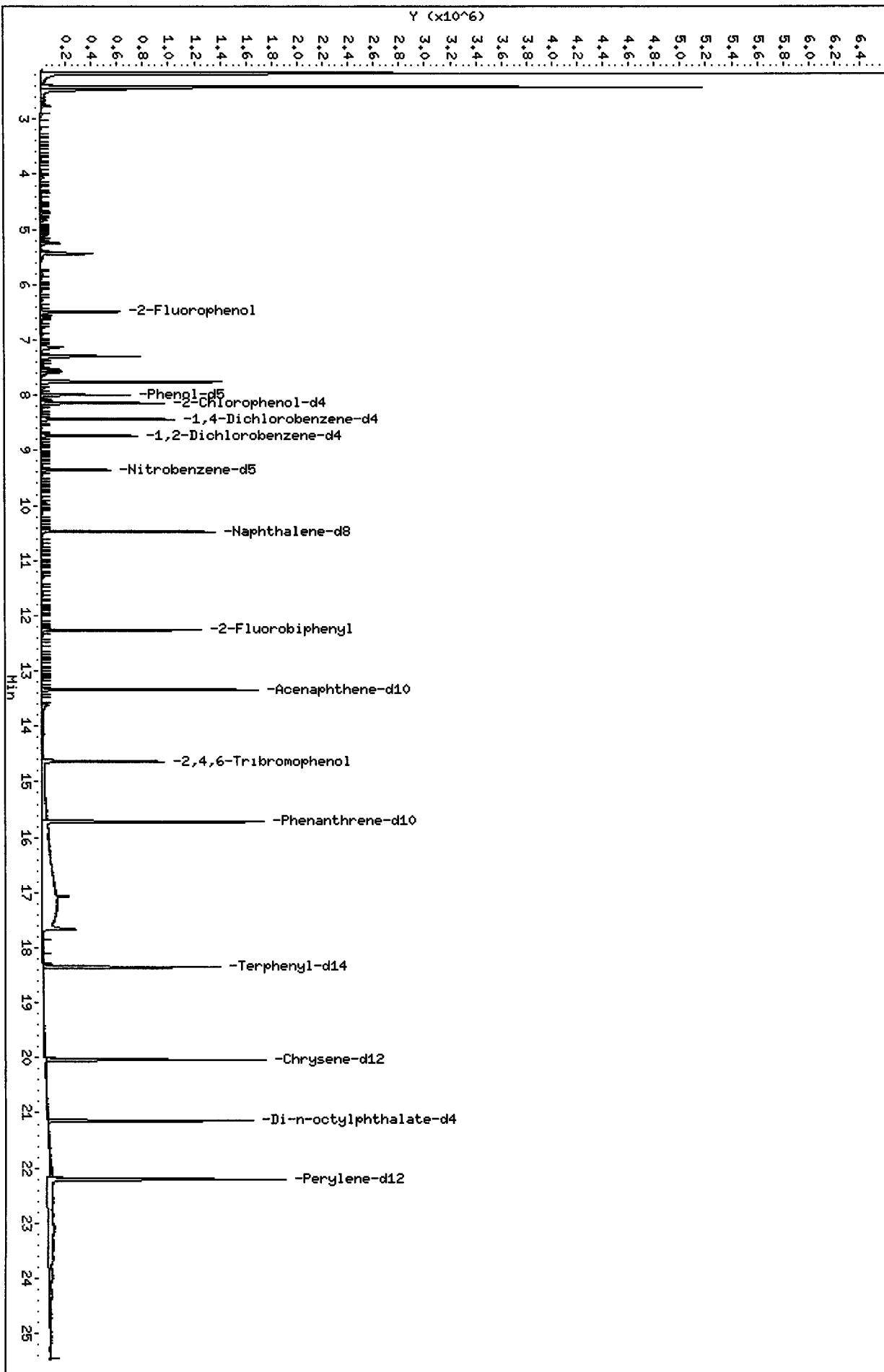
Client SDG: ZW42  
Fraction: SV  
Client Smp ID: LMW-4-0215  
Operator: JZ  
SampleType: SAMPLE  
Quant Type: ISTD

| SURROGATE COMPOUND       | CONC<br>ADDED<br>ug/L | CONC<br>RECOVERED<br>ug/L | %<br>RECOVERED | LIMITS |
|--------------------------|-----------------------|---------------------------|----------------|--------|
| \$ 1 2-Fluorophenol      | 37.50                 | 22.95                     | 61.19          | 23-100 |
| \$ 2 Phenol-d5           | 37.50                 | 22.80                     | 60.79          | 16-100 |
| \$ 5 2-Chlorophenol-d4   | 37.50                 | 24.52                     | 65.40          | 33-100 |
| \$ 10 1,2-Dichlorobenzen | 25.00                 | 14.88                     | 59.51          | 27-100 |
| \$ 18 Nitrobenzene-d5    | 25.00                 | 16.04                     | 64.17          | 34-101 |
| \$ 36 2-Fluorobiphenyl   | 25.00                 | 16.04                     | 64.18          | 38-100 |
| \$ 55 2,4,6-Tribromophen | 37.50                 | 26.15                     | 69.73          | 31-128 |
| \$ 66 Terphenyl-d14      | 25.00                 | 20.01                     | 80.02          | 27-122 |

Data File: /chem2/nt6.i/20150218.b/02181506.d  
Date : 18-FEB-2015 12:04  
Client ID: LHM-4-0215  
Sample Info: ZM42A  
Volume Injected (uL): 1.0  
Column phase: ZB-5msi

Instrument: nt6.i  
Operator: JZ  
Column diameter: 0.32

/chem2/nt6.i/20150218.b/02181506.d



CO-ELUTION SUMMARY FOR FILE - 02181506.d

Lab ID: ZW42A, Method: SW846021115.m, Instrument: nt6.i, Date: 18-FEB-2015

| RT | CO-ELUTION COMPOUNDS |
|----|----------------------|
|----|----------------------|

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NO CO-ELUTIONS

Quant Method: ICAL

Analytical Resources, Inc.

Semivolatiles Report SW846 Method 8270D

Data file : /chem2/nt6.i/20150218.b/02181502.d  
Lab Smp Id: ZV95MBW1 *ZW42-BMW1* Client Smp ID: ZV95MBW1  
Inj Date : 18-FEB-2015 09:51  
Operator : JZ Inst ID: nt6.i  
Smp Info : ZV95MBW1,  
Misc Info : 15-2549  
Comment : 1ul Injection *It-7811*  
Method : /chem2/nt6.i/20150218.b/SW846021115.m  
Meth Date : 18-Feb-2015 14:07 jianqing Quant Type: ISTD  
Cal Date : 11-FEB-2015 12:58 Cal File: 02111502.d  
Als bottle: 2 QC Sample: BLANK  
Dil Factor: 1.00000  
Integrator: HP RTE Compound Sublist: LL.sub  
Target Version: 3.50

Concentration Formula:  $\text{Amt} * \text{DF} * \text{Vt}/\text{Vo} * \text{CpndVariable}$  *02/18/15*

| Name | Value     | Description                     |
|------|-----------|---------------------------------|
| DF   | 1.00000   | Dilution Factor                 |
| Vt   | 500.00000 | Volume of final extract (uL)    |
| Vo   | 500.00000 | Volume of sample extracted (mL) |

Cpnd Variable

Local Compound Variable

| Compounds                       | QUANT | SIG   |       |        |         |                        |           | CONCENTRATIONS |  |
|---------------------------------|-------|-------|-------|--------|---------|------------------------|-----------|----------------|--|
|                                 |       |       | RT    | EXP RT | REL RT  | RESPONSE               | ON-COLUMN | FINAL          |  |
|                                 |       |       |       |        |         |                        | (ug/mL)   | ( ug/L)        |  |
| =====                           | ===== | ===== | ==    | =====  | =====   | =====                  | =====     | =====          |  |
| \$ 1 2-Fluorophenol             | 112   |       | 6.485 | 6.491  | (0.769) | 212405                 | 20.3091   | 20.31          |  |
| \$ 2 Phenol-d5                  | 99    |       | 7.986 | 7.998  | (0.947) | 265993                 | 21.0435   | 21.04          |  |
| 3 Phenol                        | 94    |       |       |        |         | Compound Not Detected. |           |                |  |
| \$ 5 2-Chlorophenol-d4          | 132   |       | 8.131 | 8.137  | (0.964) | 275180                 | 22.8738   | 22.87          |  |
| 4 Bis(2-Chloroethyl) ether      | 93    |       |       |        |         | Compound Not Detected. |           |                |  |
| 6 2-Chlorophenol                | 128   |       |       |        |         | Compound Not Detected. |           |                |  |
| 7 1,3-Dichlorobenzene           | 146   |       |       |        |         | Compound Not Detected. |           |                |  |
| * 8 1,4-Dichlorobenzene-d4      | 152   |       | 8.430 | 8.431  | (1.000) | 202867                 | 20.0000   |                |  |
| 9 1,4-Dichlorobenzene           | 146   |       |       |        |         | Compound Not Detected. |           |                |  |
| \$ 10 1,2-Dichlorobenzene-d4    | 152   |       | 8.729 | 8.735  | (1.035) | 133705                 | 13.7542   | 13.75          |  |
| 12 1,2-Dichlorobenzene          | 146   |       |       |        |         | Compound Not Detected. |           |                |  |
| 11 Benzyl alcohol               | 108   |       |       |        |         | Compound Not Detected. |           |                |  |
| 14 2,2'-oxybis(1-Chloropropane) | 45    |       |       |        |         | Compound Not Detected. |           |                |  |
| 13 2-Methylphenol               | 108   |       |       |        |         | Compound Not Detected. |           |                |  |
| 17 Hexachloroethane             | 117   |       |       |        |         | Compound Not Detected. |           |                |  |
| 16 N-Nitroso-di-n-propylamine   | 70    |       |       |        |         | Compound Not Detected. |           |                |  |

| Compounds                     | QUANT SIG | CONCENTRATIONS         |        |         |          |                      |                  |
|-------------------------------|-----------|------------------------|--------|---------|----------|----------------------|------------------|
|                               |           | RT                     | EXP RT | REL RT  | RESPONSE | ON-COLUMN<br>(ug/mL) | FINAL<br>( ug/L) |
| =====                         | =====     | ==                     | =====  | =====   | =====    | =====                | =====            |
| 15 4-Methylphenol             | 108       | Compound Not Detected. |        |         |          |                      |                  |
| \$ 18 Nitrobenzene-d5         | 82        | 9.354                  | 9.360  | (0.893) | 190365   | 14.9548              | 14.95            |
| 19 Nitrobenzene               | 77        | Compound Not Detected. |        |         |          |                      |                  |
| 20 Isophorone                 | 82        | Compound Not Detected. |        |         |          |                      |                  |
| 21 2-Nitrophenol              | 139       | Compound Not Detected. |        |         |          |                      |                  |
| 22 2,4-Dimethylphenol         | 107       | Compound Not Detected. |        |         |          |                      |                  |
| 23 Bis(2-Chloroethoxy)methane | 93        | Compound Not Detected. |        |         |          |                      |                  |
| 24 Benzoic acid               | 105       | Compound Not Detected. |        |         |          |                      |                  |
| 25 2,4-Dichlorophenol         | 162       | Compound Not Detected. |        |         |          |                      |                  |
| 26 1,2,4-Trichlorobenzene     | 180       | Compound Not Detected. |        |         |          |                      |                  |
| * 27 Naphthalene-d8           | 136       | 10.470                 | 10.477 | (1.000) | 714453   | 20.0000              |                  |
| 28 Naphthalene                | 128       | Compound Not Detected. |        |         |          |                      |                  |
| 29 4-Chloroaniline            | 127       | Compound Not Detected. |        |         |          |                      |                  |
| 30 Hexachlorobutadiene        | 225       | Compound Not Detected. |        |         |          |                      |                  |
| 31 4-Chloro-3-methylphenol    | 107       | Compound Not Detected. |        |         |          |                      |                  |
| 32 2-Methylnaphthalene        | 141       | Compound Not Detected. |        |         |          |                      |                  |
| 33 Hexachlorocyclopentadiene  | 237       | Compound Not Detected. |        |         |          |                      |                  |
| 34 2,4,6-Trichlorophenol      | 196       | Compound Not Detected. |        |         |          |                      |                  |
| 35 2,4,5-Trichlorophenol      | 196       | Compound Not Detected. |        |         |          |                      |                  |
| \$ 36 2-Fluorobiphenyl        | 172       | 12.260                 | 12.266 | (0.919) | 457045   | 14.7203              | 14.72            |
| 37 2-Chloronaphthalene        | 162       | Compound Not Detected. |        |         |          |                      |                  |
| 38 2-Nitroaniline             | 65        | Compound Not Detected. |        |         |          |                      |                  |
| 39 Dimethylphthalate          | 163       | Compound Not Detected. |        |         |          |                      |                  |
| 40 Acenaphthylene             | 152       | Compound Not Detected. |        |         |          |                      |                  |
| 41 2,6-Dinitrotoluene         | 165       | Compound Not Detected. |        |         |          |                      |                  |
| * 42 Acenaphthene-d10         | 164       | 13.334                 | 13.340 | (1.000) | 494033   | 20.0000              |                  |
| 43 3-Nitroaniline             | 138       | Compound Not Detected. |        |         |          |                      |                  |
| 44 Acenaphthene               | 153       | Compound Not Detected. |        |         |          |                      |                  |
| 45 2,4-Dinitrophenol          | 184       | Compound Not Detected. |        |         |          |                      |                  |
| 46 Dibenzofuran               | 168       | Compound Not Detected. |        |         |          |                      |                  |
| 47 4-Nitrophenol              | 109       | Compound Not Detected. |        |         |          |                      |                  |
| 48 2,4-Dinitrotoluene         | 165       | Compound Not Detected. |        |         |          |                      |                  |
| 50 Diethylphthalate           | 149       | Compound Not Detected. |        |         |          |                      |                  |
| 49 Fluorene                   | 166       | Compound Not Detected. |        |         |          |                      |                  |
| 51 4-Chlorophenyl-phenylether | 204       | Compound Not Detected. |        |         |          |                      |                  |
| 52 4-Nitroaniline             | 138       | Compound Not Detected. |        |         |          |                      |                  |
| 53 4,6-Dinitro-2-methylphenol | 198       | Compound Not Detected. |        |         |          |                      |                  |
| 54 N-Nitrosodiphenylamine     | 169       | Compound Not Detected. |        |         |          |                      |                  |
| \$ 55 2,4,6-Tribromophenol    | 330       | 14.627                 | 14.638 | (1.097) | 130284   | 22.7998              | 22.80            |
| 56 4-Bromophenyl-phenylether  | 248       | Compound Not Detected. |        |         |          |                      |                  |
| 57 Hexachlorobenzene          | 284       | Compound Not Detected. |        |         |          |                      |                  |
| 58 Pentachlorophenol          | 266       | Compound Not Detected. |        |         |          |                      |                  |
| * 59 Phenanthrene-d10         | 188       | 15.716                 | 15.723 | (1.000) | 904197   | 20.0000              |                  |
| 60 Phenanthrene               | 178       | Compound Not Detected. |        |         |          |                      |                  |
| 61 Anthracene                 | 178       | Compound Not Detected. |        |         |          |                      |                  |
| 62 Carbazole                  | 167       | Compound Not Detected. |        |         |          |                      |                  |
| 63 Di-n-butylphthalate        | 149       | Compound Not Detected. |        |         |          |                      |                  |

| Compounds                         | QUANT SIG |       |        |        |         |                        |  | CONCENTRATIONS       |                  |
|-----------------------------------|-----------|-------|--------|--------|---------|------------------------|--|----------------------|------------------|
|                                   |           | MASS  | RT     | EXP RT | REL RT  | RESPONSE               |  | ON-COLUMN<br>(ug/mL) | FINAL<br>( ug/L) |
| =====                             | =====     | ===== | ==     | =====  | =====   | =====                  |  | =====                | =====            |
| 64 Fluoranthene                   | 202       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 65 Pyrene                         | 202       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| \$ 66 Terphenyl-d14               | 244       |       | 18.355 | 18.356 | (0.916) | 657376                 |  | 21.9811              | 21.98            |
| 67 Butylbenzylphthalate           | 149       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 68 Benzo(a)anthracene             | 228       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| * 69 Chrysene-d12                 | 240       |       | 20.038 | 20.050 | (1.000) | 982691                 |  | 20.0000              |                  |
| 70 3,3'-Dichlorobenzidine         | 252       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 71 Chrysene                       | 228       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 72 bis(2-Ethylhexyl)phthalate     | 149       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| * 134 Di-n-octylphthalate-d4      | 153       |       | 21.144 | 21.150 | (1.000) | 1189121                |  | 20.0000              |                  |
| 73 Di-n-octylphthalate            | 149       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 74 Benzo(b)fluoranthene           | 252       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 75 Benzo(k)fluoranthene           | 252       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 76 Benzo(a)pyrene                 | 252       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| * 77 Perylene-d12                 | 264       |       | 22.207 | 22.213 | (1.000) | 1021670                |  | 20.0000              |                  |
| 78 Indeno(1,2,3-cd)pyrene         | 276       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 79 Dibenzo(a,h)anthracene         | 278       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 80 Benzo(g,h,i)perylene           | 276       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 90 N-Nitrosodimethylamine         | 74        |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 91 Aniline                        | 93        |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 93 Benzidine                      | 184       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 103 Pyridine                      | 79        |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 105 1-methylnaphthalene           | 141       |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 111 Azobenzene (1,2-DP-Hydrazine) | 77        |       |        |        |         | Compound Not Detected. |  |                      |                  |
| 187 Total Benzofluoranthenes      | 252       |       |        |        |         | Compound Not Detected. |  |                      |                  |

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS  
 AREA AND RT SUMMARY

|                                                    |                               |
|----------------------------------------------------|-------------------------------|
| Instrument ID: nt6.i                               | Calibration Date: 18-FEB-2015 |
| Lab File ID: 02181502.d                            | Calibration Time: 09:17       |
| Lab Smp Id: ZV95MBW1                               | Client Smp ID: ZV95MBW1       |
| Analysis Type: SV                                  | Level: LOW                    |
| Quant Type: ISTD                                   | Sample Type: Liquid           |
| Operator: JZ                                       |                               |
| Method File: /chem2/nt6.i/20150218.b/SW846021115.m |                               |
| Misc Info: 15-2549                                 |                               |

Test Mode:  
 Use Initial Calibration Level 4.

| COMPOUND              | STANDARD | AREA LIMIT |         | SAMPLE  | %DIFF  |
|-----------------------|----------|------------|---------|---------|--------|
|                       |          | LOWER      | UPPER   |         |        |
| =====                 | =====    | =====      | =====   | =====   | =====  |
| 8 1,4-Dichlorobenze   | 274359   | 137180     | 548718  | 202867  | -26.06 |
| 27 Naphthalene-d8     | 982337   | 491168     | 1964674 | 714453  | -27.27 |
| 42 Acenaphthene-d10   | 661495   | 330748     | 1322990 | 494033  | -25.32 |
| 59 Phenanthrene-d10   | 1274103  | 637052     | 2548206 | 904197  | -29.03 |
| 69 Chrysene-d12       | 1469248  | 734624     | 2938496 | 982691  | -33.12 |
| 134 Di-n-octylphthala | 1764428  | 882214     | 3528856 | 1189121 | -32.61 |
| 77 Perylene-d12       | 1586921  | 793460     | 3173842 | 1021670 | -35.62 |

| COMPOUND              | STANDARD | RT LIMIT |       | SAMPLE | %DIFF |
|-----------------------|----------|----------|-------|--------|-------|
|                       |          | LOWER    | UPPER |        |       |
| =====                 | =====    | =====    | ===== | =====  | ===== |
| 8 1,4-Dichlorobenze   | 8.43     | 7.93     | 8.93  | 8.43   | -0.01 |
| 27 Naphthalene-d8     | 10.48    | 9.98     | 10.98 | 10.47  | -0.06 |
| 42 Acenaphthene-d10   | 13.34    | 12.84    | 13.84 | 13.33  | -0.05 |
| 59 Phenanthrene-d10   | 15.72    | 15.22    | 16.22 | 15.72  | -0.04 |
| 69 Chrysene-d12       | 20.05    | 19.55    | 20.55 | 20.04  | -0.06 |
| 134 Di-n-octylphthala | 21.15    | 20.65    | 21.65 | 21.14  | -0.03 |
| 77 Perylene-d12       | 22.21    | 21.71    | 22.71 | 22.21  | -0.03 |

AREA UPPER LIMIT = +100% of internal standard area.  
 AREA LOWER LIMIT = - 50% of internal standard area.  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Analytical Resources, Inc.

RECOVERY REPORT

Client Name: CH2M Hill  
Sample Matrix: LIQUID  
Lab Smp Id: ZV95MBW1  
Level: LOW  
Data Type: MS DATA  
SpikeList File: LLLCS.spk  
Sublist File: LL.sub  
Method File: /chem2/nt6.i/20150218.b/SW846021115.m  
Misc Info: 15-2549

Client SDG: ZV95  
Fraction: SV  
Client Smp ID: ZV95MBW1  
Operator: JZ  
SampleType: BLANK  
Quant Type: ISTD

| SURROGATE COMPOUND       | CONC<br>ADDED<br>ug/L | CONC<br>RECOVERED<br>ug/L | %<br>RECOVERED | LIMITS |
|--------------------------|-----------------------|---------------------------|----------------|--------|
| \$ 1 2-Fluorophenol      | 37.50                 | 20.31                     | 54.16          | 23-100 |
| \$ 2 Phenol-d5           | 37.50                 | 21.04                     | 56.12          | 16-100 |
| \$ 5 2-Chlorophenol-d4   | 37.50                 | 22.87                     | 61.00          | 33-100 |
| \$ 10 1,2-Dichlorobenzen | 25.00                 | 13.75                     | 55.02          | 27-100 |
| \$ 18 Nitrobenzene-d5    | 25.00                 | 14.95                     | 59.82          | 34-101 |
| \$ 36 2-Fluorobiphenyl   | 25.00                 | 14.72                     | 58.88          | 38-100 |
| \$ 55 2,4,6-Tribromophen | 37.50                 | 22.80                     | 60.80          | 31-128 |
| \$ 66 Terphenyl-d14      | 25.00                 | 21.98                     | 87.92          | 27-122 |

Date : 18-FEB-2015 09:51

Client ID: ZV95MBW1

Instrument: nt6.i

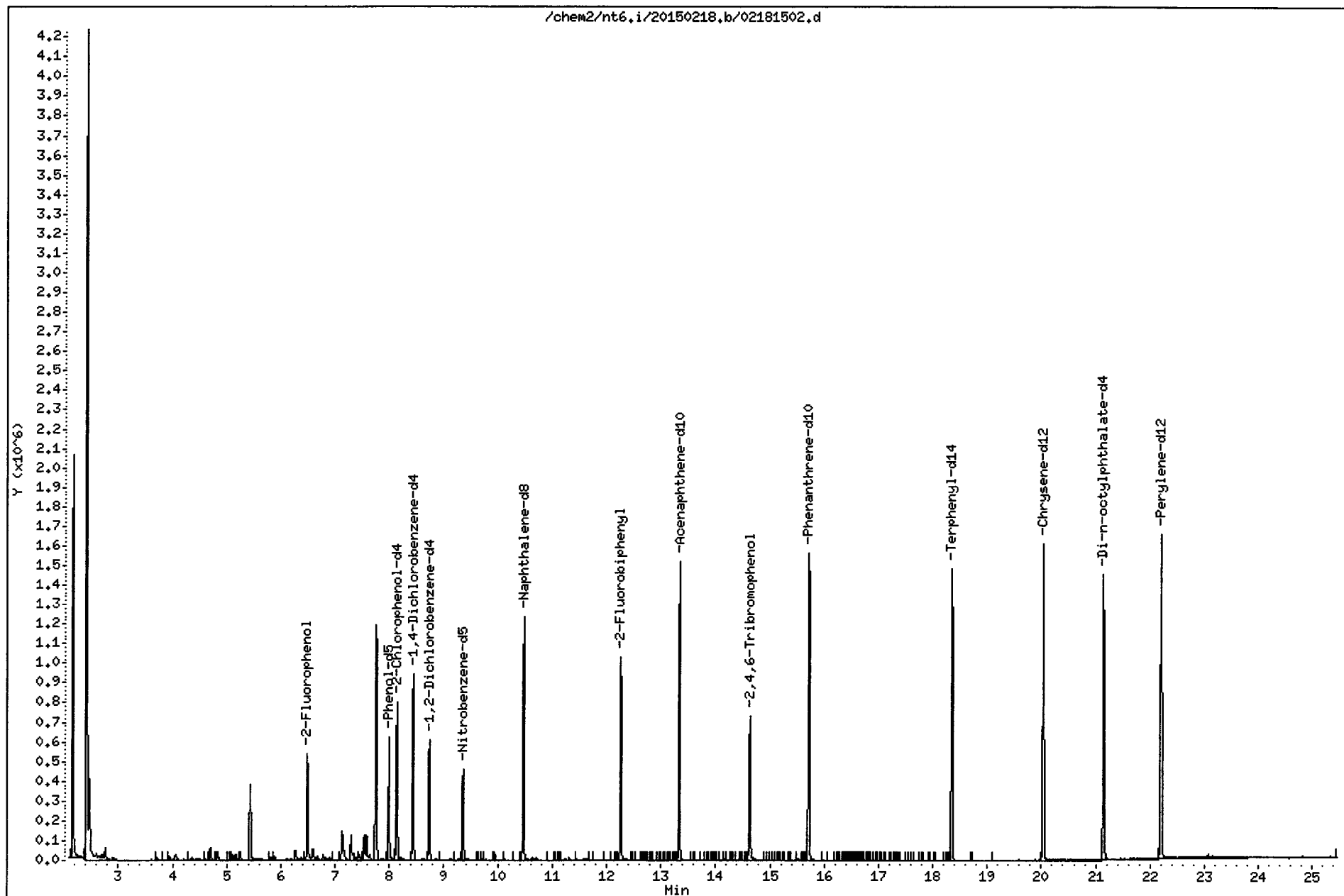
Sample Info: ZV95MBW1,

Volume Injected (uL): 1.0

Operator: JZ

Column phase: ZB-5msi

Column diameter: 0.32



ZW42:00021

CO-ELUTION SUMMARY FOR FILE - 02181502.d

Lab ID: ZV95MBW1, Method: SW846021115.m, Instrument: nt6.i, Date: 18-FEB-2015

| RT | CO-ELUTION COMPOUNDS |
|----|----------------------|
|----|----------------------|

-----

NO CO-ELUTIONS

Quant Method: ICAL

ZW42:00022

**APPENDIX B**  
**SAMPLE INTEGRITY DATA SHEETS (SIDS)**

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002  
 Site Location Ravensdale, WA Sample ID LMW-2-1113 1114  
 Sampling Location Groundwater Monitoring Well End of dedicated sampling tube *mw*

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Dedicated Pump Grundfos

Date 11/18/2014 Time 1140

Media Water Station LMW-2

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL - 7.62 ft below TOC (monument at elev. X) (bottom at 38.1 ft bgs, 4-in casing) @ 9 19 m 11/18/14 <sup>17</sup>

Screen Interval - 27.9-38.1 ft bgs Monument: 2.94 ags

Sand Pack Interval - 24.8-38.1 ft bgs (8-in hole) (~7.8 gal/sand pack vol)

Packer Depth - NA (~22.3 gal/casing vol) (~30.1 gal/total well vol)

Sample Description clear, sulfur odor

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                                     | Analysis         | Container             | Preservation / Amount |
|----------------------------------------------------|------------------|-----------------------|-----------------------|
| 3 - 40 mL                                          | VOA              | VOA Vial              | HCl                   |
| 1 - 500 ml                                         | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 - 500 ml                                         | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 - <sup>500ml</sup> <del>1</del> Liter, 2 - 40 ml | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 - 1 Liter, 2 - 500 ml                            | PCBs/Pest        | Glass Amber           | none                  |
| 2 - 500 mL                                         | SVOCs            | Glass Amber           | none                  |

(17 total bottles)

Sampler (signature) *[Signature]* Date 11/18/2-14

Supervisor (signature) *[Signature]* Date 11/21/2014

## FIELD PARAMETERS SHEET

Well ID LMW-2

Date 11/18/2014

Time Begin Purge 1022

Time Collect Sample 1140

(pH)

[illegible]

Comments:

Comments:

$$\frac{5 \text{ gal}}{4.3 \text{ min}} = 1.15 \text{ gpm} \rightarrow \frac{30 \text{ gal/well}}{1.15 \text{ gpm}} = 26 \text{ min/well} \times 3 = 78 \text{ min purge}$$

$PID = 0.2$   
 $PID = 0.2 \text{ ppm}$

Sulfur odor

Grundfos @ 111 Hz initially, then dialed up to 120 Hz.

match previous purge rate @ 18THz  
151 Hz D:led down @ 1030 b/c 181 Hz killed generator power,

Sampler's Initials JS

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002  
 Site Location Ravensdale, WA Sample ID LMW-3-1113 1114  
 Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Dedicated Pump Grundfos

Date 11/19/2014 Time 0920

Media Water Station LMW-3

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL - 13.65 ft below TOC (monument at elev. X) (bottom at 64.8 ft bgs, 4-in casing) 16:48 on 11/17/14

Screen Interval - 49.8-64.8 ft bgs Monument: 3.08 ags

Sand Pack Interval - 47.1-64.8 ft bgs (8-in hole) (~10.4 gal/sand pack)

Packer Depth - 39.33 ft bgs (~36.1 gal/casing vol) (~16.6 gal/packer casing volume)

(~27.0 gal/total well vol below packer)

Sample Description clear, no odor

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                           | Analysis         | Container             | Preservation / Amount |
|------------------------------------------|------------------|-----------------------|-----------------------|
| 3 - 40 mL                                | VOA              | VOA Vial              | HCl                   |
| 1 - 500 ml                               | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 - 500 ml                               | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 - <sup>500 mL</sup> 1 Liter, 2 - 40 ml | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 - 1 Liter, 2 - 500 ml                  | PCBs/Pest        | Glass Amber           | none                  |
| 2 - 500 mL                               | SVOCs            | Glass Amber           | none                  |

17  
Bottles  
Total

Sampler (signature) [Signature] Date 11/19/2014

Supervisor (signature) [Signature] Date 11/21/2014

## FIELD PARAMETERS SHEET

Well ID LMW-3  
Date 11/19/2014  
Time Begin Purge 0815  
Time Collect Sample 0920

(p14)

[illegible]

Comments:

packer 110 psi  
Grundfos 185 Hz

$P_{10} = 0.0 \text{ ppm}$

$$\frac{5 \text{ gal}}{4 \text{ min}} = 1.25 \text{ gpm} \rightarrow \frac{\frac{27 \text{ gal}}{\text{well vol}}}{1.25 \text{ gpm}} = 22 \frac{\text{min}}{\text{well vol}} = 65 \text{ min purge}$$

Sampler's Initials JS

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002

Site Location Ravensdale, WA Sample ID LMW-4-1113 1114

Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Dedicated Pump Grundfos

Date 11/18/2014 Time 1310

Media Water Station LMW-4

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL - 9.67 ft below TOC (monument at elev. X) (bottom at 209.7 ft bgs, 4-in casing) @ 9:22 on 11/17/2014  
(not corrected for angle)

Screen Interval - 195-209.7 ft bgs Monument: 2.76 ags

Sand Pack Interval - 189-209.7 ft bgs (8-in hole) (~12.3 gal/sand pack)

Packer Depth - 187.3 ft bgs (~133.3 gal/casing vol) (~14.6 gal/packer casing volume)

(~26.9 gal/total well vol below packer)

\*\* Depths corrected for 70° inclination

Sample Description clear, no odor sulfur odor,

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount          | Analysis         | Container             | Preservation / Amount |
|-------------------------|------------------|-----------------------|-----------------------|
| 3 - 40 mL               | VOA              | VOA Vial              | HCl                   |
| 1 - 500 ml              | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 - 500 ml              | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 - 1 Liter, 2 - 40 ml  | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 - 1 Liter, 2 - 500 ml | PCBs/Pest        | Glass Amber           | none                  |
| 2 - 500 mL              | SVOCs            | Glass Amber           | none                  |

Sampler (signature) [Signature] Date 11/18/2014

Supervisor (signature) [Signature] Date 11/21/2014

17  
Bottles  
Total

# FIELD PARAMETERS SHEET

Well ID LMW-4  
Date 11/18/2014  
Time Begin Purge 1215  
Time Collect Sample 1310

(pH)

[illegible]

Comments:

Grundfvs: 191 Hz

Packer: 140 psi  $\rightarrow$  ~~340-149~~

$$\frac{5 \text{ gal}}{6 \text{ min}} = \frac{0.83}{1.5} \text{ gpm}$$

$$\frac{27 \text{ gpl/welt vsl}}{0.05 \text{ gram}} = \frac{18}{32} \text{ min/vsl under pressure}$$

$x3 = \cancel{98}$  min purge  
54

PID = 0.0 ppm

• میں نے

Sampler's Initials JS

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002  
 Site Location Ravensdale, WA Sample ID LMW-5-11184  
 Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Dedicated Pump Grundfos

Date 11/19/2014 Time 1130

Media Water Station LMW-5

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL -15.08 ft below TOC (monument at elev. X) (bottom at 241.8 ft bgs, 4-in casing) 11/17/14

Screen Interval -231.8-241.8 ft bgs Monument: 3.24 ags

Sand Pack Interval -231.8-241.8 ft bgs (8-in hole) (~5.9 gal/sand pack)

Packer Depth -222.11 ft bgs (~150.8 gal/casing vol) (~12.9 gal/packer casing volume)  
 (~18.7 gal/total well vol below packer)

Sample Description clear, sulfur odor

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                           | Analysis         | Container             | Preservation / Amount |
|------------------------------------------|------------------|-----------------------|-----------------------|
| 3 – 40 mL                                | VOA              | VOA Vial              | HCl                   |
| 1 – 500 ml                               | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 – 500 ml                               | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 – <sup>500 mL</sup> 1 Liter, 2 – 40 ml | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 – 1 Liter, 2 – 500 ml                  | PCBs/Pest        | Glass Amber           | none                  |
| 2 – 500 mL                               | SVOCs            | Glass Amber           | none                  |

17  
Bottles  
Total

Sampler (signature) [Signature] Date 11/19/2014

Supervisor (signature) [Signature] Date 11/21/2014

# FIELD PARAMETERS SHEET

Well ID LMW-5  
 Date 11/19/2014  
 Time Begin Purge 1043  
 Time Collect Sample 1130

| Water Level<br>feet bmp | Time                 | Volume<br>Purged | pH   | Conductivity<br>uS/cm | Temp.<br>°C | DO<br>mg/L | Turbidity<br>NTU | Eh<br>Rel mV |
|-------------------------|----------------------|------------------|------|-----------------------|-------------|------------|------------------|--------------|
|                         | 1055                 |                  | 7.05 | 696                   | 11.1        | 0.02       | 1.14             | -83.9        |
|                         | <del>10</del> 1105   |                  | 7.04 | 696                   | 11.1        | 0.01       | 0.59             | -88.8        |
|                         | 1115                 |                  | 7.04 | 692                   | 11.1        | 0.00       | 0.88             | -92.5        |
|                         | <del>1125</del> 1120 |                  | 7.03 | 688                   | 11.1        | 0.00       | 0.94             | -93.6        |
|                         | <del>1125</del> 1125 |                  | 7.03 | 686                   | 11.2        | 0.00       | 0.92             | -94.9        |
|                         | <del>1125</del> 1130 |                  | 7.03 | 691                   | 11.2        | 0.00       | 0.82             | -95.5        |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |
|                         |                      |                  |      |                       |             |            |                  |              |

## Comments:

Grund fvs: 201 Hz

Packer: 130 psi

$$\frac{5 \text{ gal}}{4 \text{ min}} = 1.25 \text{ gpm} \rightarrow$$

$$\frac{19 \text{ gal/well vol}}{1.25 \text{ gpm}} = 15.2 \text{ min/well vol}$$

$$\times 3 = 46 \text{ min purge}$$

PID = 0.1 ppm

Sampler's Initials jsl / AMR

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002  
 Site Location Ravensdale, WA Sample ID LMW-6-11184  
 Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Dedicated Pump Grundfos

Date 11/20/2014 Time 1240

Media Water Station LMW-6

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL -28.80 ft below TOC (monument at elev. X) (bottom at 105.9 ft bgs, 4-in casing) @ 1016 cm 11/17/14

Screen Interval – 90.9-105.9 ft bgs Monument: 3.05 ags

Sand Pack Interval – 82.5-105.9 ft bgs (8-in hole) (~13.7 gal/sand pack)

Packer Depth – 81.22 ft bgs (~53 gal/casing vol) (~16.1 gal/packer casing volume)

(~29.9 gal/total well vol below packer)

Sample Description clear, no odor

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                                     | Analysis         | Container             | Preservation / Amount |
|----------------------------------------------------|------------------|-----------------------|-----------------------|
| 3 – 40 mL                                          | VOA              | VOA Vial              | HCl                   |
| 1 – 500 ml                                         | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 – 500 ml                                         | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 – <sup>500 mL</sup> <del>Liter</del> , 2 – 40 ml | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 – 1 Liter, 2 – 500 ml                            | PCBs/Pest        | Glass Amber           | none                  |
| 2 – 500 mL                                         | SVOCs            | Glass Amber           | none                  |

Sampler (signature) [Signature] Date 11/20/2014

Supervisor (signature) [Signature] Date 11/21/2014

## FIELD PARAMETERS SHEET

Well ID 1 MW-6

Date 11/20/2014

Time Begin Purge 1141

Time Collect Sample 1240

[illegible]

Comments:

Grundfos: 299 Hz

Bladder = 110 psi

$$\frac{5 \text{ gal}}{53 \text{ min}} = \frac{1 \text{ gpm}}{1.67} \rightarrow \frac{30 \text{ gal/well vol}}{1.67 \text{ gpm}} = 18 \text{ min/well vol} \times 3 = 54 \text{ min purge}$$

PID = 20 ppm

Sampler's Initials JS/AMR

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002  
 Site Location Ravensdale, WA Sample ID LMW-7-1113, LMW-7-1113-D  
 Sampling Location Groundwater Monitoring Well End of dedicated sampling tube *No Sample*

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Dedicated Pump Grundfos

Date 11/20/14

Time NA - No Sample

Media Water

Station LMW-7

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL 226.30 *(not corrected)* ft below TOC (monument at elev. X) (bottom at 253.7 ft bgs, 4-in casing) *Q 934 on 11/17/14*

Screen Interval - 239.6-253.7 ft bgs Monument: 3.09 ags

Sand Pack Interval - NA

Packer Depth - NA (~28.3 gal/casing vol) \*\* Depths corrected for 70° inclination

Sample Description NA - No Sample

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                                     | Analysis         | Container             | Preservation / Amount |
|----------------------------------------------------|------------------|-----------------------|-----------------------|
| 6 - 40 mL                                          | VOA              | VOA Vial              | HCl                   |
| 2 - 500 ml                                         | Total Metals     | HDPE                  | HNO3 (non)            |
| 2 - 500 ml                                         | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 8 - <del>1</del> <sup>500mL</sup> Liter, 4 - 40 ml | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 4 - 1 Liter, 4 - 500 ml                            | PCBs/Pest        | Glass Amber           | none                  |
| 4 - 500 mL                                         | SVOCs            | Glass Amber           | none                  |

Sampler (signature) [Signature] Date 11/20/14

Supervisor (signature) [Signature] Date 11/21/2014

## FIELD PARAMETERS SHEET

Well ID LMW-7

Date 1/20/14

Time Begin Purge NA

Time Collect Sample NA[illegible]

## Comments:

Comments:  
Grundfos @ 400Hz. No water. Performed troubleshooting: Pump works but is unable to bring water to the surface. Pulled pump. Taking to Geo tech for further troubleshooting.

Sampler's Initials JSF

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002  
 Site Location Ravensdale, WA Sample ID LMW-8-11124  
 Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Dedicated Tubing and Peristaltic Pump, Bailer for VOC samples

Date 11/19/2014 Time 1600 1005

Media Water Station LMW-8

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL - ~~5.68~~ <sup>4.70' js1</sup> ft below TOC (PVC at black notch) (bottom at 13 ft bgs, 2-in casing) C 1045 on 11/17/14

Screen Interval - 8-13 ft bgs PVC stickup: 1.72 ags

Sand Pack Interval - 6-13 ft bgs (8-in hole) (~5.1 gal/sand pack)

Packer Depth - NA (~1.9 gal/casing vol) (~7.0 gal/total well vol)

Sample Description Clear, no odor

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                                      | Analysis         | Container             | Preservation / Amount |
|-----------------------------------------------------|------------------|-----------------------|-----------------------|
| 3 - 40 mL                                           | VOA              | VOA Vial              | HCl                   |
| 1 - 500 ml                                          | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 - 500 ml                                          | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 - <del>1</del> <sup>500 mL</sup> Liter, 2 - 40 ml | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 - 1 Liter, 2 - 500 ml                             | PCBs/Pest        | Glass Amber           | none                  |
| 2 - 500 mL                                          | SVOCs            | Glass Amber           | none                  |

Sampler (signature) [Signature] Date 11/19/2014

Supervisor (signature) [Signature] Date 11/21/2014

Time Collect Sample 1005

[illegible]

Comments:

PID = 0.0 ppm

Flowrate:  $\sim 280 \text{ mL/min}$

WCS + ACID WALS collected using bucket

0900 collected EB - Field blank before purge. Through tubing (and filter for diss metals) using lab-provided DI water: EB-1114

Sampler's Initials DWL/JSL

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002

Site Location Ravensdale, WA Sample ID LMW-9-1113

Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Pump Grundfos and Dedicated Tubing

Date 11/19/2014 Time 1315

Media Water Station LMW-9

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL - 100.18 ft below TOC (PVC at black notch) (bottom at 159 ft bgs, 2-in casing) 10.54 on 11/17/14

Screen Interval - 149-159 ft bgs PVC stickup: 2.86 ags

Sand Pack Interval - 143.5-159 ft bgs (8-in hole) (~11.4 gal/sand pack)

Packer Depth - NA (~10.2 gal/casing vol) (~21.6 gal/total well vol)

Sample Description clear, no odor

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount          | Analysis         | Container             | Preservation / Amount |
|-------------------------|------------------|-----------------------|-----------------------|
| 3 - 40 mL               | VOA              | VOA Vial              | HCl                   |
| 1 - 500 ml              | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 - 500 ml              | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 - 1 Liter, 2 - 40 ml  | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 - 1 Liter, 2 - 500 ml | PCBs/Pest        | Glass Amber           | none                  |
| 2 - 500 mL              | SVOCs            | Glass Amber           | none                  |

17  
Bottles  
total!

Sampler (signature) [Signature] Date 11/19/2014

Supervisor (signature) [Signature] Date 11/21/2014

(plt)

Comments:

PID: 0.0 ppm

Grundfos: 359 Hz

$$\frac{5 \text{ gal}}{4 \text{ minutes}} = 1.25 \text{ gpm} \rightarrow \frac{21.6 \text{ gal/well volume}}{1.25 \text{ gpm}} = 17.28 \text{ minutes/well volume}$$

∴ ~17 minute purge

## Golder Associates

## SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002

Site Location Ravensdale, WA Sample ID LMW-10-1113-1114

Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler QED Bladder

Date 11/18/2014 Time 0945

Media Water Station LMW-10

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL - 1.67 ft below TOC (PVC) (bottom at 289 ft bgs, 4-in casing) 01:09 on 11/17/2014

Screen Interval - 267-289 ft bgs PVC stickup: 3.12 ags

Sand Pack Interval - 258-289 ft bgs (9-in hole) (~18.2 gal/sand pack)

Packer Depth - NA (~191 gal/casing vol) (~209 gal/total well vol)

Sample Description clear, no odor

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                           | Analysis         | Container             | Preservation / Amount |
|------------------------------------------|------------------|-----------------------|-----------------------|
| 3 - 40 mL                                | VOA              | VOA Vial              | HCl                   |
| 1 - 500 ml                               | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 - 500 ml                               | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 - 1 <sup>500 mL</sup> Liter, 2 - 40 ml | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 - 1 Liter, 2 - 500 ml                  | PCBs/Pest        | Glass Amber           | none                  |
| 2 - 500 mL                               | SVOCs            | Glass Amber           | none                  |

Sampler (signature) [Signature] Date 11/18/2014

Supervisor (signature) [Signature] Date 11/21/2014

## FIELD PARAMETERS SHEET

Well ID LMW-10  
Date 11/19/2014  
Time Begin Purge 0900  
Time Collect Sample 0945

(PH)

[illegible]

Comments:

Tank: 110psi

Controller: 60psi

Cycle ID: 50 (20/20), 2 cpm

PID = 0.0 ppm

Purge rate:  $\sim 900 \text{ mLs/min}$

Sampler's Initials jsl

# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002  
 Site Location Ravensdale, WA Sample ID LMW-11-11184  
 Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Pump Grundfos and QED Bladder

Date 11/20/2014 Time 1045

Media Water Station LMW-11

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL - 158.02 ft below TOC (PVC) (bottom at 707 ft bgs, 4-in casing) At 1026 @ 11/17/2014

Screen Interval - 696-707 ft bgs PVC stickup: 2.70 ags

Sand Pack Interval - 688-707 ft bgs (8-in hole) (~11.2 gal/sand pack)

Packer Depth - NA (~360.4 gal/casing vol) (~371.6 gal/total well vol)

Sample Description clear, no odor Sulfur odor

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                         | Analysis         | Container             | Preservation / Amount |
|----------------------------------------|------------------|-----------------------|-----------------------|
| 3 - 40 mL                              | VOA              | VOA Vial              | HCl                   |
| 1 - 500 ml                             | Total Metals     | HDPE                  | HNO3 (non)            |
| 1 - 500 ml                             | Dissolved Metals | HDPE                  | HNO3 (filter)         |
| 4 - <sup>500 mL</sup> Liter, 2 - 40 ml | TPH-HCID         | Glass Amber, VOA Vial | HCl                   |
| 2 - 1 Liter, 2 - 500 ml                | PCBs/Pest        | Glass Amber           | none                  |
| 2 - 500 mL                             | SVOCs            | Glass Amber           | none                  |

Sampler (signature) [Signature] Date 11/20/2014

Supervisor (signature) [Signature] Date 11/21/2014

## FIELD PARAMETERS SHEET

Well ID MW-11

Date 11/20/2014

Date 11/20/2014  
Time Begin Purge 0905 (Grindfish), Star 1005 (bladder)

Time Collect Sample 1045

[illegible]

Comments:

Comments: 0905 - start pump (Grundfos) @ 170' hloc. 390 Hz. Purge rate:  $\frac{59 \text{ gal}}{10 \text{ min}} = 0.5 \text{ gpm}$

1005 - Start bladder pump after 1 hr: 230 gal  
purged

$PID = 0.0 \text{ ppm}$

Tank 11 ops<sup>i</sup>

Controller 110 psi

ID: ~~12~~<sup>20</sup>cpm (#30, 30s/30s)

Rate: 600 mLs/min  
sp 450

Sampler's Initials Jsl / AM12



# SAMPLE INTEGRITY DATA SHEET

Plant/Site Landsburg Mine Site Project No. 923-1000-002

Site Location Ravensdale, WA Sample ID LMW-4-0215

Sampling Location Groundwater Monitoring Well End of dedicated sampling tube

Technical Procedure Reference(s) TP-1.4-6A, TP-1.2-20, TP-1.2-23

Type of Sampler Dedicated Pump Grundfos

Date 2/12/2015 Time 1450

Media Water Station LMW-4

Sample Type: grab time composite space composite

Sample Acquisition Measurements (depth, volume of static well water and purged water, etc.)

SWL - 7.65 ft below TOC (monument at elev. X) (bottom at 209.7 ft bgs, 4-in casing) (not converted) (PVC) 2/15/2015

Screen Interval - 195-209.7 ft bgs Monument: 2.76 ags

Sand Pack Interval - 189-209.7 ft bgs (8-in hole) (~12.3 gal/sand pack)

Packer Depth - 187.3 ft bgs (~133.3 gal/casing vol) (~14.6 gal/packer casing volume)

(~26.9 gal/total well vol below packer)

\*\* Depths corrected for 70° inclination

Sample Description clear, no odor - slight sulfur

Field Measurements on Sample (pH, conductivity, etc.)

SEE FIELD PARAMETERS SHEET

| Aliquot Amount                   | Analysis | Container   | Preservation / Amount |
|----------------------------------|----------|-------------|-----------------------|
| 3 <u>2</u> - 1 Liter - to Onsite | SVOCs    | Glass Amber | none                  |
| 4 <u>2</u> - 500 mL - to ARI     | SVOCs    | Glass Amber | none                  |

BEHP only

Sampler (signature) [Signature] Date 2/12/2015

Supervisor (signature) [Signature] Date 2/18/2015

# FIELD PARAMETERS SHEET

Well ID LMW-4  
 Date 2/12/2015  
 Time Begin Purge 1412  
 Time Collect Sample 1450

(pH)

| Water Level<br>feet bmp | Time            | Volume<br>Purged | pH   | Conductivity<br>uS/cm | Temp.<br>°C | DO<br>mg/L | Turbidity<br>NTU | Eh<br>Rel mV |
|-------------------------|-----------------|------------------|------|-----------------------|-------------|------------|------------------|--------------|
|                         | 1420            |                  | 7.03 | 409                   | 10.7        | 0.01       | 1.11             | -42.9        |
|                         | 1425            |                  | 7.02 | 405                   | 10.7        | 0.00       | 0.58             | -57.3        |
|                         | 1430            |                  | 7.01 | 410                   | 10.7        | 0.00       | 0.67             | -67.4        |
|                         | 1435            |                  | 7.01 | 408                   | 10.7        | 0.00       | 0.41             | -71.1        |
|                         | 1440            |                  | 7.01 | 405                   | 10.7        | 0.00       | 0.24             | -74.9        |
|                         | 1445            |                  | 7.00 | 410                   | 10.7        | 0.00       | 0.39             | -78.0        |
|                         | <del>1450</del> |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |
|                         |                 |                  |      |                       |             |            |                  |              |

## Comments:

Ground fvs: 197 wtz

Packer: 140 psi

$$\frac{5 \text{ gal}}{2 \text{ min}} = 2.5 \text{ gpm}$$

$$\frac{27 \text{ gal/well vol}}{2.5 \text{ gpm}} = 10.8 \text{ min/vol under packer}$$

$$\times 3 = 33 \text{ min purge}$$

Sampler's Initials JSL

**APPENDIX C**

**LANDSBURG MINE SITE NOVEMBER 2014 DATA VALIDATION  
AND QUALITY ASSURANCE / QUALITY CONTROL REVIEW MEMORANDUM**

**TECHNICAL REVIEW OF LABORATORY RAW DATA DELIVERABLE  
FOR SEMI-VOLATILE ANALYSIS**

## TECHNICAL MEMORANDUM

**Date:** February 24, 2015

**To:** Bill Kombol

**From:** Jill Lamberts, Project Environmental Scientist

**Email:** Jill\_Lamberts@golder.com

**Project No.:** 923-1000-002.R273

**Company:** Palmer Coking Coal Company



**RE: LANDSBURG MINE SITE NOVEMBER 2014 DATA VALIDATION & QUALITY ASSURANCE /  
QUALITY CONTROL REVIEW**

A total of 13 water samples (including two Trip Blanks and one Equipment Blank) were collected by Golder Associates Inc. (Golder) on November 18 to 20, 2014 and February 12, 2015 as part of the Landsburg sampling project. Samples were analyzed by Analytical Resources Inc. of Tukwila, Washington for the following:

- Volatile Organic Compounds (VOCs) by United States Environmental Protection Agency (EPA) Method 8260C;
- Semi-Volatile Organic Compounds (SVOCs) by GC/MS by EPA 8270D;
- Pesticides by EPA 8081B;
- Polychlorinated Biphenyls (PCBs) by EPA 8082A;
- Northwest Total Petroleum Hydrocarbon Identification Scan (NWTPH-HCID) by NWTPH-HCID Method; and,
- Total Metals EPA Method 6010C and 200.8; and Mercury by EPA 7470A.

Samples were analyzed in accordance with procedures described in Test Methods for Evaluating Solid Waste, Physical/Chemical Methods (EPA SW-846, 3rd edition; methods 6010C, 7470A, 8081B, 8082A, 8260C, 200.8 and 8270D), and for Northwest Total Petroleum Hydrocarbon Methods. Quality assurance / quality control (QA/QC) reviews of laboratory data were performed in the laboratory in accordance with the laboratory quality assurance program plan. The data validation QA/QC review focused primarily on laboratory result summary sheets and quality control summary sheets to ensure that work plan data quality objectives were met for the project. Data validation was conducted in accordance with the criteria outlined in the National Function Guidelines for Inorganic Review (EPA 2014a) and National Functional Guidelines for Organic Review (EPA 2014b), modified to include method specific requirements of the laboratory analytical methods and laboratory standard operating procedures (SOPs).

The validation level for the data is Tier II, and included the following:

- Data Package Completeness
- Verification of required deliverables
- Evaluation of holding times



- Laboratory narrative evaluation
- Evaluation and qualification of quality control elements for: Surrogates, Matrix Spike, Laboratory Control samples, Laboratory Duplicates, Method Blanks, and Field Blank and Field Duplicate evaluation as applicable
- Evaluation of detection limits

Raw data was not provided and calibration elements, including Gas Chromatograph (GC) instrument tuning and performance check, initial and continuing calibration, internal standard performance, and compound identification, were not evaluated unless information was provided by the lab in the case narratives. Data review and validation was performed by an experienced quality assurance chemist independent of the analytical laboratory and not directly involved in the project. Data qualifiers that were applied by the laboratory have been removed from the data summary report sheets, when applicable, and superseded by data validation qualifiers. Overall, the data review showed that data are acceptable for use except where indicated by data qualifiers. For details about the data validation, refer to the discussion below and see the validated analytical report pages in Attachment 1. Table 1 is a summary of the qualifiers applied to the data.

#### Data Qualifier Definitions

|     |                                                                                                                                                                                                                                                                                                          |
|-----|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| U   | The constituent was analyzed for, but was not detected above the reported sample quantitation limit.                                                                                                                                                                                                     |
| J   | The constituent was positively identified and detected; however, the concentration reported is an estimated value because the result is less than the quantitation limit or quality control criteria were not met.                                                                                       |
| J+  | The constituent was positively identified and detected; however, the concentration reported is an estimated value because the result may be biased high.                                                                                                                                                 |
| J-  | The constituent was positively identified and detected; however, the concentration reported is an estimated value because the result may be biased low.                                                                                                                                                  |
| UJ  | The constituent was not detected; the associated quantitation limit is an estimated value because quality control criteria were not met.                                                                                                                                                                 |
| R   | Data are rejected due to significant exceedance of quality control criteria. The analyte may or may not be present. Additional sampling and analysis may be required to determine the presence or absence of the constituent. For statistical reasons, rejected values are not included in the database. |
| UR  | The constituent is rejected at the reported quantitation limit.                                                                                                                                                                                                                                          |
| DNR | Do Not Report. More than one set of results are reported due to re-analyses or re-reporting (below reporting level). This result should not be reported.                                                                                                                                                 |

### Data Validation Comments

- Several samples had small- or pea-sized bubbles in the vials to be analyzed for VOCs (LMW-2-1114, LMW-4-1114, LMW-10-1114, Trip Blank-111914, Trip Blank-112014). Small- and pea-sized bubbles (< 4 millimeters [mm]) are not considered to affect VOC results, especially if analyzed within seven days. Protocol dictates that the lab will choose vials without bubbles or headspace for analysis.
- Sample LMW-5-1114 had large bubbles in the vials to be analyzed for VOCs. Large bubbles (4 to < 6 mm in size) may have a minor effect on VOC results, but protocol dictates that the lab will select vials without bubbles for analysis. Sample LMW-5-1114 had no detections of VOCs, which is consistent with historical results. No action was taken other than to note.
- For sample data group (SDG) ZL65/ZL66: The method blank for VOCs contained methylene chloride and hexachlorobutadiene. The associated Trip Blank-112014 contained methylene chloride greater than two times the blank contamination amount and is qualified as estimated with a high bias (J+). There were no detects of methylene chloride in the remaining associated samples. There were no detects of hexachlorobutadiene in the associated samples.
- SDG ZL65/ZL66: The continuing calibration standard (CCAL) was out of control low for VOCs analytes vinyl acetate, 2-chlorovinylether, and naphthalene. The associated sample results were qualified as UJ (The analyte was not detected above the reporting limit and is estimated).
- SDG ZL65/ZL66: The CCAL was out of control high for SVOCs analyte 4-nitroaniline. No action was taken since the associated sample results were non-detect for the analytes.
- SDG ZL65/ZL66: The laboratory control sample duplicate (LCSD) was out of control low for 3,3'-dichlorobenzidine and the relative percent difference (RPD) for 4-chloroaniline was out of control limits. No action was taken since the associated sample results were non-detect for the analytes.
- SDG ZL38/ZL39: The method blank for VOCs contained methylene chloride, naphthalene, n-butylbenzene, 1,2,3-trichlorobenzene, and hexachlorobutadiene. Sample LMW-10-1114 contained hexachlorobutadiene at 0.33 micrograms per liter (µg/L) and is qualified as U (non-detect) since the result is less than the method blank contamination value. EB-1114 contained methylene chloride at 1.4 µg/L and is qualified as U (non-detect) since the result is less than two times the method blank contamination (allowable for methylene chloride since it is a common lab contaminant). The Trip Blank-111814 contained methylene chloride greater than two times the blank contamination amount and is qualified as estimated with a high bias (J+). No other samples were affected by the method blank contamination.
- SDG ZL38/ZL39: The surrogate 2,4,6-tribromophenol (TBP) for SVOCs was out of control low for sample LMW-2-1114. No action was taken since the surrogate was within advisory control limits.
- SDG ZL38/ZL39: The Trip Blank-111914 was not analyzed since Trip Blank-111814 was submitted with the same group of samples and was non-detect for all analytes (except as discussed above with method blank contamination).
- The Equipment Blank sample (EB-1114) contained a low level detection of methylene chloride, just above the limit of quantitation (LOQ, 1.0 µg/L), at 1.4 µg/L. This trace detection was due to method blank contamination and was qualified as U, as discussed above.
- There were no other QA/QC issues noted. The validated laboratory data sheets are provided in Attachment 1.

## Attachments

|              |                                   |
|--------------|-----------------------------------|
| Table 1      | Summary of Qualifiers Applied     |
| Attachment 1 | Validated Analytical Report Pages |

## References

United States Environmental Protection Agency (EPA). 2014a. USEPA Contract Laboratory Program, National Functional Guidelines for Inorganic Superfund Data Review. OSWER 9355.0-131.EPA-540-R-013-001, August.

EPA. 2014b. USEPA Contract Laboratory Program, National Functional Guidelines for Superfund Organic Methods Data Review. OSWER 9355.0-132.EPA-540-R-014-002, August.

## TABLE

**Table 1: Summary of Qualifiers Applied**

| Method        | Sample ID(s)                                   | Parameter                                          | Qualifier  | Reason                                                     |
|---------------|------------------------------------------------|----------------------------------------------------|------------|------------------------------------------------------------|
| VOCs by 8260B | Trip Blank-111814<br>Trip Blank-112014         | Methylene Chloride                                 | J+         | Method blank contamination.                                |
| VOCs by 8260B | EB-1114                                        | Methylene Chloride                                 | 1.4 U µg/L | Method blank contamination.                                |
| VOCs by 8260B | Trip Blank-112014<br>LMW-11-1114<br>LMW-6-1114 | Vinyl Acetate<br>2-Chlorovinylether<br>Naphthalene | UJ         | Continuing calibration standard (CCAL) out of control low. |

Note: µg/L = micrograms per liter

**ATTACHMENT 1**  
**VALIDATED ANALYTICAL REPORT PAGES**

# Chain of Custody Record & Laboratory Analysis Request

|                                                      |                                           |                             |                                  |
|------------------------------------------------------|-------------------------------------------|-----------------------------|----------------------------------|
| ARI Assigned Number:<br><b>ZL38</b>                  | Turn-around Requested:<br><b>Standard</b> | Page: <b>1</b> of <b>1</b>  |                                  |
| ARI Client Company:<br><b>Golder Associates</b>      |                                           | Date:<br><b>11/18/2014</b>  | Ice Present?<br><b>Y</b>         |
| Client Contact:<br><b>Doug Morell, Jill Lamberts</b> |                                           | No. of Coolers:<br><b>7</b> | Cooler Temps:<br><b>0.41-4.7</b> |
| Client Project Name:<br><b>Landsburg</b>             |                                           |                             |                                  |
| Client Project #:<br><b>923-1000-002.R273</b>        | Samplers:<br><b>Rydecki, Lamberts</b>     |                             |                                  |



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants  
4611 South 134th Place, Suite 100  
Tukwila, WA 98168  
206-695-6200 206-695-6201 (fax)  
www.arilabs.com

| Sample ID                                         | Date     | Time | Matrix | No. Containers | Analysis Requested                                 |                               |            |            |            |                                                |          |       |               |       | Notes/Comments                  |             |  |  |  |                             |  |  |  |  |
|---------------------------------------------------|----------|------|--------|----------------|----------------------------------------------------|-------------------------------|------------|------------|------------|------------------------------------------------|----------|-------|---------------|-------|---------------------------------|-------------|--|--|--|-----------------------------|--|--|--|--|
|                                                   |          |      |        |                | VOCs                                               | Client list                   | PCBs (CLL) | Pesticides | SVOCs 8270 | Client list                                    | TPH-HClD | TAM/L | -Total Metals | TAM/L |                                 | Diss Metals |  |  |  |                             |  |  |  |  |
| Trip Blank-111814                                 | 11/18/14 | -    | W      | 9-6            | X                                                  |                               |            |            |            |                                                |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-10-1114                                       |          | 945  | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              | HOLD     |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-2-1114                                        |          | 1140 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-4-1114                                        |          | 1310 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| Trip Blank-111914                                 | 11/19/14 | -    | W      | 6              | X                                                  | HOLD unless there are detects |            |            |            |                                                |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-3-1114                                        |          | 0920 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              | HOLD     |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| EB-1114                                           |          | 0900 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-8-1114                                        |          | 1005 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-5-1114                                        |          | 1130 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| LMW-9-1114                                        |          | 1315 | W      | 17             | X                                                  | X                             | X          | X          | X          | X                                              |          |       |               |       |                                 |             |  |  |  |                             |  |  |  |  |
| Comments/Special Instructions<br>-Ecology ETM EDD |          |      |        |                | Relinquished by:<br>(Signature) <i>[Signature]</i> |                               |            |            |            | Received by:<br>(Signature) <i>[Signature]</i> |          |       |               |       | Relinquished by:<br>(Signature) |             |  |  |  | Received by:<br>(Signature) |  |  |  |  |
| Printed Name:<br><b>J. Lamberts</b>               |          |      |        |                | Printed Name:<br><b>J. Lamberts</b>                |                               |            |            |            | Printed Name:<br><b>Taylor Stricker</b>        |          |       |               |       | Printed Name:                   |             |  |  |  | Printed Name:               |  |  |  |  |
| Company:<br><b>Golder</b>                         |          |      |        |                | Company:<br><b>Golder</b>                          |                               |            |            |            | Company:<br><b>ARL</b>                         |          |       |               |       | Company:                        |             |  |  |  | Company:                    |  |  |  |  |
| Date & Time:<br><b>11/19/14 1442</b>              |          |      |        |                | Date & Time:<br><b>11/19/14 1442</b>               |                               |            |            |            | Date & Time:                                   |          |       |               |       | Date & Time:                    |             |  |  |  |                             |  |  |  |  |

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, notwithstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

Validated by J. Lamberts 1/26/2015

ZL38: 00002

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 2Sample ID: LMW-10-1114  
SAMPLE

Lab Sample ID: ZL38A

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: *mm*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Date Analyzed: 11/28/14 17:40

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

| CAS Number     | Analyte                               | DL          | LOQ         | Result        |
|----------------|---------------------------------------|-------------|-------------|---------------|
| 74-87-3        | Chloromethane                         | 0.09        | 0.50        | < 0.50 U      |
| 74-83-9        | Bromomethane                          | 0.25        | 1.0         | < 1.0 U       |
| 75-01-4        | Vinyl Chloride                        | 0.06        | 0.10        | < 0.10 U      |
| 75-00-3        | Chloroethane                          | 0.09        | 0.20        | < 0.20 U      |
| 75-09-2        | Methylene Chloride                    | 0.48        | 1.0         | < 1.0 U       |
| 67-64-1        | Acetone                               | 2.1         | 5.0         | < 5.0 U       |
| <b>75-15-0</b> | <b>Carbon Disulfide</b>               | <b>0.04</b> | <b>0.20</b> | <b>0.14 J</b> |
| 75-35-4        | 1,1-Dichloroethene                    | 0.05        | 0.20        | < 0.20 U      |
| 75-34-3        | 1,1-Dichloroethane                    | 0.05        | 0.20        | < 0.20 U      |
| 156-60-5       | trans-1,2-Dichloroethene              | 0.05        | 0.20        | < 0.20 U      |
| 156-59-2       | cis-1,2-Dichloroethene                | 0.04        | 0.20        | < 0.20 U      |
| 67-66-3        | Chloroform                            | 0.03        | 0.20        | < 0.20 U      |
| 107-06-2       | 1,2-Dichloroethane                    | 0.07        | 0.20        | < 0.20 U      |
| 78-93-3        | 2-Butanone                            | 0.81        | 5.0         | < 5.0 U       |
| 71-55-6        | 1,1,1-Trichloroethane                 | 0.04        | 0.20        | < 0.20 U      |
| 56-23-5        | Carbon Tetrachloride                  | 0.04        | 0.20        | < 0.20 U      |
| 108-05-4       | Vinyl Acetate                         | 0.07        | 0.20        | < 0.20 U      |
| 75-27-4        | Bromodichloromethane                  | 0.05        | 0.20        | < 0.20 U      |
| 78-87-5        | 1,2-Dichloropropane                   | 0.04        | 0.20        | < 0.20 U      |
| 10061-01-5     | cis-1,3-Dichloropropene               | 0.06        | 0.20        | < 0.20 U      |
| 79-01-6        | Trichloroethene                       | 0.05        | 0.20        | < 0.20 U      |
| 124-48-1       | Dibromochloromethane                  | 0.05        | 0.20        | < 0.20 U      |
| 79-00-5        | 1,1,2-Trichloroethane                 | 0.13        | 0.20        | < 0.20 U      |
| 71-43-2        | Benzene                               | 0.03        | 0.20        | < 0.20 U      |
| 10061-02-6     | trans-1,3-Dichloropropene             | 0.08        | 0.20        | < 0.20 U      |
| 110-75-8       | 2-Chloroethylvinylether               | 0.25        | 0.50        | < 0.50 U      |
| 75-25-2        | Bromoform                             | 0.06        | 0.20        | < 0.20 U      |
| 108-10-1       | 4-Methyl-2-Pentanone (MIBK)           | 0.97        | 2.5         | < 2.5 U       |
| 591-78-6       | 2-Hexanone                            | 0.90        | 5.0         | < 5.0 U       |
| 127-18-4       | Tetrachloroethene                     | 0.05        | 0.20        | < 0.20 U      |
| 79-34-5        | 1,1,2,2-Tetrachloroethane             | 0.06        | 0.10        | < 0.10 U      |
| 108-88-3       | Toluene                               | 0.04        | 0.20        | < 0.20 U      |
| 108-90-7       | Chlorobenzene                         | 0.02        | 0.20        | < 0.20 U      |
| 100-41-4       | Ethylbenzene                          | 0.04        | 0.20        | < 0.20 U      |
| 100-42-5       | Styrene                               | 0.05        | 0.20        | < 0.20 U      |
| 75-69-4        | Trichlorofluoromethane                | 0.04        | 0.20        | < 0.20 U      |
| 76-13-1        | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04        | 0.20        | < 0.20 U      |
| 179601-23-1    | m,p-Xylene                            | 0.05        | 0.40        | < 0.40 U      |
| 95-47-6        | o-Xylene                              | 0.03        | 0.20        | < 0.20 U      |
| 95-50-1        | 1,2-Dichlorobenzene                   | 0.04        | 0.20        | < 0.20 U      |
| 541-73-1       | 1,3-Dichlorobenzene                   | 0.04        | 0.20        | < 0.20 U      |
| 106-46-7       | 1,4-Dichlorobenzene                   | 0.04        | 0.20        | < 0.20 U      |
| 107-02-8       | Acrolein                              | 2.5         | 2.5         | < 2.5 U       |
| 74-88-4        | Iodomethane                           | 0.23        | 0.50        | < 0.50 U      |
| 107-13-1       | Acrylonitrile                         | 0.60        | 1.0         | < 1.0 U       |
| 563-58-6       | 1,1-Dichloropropene                   | 0.03        | 0.10        | < 0.10 U      |
| 74-95-3        | Dibromomethane                        | 0.14        | 0.20        | < 0.20 U      |
| 630-20-6       | 1,1,1,2-Tetrachloroethane             | 0.04        | 0.20        | < 0.20 U      |
| 96-12-8        | 1,2-Dibromo-3-chloropropane           | 0.04        | 0.50        | < 0.50 U      |
| 96-18-4        | 1,2,3-Trichloropropane                | 0.13        | 0.20        | < 0.20 U      |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-10-1114  
SAMPLE

Lab Sample ID: ZL38A

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25246

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 17:40

| CAS Number | Analyte                     | DL   | LOQ  | Result              |
|------------|-----------------------------|------|------|---------------------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U             |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U            |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U            |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | 0.33 <del>B</del> U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U            |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U            |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U            |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U            |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U            |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U            |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U            |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U            |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U            |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U            |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U            |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U            |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U            |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U            |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U            |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U            |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 102%  |
| d8-Toluene             | 100%  |
| Bromofluorobenzene     | 96.5% |
| d4-1,2-Dichlorobenzene | 101%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-2-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38B

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25247

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: 

Date Sampled: 11/18/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 18:07

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-2-1114  
SAMPLE

Lab Sample ID: ZL38B

LIMS ID: 14-25247

Matrix: Water

Date Analyzed: 11/28/14 18:07

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 105%  |
| d8-Toluene             | 98.0% |
| Bromofluorobenzene     | 96.5% |
| d4-1,2-Dichlorobenzene | 103%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-4-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38C

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25248

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: 

Date Sampled: 11/18/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 18:34

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-4-1114

Page 2 of 2

SAMPLE

Lab Sample ID: ZL38C

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25248

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 18:34

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 106%  |
| d8-Toluene             | 97.8% |
| Bromofluorobenzene     | 98.1% |
| d4-1,2-Dichlorobenzene | 106%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-3-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38D

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25249

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: 

Date Sampled: 11/19/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 19:00

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-3-1114  
SAMPLE

Lab Sample ID: ZL38D

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25249

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 19:00

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 103%  |
| d8-Toluene             | 102%  |
| Bromofluorobenzene     | 95.1% |
| d4-1,2-Dichlorobenzene | 105%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 2Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E

LIMS ID: 14-25250

Matrix: Water

Data Release Authorized: 

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 19:27

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result             |
|-------------|---------------------------------------|------|------|--------------------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U           |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U            |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U           |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U           |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | 1.4 <del>B</del> U |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U            |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U           |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U           |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U           |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U           |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U           |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U           |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U           |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U            |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U           |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U           |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U           |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U           |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U           |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U           |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U           |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U           |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U           |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U           |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U           |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U           |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U           |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U            |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U            |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U           |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U           |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U           |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U           |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U           |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U           |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U           |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U           |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U           |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U           |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U           |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U           |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U           |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U            |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U           |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U            |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U           |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U           |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U           |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U           |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U           |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: EB-1114

Page 2 of 2

SAMPLE

Lab Sample ID: ZL38E

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25250

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 19:27

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 104%  |
| d8-Toluene             | 98.6% |
| Bromofluorobenzene     | 95.7% |
| d4-1,2-Dichlorobenzene | 104%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 2Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F

LIMS ID: 14-25251

Matrix: Water

Data Release Authorized: *mmw*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Date Analyzed: 11/28/14 19:53

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25251

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 19:53

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 104%  |
| d8-Toluene             | 102%  |
| Bromofluorobenzene     | 94.7% |
| d4-1,2-Dichlorobenzene | 103%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 2Sample ID: LMW-5-1114  
SAMPLE

Lab Sample ID: ZL38G

LIMS ID: 14-25252

Matrix: Water

Data Release Authorized: 

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Date Analyzed: 11/28/14 20:20

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-5-1114

Page 2 of 2

SAMPLE

Lab Sample ID: ZL38G

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25252

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 20:20

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 108%  |
| d8-Toluene             | 99.6% |
| Bromofluorobenzene     | 95.8% |
| d4-1,2-Dichlorobenzene | 104%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-9-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38H

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25253

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: 

Date Sampled: 11/19/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 20:47

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result   |
|-------------|---------------------------------------|------|------|----------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U  |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U  |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U  |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U  |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U  |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U  |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U  |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U  |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 2 of 2Sample ID: LMW-9-1114  
SAMPLE

Lab Sample ID: ZL38H

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25253

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 20:47

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 107%  |
| d8-Toluene             | 100%  |
| Bromofluorobenzene     | 94.9% |
| d4-1,2-Dichlorobenzene | 105%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: TRIP BLANK-111814

Page 1 of 2

SAMPLE

Lab Sample ID: ZL38I

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25254

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: 

Date Sampled: 11/18/14

Reported: 12/03/14

Date Received: 11/19/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 11/28/14 21:14

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result              |
|-------------|---------------------------------------|------|------|---------------------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U            |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U             |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U            |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U            |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | 2.3 <del>B</del> J+ |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U             |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U            |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U            |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U            |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U            |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U            |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U            |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U            |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U             |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U            |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U            |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U            |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U            |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U            |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U            |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U            |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U            |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U            |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U            |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U            |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U            |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U            |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U             |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U             |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U            |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U            |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U            |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U            |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U            |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U            |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U            |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U            |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U            |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U            |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U            |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U            |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U            |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U             |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U            |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U             |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U            |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U            |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U            |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U            |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U            |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: TRIP BLANK-111814

Page 2 of 2

SAMPLE

Lab Sample ID: ZL38I

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25254

Project: Landsburg

Matrix: Water

923-1000-002.R273

Date Analyzed: 11/28/14 21:14

| CAS Number | Analyte                     | DL   | LOQ  | Result   |
|------------|-----------------------------|------|------|----------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U  |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 108%  |
| d8-Toluene             | 100%  |
| Bromofluorobenzene     | 94.2% |
| d4-1,2-Dichlorobenzene | 105%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Semivolatiles by SW8270D GC/MS

Extraction Method: SW3520C

Page 1 of 2

Sample ID: LMW-10-1114

SAMPLE

Lab Sample ID: ZL38A

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: 

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/24/14

Date Analyzed: 11/28/14 17:17

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-10-1114  
SAMPLE

Lab Sample ID: ZL38A  
LIMS ID: 14-25246  
Matrix: Water  
Date Analyzed: 11/28/14 17:17

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 78.4% | 2-Fluorobiphenyl       | 83.2% |
| d14-p-Terphenyl      | 86.0% | d4-1,2-Dichlorobenzene | 65.2% |
| d5-Phenol            | 83.7% | 2-Fluorophenol         | 69.6% |
| 2,4,6-Tribromophenol | 86.1% | d4-2-Chlorophenol      | 76.8% |

## ORGANICS ANALYSIS DATA SHEET

Semivolatiles by SW8270D GC/MS

Extraction Method: SW3520C

Page 1 of 2

Sample ID: LMW-2-1114

SAMPLE

Lab Sample ID: ZL38B

LIMS ID: 14-25247

Matrix: Water

Data Release Authorized: *AB*

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/24/14

Date Analyzed: 12/02/14 19:49

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-2-1114  
SAMPLE

Lab Sample ID: ZL38B  
LIMS ID: 14-25247  
Matrix: Water  
Date Analyzed: 12/02/14 19:49

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 48.4% | 2-Fluorobiphenyl       | 48.4% |
| d14-p-Terphenyl      | 53.6% | d4-1,2-Dichlorobenzene | 47.2% |
| d5-Phenol            | 58.7% | 2-Fluorophenol         | 51.2% |
| 2,4,6-Tribromophenol | 49.9% | d4-2-Chlorophenol      | 54.4% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-4-1114  
SAMPLE

Lab Sample ID: ZL38C  
LIMS ID: 14-25248  
Matrix: Water  
Data Release Authorized: *A*  
Reported: 12/03/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/18/14  
Date Received: 11/19/14

Date Extracted: 11/24/14  
Date Analyzed: 11/28/14 18:25  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-4-1114  
SAMPLE

Lab Sample ID: ZL38C  
LIMS ID: 14-25248  
Matrix: Water  
Date Analyzed: 11/28/14 18:25

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number      | Analyte                           | DL         | LOQ        | Result                                       |
|-----------------|-----------------------------------|------------|------------|----------------------------------------------|
| 84-66-2         | Diethylphthalate                  | 0.27       | 1.0        | < 1.0 U                                      |
| 7005-72-3       | 4-Chlorophenyl-phenylether        | 0.27       | 1.0        | < 1.0 U                                      |
| 86-73-7         | Fluorene                          | 0.29       | 1.0        | < 1.0 U                                      |
| 100-01-6        | 4-Nitroaniline                    | 2.0        | 3.0        | < 3.0 U                                      |
| 534-52-1        | 4,6-Dinitro-2-Methylphenol        | 3.6        | 10         | < 10 U                                       |
| 86-30-6         | N-Nitrosodiphenylamine            | 0.30       | 1.0        | < 1.0 U                                      |
| 101-55-3        | 4-Bromophenyl-phenylether         | 0.24       | 1.0        | < 1.0 U                                      |
| 118-74-1        | Hexachlorobenzene                 | 0.28       | 1.0        | < 1.0 U                                      |
| 87-86-5         | Pentachlorophenol                 | 1.9        | 10         | < 10 U                                       |
| 85-01-8         | Phenanthrene                      | 0.32       | 1.0        | < 1.0 U                                      |
| 86-74-8         | Carbazole                         | 0.31       | 1.0        | < 1.0 U                                      |
| 120-12-7        | Anthracene                        | 0.26       | 1.0        | < 1.0 U                                      |
| 84-74-2         | Di-n-Butylphthalate               | 0.29       | 1.0        | < 1.0 U                                      |
| 206-44-0        | Fluoranthene                      | 0.30       | 1.0        | < 1.0 U                                      |
| 129-00-0        | Pyrene                            | 0.28       | 1.0        | < 1.0 U                                      |
| 85-68-7         | Butylbenzylphthalate              | 0.30       | 1.0        | < 1.0 U                                      |
| 91-94-1         | 3,3'-Dichlorobenzidine            | 1.8        | 5.0        | < 5.0 U                                      |
| 56-55-3         | Benzo(a)anthracene                | 0.29       | 1.0        | < 1.0 U                                      |
| <b>117-81-7</b> | <b>bis(2-Ethylhexyl)phthalate</b> | <b>2.1</b> | <b>3.0</b> | <b>12</b> <span style="color: red;">R</span> |
| 218-01-9        | Chrysene                          | 0.32       | 1.0        | < 1.0 U                                      |
| 117-84-0        | Di-n-Octyl phthalate              | 0.27       | 1.0        | < 1.0 U                                      |
| 205-99-2        | Benzo(b)fluoranthene              | 0.32       | 1.0        | < 1.0 U                                      |
| 207-08-9        | Benzo(k)fluoranthene              | 0.34       | 1.0        | < 1.0 U                                      |
| 50-32-8         | Benzo(a)pyrene                    | 0.30       | 1.0        | < 1.0 U                                      |
| 193-39-5        | Indeno(1,2,3-cd)pyrene            | 0.36       | 1.0        | < 1.0 U                                      |
| 53-70-3         | Dibenz(a,h)anthracene             | 0.39       | 1.0        | < 1.0 U                                      |
| 191-24-2        | Benzo(g,h,i)perylene              | 0.39       | 1.0        | < 1.0 U                                      |
| 108-39-4        | 3&4-Methylphenol                  | 0.80       | 2.0        | < 2.0 U                                      |
| 90-12-0         | 1-Methylnaphthalene               | 0.26       | 1.0        | < 1.0 U                                      |
| TOTBFA          | Total Benzofluoranthenes          | 0.80       | 2.0        | < 2.0 U                                      |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 72.8% | 2-Fluorobiphenyl       | 68.8% |
| d14-p-Terphenyl      | 72.4% | d4-1,2-Dichlorobenzene | 63.2% |
| d5-Phenol            | 77.3% | 2-Fluorophenol         | 69.3% |
| 2,4,6-Tribromophenol | 69.6% | d4-2-Chlorophenol      | 71.5% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-3-1114  
SAMPLE

Lab Sample ID: ZL38D  
LIMS ID: 14-25249  
Matrix: Water  
Data Release Authorized: *RB*  
Reported: 12/03/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/24/14  
Date Analyzed: 11/28/14 18:59  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-3-1114  
SAMPLE

Lab Sample ID: ZL38D  
LIMS ID: 14-25249  
Matrix: Water  
Date Analyzed: 11/28/14 18:59

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 66.8% | 2-Fluorobiphenyl       | 64.4% |
| d14-p-Terphenyl      | 70.0% | d4-1,2-Dichlorobenzene | 60.8% |
| d5-Phenol            | 69.9% | 2-Fluorophenol         | 61.1% |
| 2,4,6-Tribromophenol | 63.7% | d4-2-Chlorophenol      | 65.1% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E  
LIMS ID: 14-25250  
Matrix: Water  
Data Release Authorized:   
Reported: 12/03/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/24/14  
Date Analyzed: 11/28/14 19:33  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E  
LIMS ID: 14-25250  
Matrix: Water  
Date Analyzed: 11/28/14 19:33

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 74.0% | 2-Fluorobiphenyl       | 71.2% |
| d14-p-Terphenyl      | 81.2% | d4-1,2-Dichlorobenzene | 66.0% |
| d5-Phenol            | 76.8% | 2-Fluorophenol         | 68.0% |
| 2,4,6-Tribromophenol | 69.3% | d4-2-Chlorophenol      | 72.8% |

## ORGANICS ANALYSIS DATA SHEET

Semivolatiles by SW8270D GC/MS

Extraction Method: SW3520C

Page 1 of 2

Sample ID: LMW-8-1114

SAMPLE

Lab Sample ID: ZL38F

LIMS ID: 14-25251

Matrix: Water

Data Release Authorized: 

Reported: 12/03/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/24/14

Date Analyzed: 11/28/14 20:06

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F  
LIMS ID: 14-25251  
Matrix: Water  
Date Analyzed: 11/28/14 20:06

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 63.6% | 2-Fluorobiphenyl       | 60.8% |
| d14-p-Terphenyl      | 68.8% | d4-1,2-Dichlorobenzene | 56.0% |
| d5-Phenol            | 67.5% | 2-Fluorophenol         | 59.2% |
| 2,4,6-Tribromophenol | 62.1% | d4-2-Chlorophenol      | 61.9% |

**ORGANICS ANALYSIS DATA SHEET**  
**Semivolatiles by SW8270D GC/MS**  
**Extraction Method: SW3520C**  
 Page 1 of 2

**Sample ID: LMW-5-1114**  
**SAMPLE**

Lab Sample ID: ZL38G  
 LIMS ID: 14-25252  
 Matrix: Water  
 Data Release Authorized:   
 Reported: 02/04/15

QC Report No: ZL38-Golder Associates  
 Project: Landsburg  
 923-1000-002.R273  
 Date Sampled: 11/19/14  
 Date Received: 11/19/14

Date Extracted: 11/24/14  
 Date Analyzed: 11/28/14 20:40  
 Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

**ORGANICS ANALYSIS DATA SHEET**  
**Semivolatiles by SW8270D GC/MS**  
**Extraction Method: SW3520C**  
 Page 2 of 2

**Sample ID: LMW-5-1114**  
**SAMPLE**

Lab Sample ID: ZL38G  
 LIMS ID: 14-25252  
 Matrix: Water  
 Date Analyzed: 11/28/14 20:40

QC Report No: ZL38-Golder Associates  
 Project: Landsburg  
 923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 54.4% | 2-Fluorobiphenyl       | 54.8% |
| d14-p-Terphenyl      | 55.6% | d4-1,2-Dichlorobenzene | 46.8% |
| d5-Phenol            | 57.6% | 2-Fluorophenol         | 50.9% |
| 2,4,6-Tribromophenol | 53.6% | d4-2-Chlorophenol      | 53.6% |

**ORGANICS ANALYSIS DATA SHEET**  
**Semivolatiles by SW8270D GC/MS**  
**Extraction Method: SW3520C**  
 Page 1 of 2

**Sample ID: LMW-9-1114**  
**SAMPLE**

Lab Sample ID: ZL38H

LIMS ID: 14-25253

Matrix: Water

Data Release Authorized: 

Reported: 02/04/15

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/24/14

Date Analyzed: 11/28/14 21:14

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

**ORGANICS ANALYSIS DATA SHEET**  
**Semivolatiles by SW8270D GC/MS**  
**Extraction Method: SW3520C**  
 Page 2 of 2

**Sample ID: LMW-9-1114**  
**SAMPLE**

Lab Sample ID: ZL38H  
 LIMS ID: 14-25253  
 Matrix: Water  
 Date Analyzed: 11/28/14 21:14

QC Report No: ZL38-Golder Associates  
 Project: Landsburg  
 923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 70.8% | 2-Fluorobiphenyl       | 63.2% |
| d14-p-Terphenyl      | 65.2% | d4-1,2-Dichlorobenzene | 57.6% |
| d5-Phenol            | 64.0% | 2-Fluorophenol         | 56.5% |
| 2,4,6-Tribromophenol | 65.9% | d4-2-Chlorophenol      | 58.9% |

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-10-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38A

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25246

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/18/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 15:01

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

**Pest/PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 58.8% |
| Tetrachlorometaxylene | 75.5% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-2-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38B

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25247

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/18/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 15:19

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 64.8% |
| Tetrachlorometaxylene | 84.8% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-4-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38C

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25248

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *mm*

Date Sampled: 11/18/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 15:36

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 69.0% |
| Tetrachlorometaxylene | 116%  |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-3-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38D

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25249

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/19/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 15:53

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

**Pest/PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 68.2% |
| Tetrachlorometaxylene | 83.8% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: EB-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38E

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25250

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: 

Date Sampled: 11/19/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 16:10

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 65.2% |
| Tetrachlorometaxylene | 82.8% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-8-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38F

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25251

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MW*

Date Sampled: 11/19/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 16:27

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 52.2% |
| Tetrachlorometaxylene | 75.5% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-5-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38G

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25252

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MM*

Date Sampled: 11/19/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 16:45

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

**Pest/PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 77.5% |
| Tetrachlorometaxylene | 109%  |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Sample ID: LMW-9-1114

Extraction Method: SW3510C

SAMPLE

Page 1 of 1

Lab Sample ID: ZL38H

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25253

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *MMW*

Date Sampled: 11/19/14

Reported: 12/17/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Sample Amount: 500 mL

Date Analyzed: 12/16/14 17:02

Final Extract Volume: 5.0 mL

Instrument/Analyst: ECD6/YZ

Dilution Factor: 1.00

GPC Cleanup: No

Silica Gel: No

Sulfur Cleanup: No

Florisil Cleanup: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 83.8% |
| Tetrachlorometaxylene | 89.2% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-10-1114  
SAMPLE

Lab Sample ID: ZL38A

LIMS ID: 14-25246

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 18:15

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: Yes

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 66.8% |
| Tetrachlorometaxylene | 46.0% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-2-1114  
SAMPLE

Lab Sample ID: ZL38B

LIMS ID: 14-25247

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 18:34

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 74.2% |
| Tetrachlorometaxylene | 56.5% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-4-1114  
SAMPLE

Lab Sample ID: ZL38C

LIMS ID: 14-25248

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 18:54

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 76.8% |
| Tetrachlorometaxylene | 55.5% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-3-1114  
SAMPLE

Lab Sample ID: ZL38D

LIMS ID: 14-25249

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 19:13

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 70.5% |
| Tetrachlorometaxylene | 55.2% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E  
LIMS ID: 14-25250  
Matrix: Water  
Data Release Authorized:   
Reported: 12/10/14

QC Report No: ZL38-Golder Associates  
Project: Landsburg  
923-1000-002.R273  
Date Sampled: 11/19/14  
Date Received: 11/19/14

Date Extracted: 11/25/14  
Date Analyzed: 12/08/14 19:33  
Instrument/Analyst: ECD5/JGR  
GPC Cleanup: No  
Sulfur Cleanup: Yes

Sample Amount: 1000 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00  
Silica Gel: No  
Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 50.0% |
| Tetrachlorometaxylene | 54.8% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F

LIMS ID: 14-25251

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 19:53

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 66.0% |
| Tetrachlorometaxylene | 57.5% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-5-1114  
SAMPLE

Lab Sample ID: ZL38G

LIMS ID: 14-25252

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 20:51

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 67.8% |
| Tetrachlorometaxylene | 56.2% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-9-1114  
SAMPLE

Lab Sample ID: ZL38H

LIMS ID: 14-25253

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

Date Extracted: 11/25/14

Date Analyzed: 12/08/14 21:11

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 74.8% |
| Tetrachlorometaxylene | 53.8% |

**ORGANICS ANALYSIS DATA SHEET**

NWTPH-HCID Method by GC/FID

Extraction Method: SW3510C

Page 1 of 2

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Matrix: Water

Data Release Authorized: *MMW*

Reported: 11/25/14

| ARI ID                | Sample ID                 | Extraction Date | Analysis Date | DL  | Range       | Result   |
|-----------------------|---------------------------|-----------------|---------------|-----|-------------|----------|
| MB-112114<br>14-25246 | Method Blank              | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 79.4%    |
| ZL38A<br>14-25246     | LMW-10-1114<br>HC ID: --- | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 91.4%    |
| ZL38B<br>14-25247     | LMW-2-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 91.5%    |
| ZL38C<br>14-25248     | LMW-4-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 88.6%    |
| ZL38D<br>14-25249     | LMW-3-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 83.6%    |
| ZL38E<br>14-25250     | EB-1114<br>HC ID: ---     | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 90.2%    |
| ZL38F<br>14-25251     | LMW-8-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 86.0%    |
| ZL38G<br>14-25252     | LMW-5-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 92.3%    |
| ZL38H<br>14-25253     | LMW-9-1114<br>HC ID: ---  | 11/21/14        | 11/22/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 78.2%    |

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LMW-10-1114

SAMPLE

Lab Sample ID: ZL38A

QC Report No: ZL38-Golder Associates

LIMS ID: 14-25246

Project: Landsburg

Matrix: Water

923-1000-002.R273

Data Release Authorized: *XellH*

Date Sampled: 11/18/14

Reported: 12/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>7,840</b>  |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>300</b>    |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>2,950</b>  |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-96-5        | Manganese        | 0.3  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>1,330</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>81,300</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET****TOTAL METALS**

Page 1 of 1

**Sample ID: LMW-2-1114**  
**SAMPLE**

Lab Sample ID: ZL38B

LIMS ID: 14-25247

Matrix: Water

Data Release Authorized: *Kelly*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result         | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|----------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000          | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3              | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>113,000</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000          | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>250</b>     |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>69,200</b>  |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>220</b>     |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>3,590</b>   |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>20,700</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20             | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET****TOTAL METALS**

Page 1 of 1

Sample ID: LMW-4-1114  
SAMPLE

Lab Sample ID: ZL38C

LIMS ID: 14-25248

Matrix: Water

Data Release Authorized: *Xe/14*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result         | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|----------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000          | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3              | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>110,000</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000          | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>1,090</b>   |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>66,300</b>  |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>160</b>     |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>3,730</b>   |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>28,400</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3              | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20             | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET****TOTAL METALS**

Page 1 of 1

**Sample ID: LMW-3-1114**  
**SAMPLE**

Lab Sample ID: ZL38D

LIMS ID: 14-25249

Matrix: Water

Data Release Authorized: *hlm*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>38,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-89-6        | Iron             | 7.5  | 200   | 200           | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>16,000</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>60</b>     |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>1,720</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>10,200</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET****TOTAL METALS**

Page 1 of 1

Sample ID: EB-1114  
SAMPLE

Lab Sample ID: ZL38E

LIMS ID: 14-25250

Matrix: Water

Data Release Authorized: *12/14*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte   | DL   | LOQ   | Result | Q |
|-----------|-----------|-----------------|---------------|------------|-----------|------|-------|--------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5  | Aluminum  | 7.6  | 1,000 | 1,000  | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0  | Antimony  | 0.01 | 3     | 3      | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2  | Arsenic   | 0.05 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3  | Barium    | 1.3  | 500   | 500    | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7  | Beryllium | 0.16 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9  | Cadmium   | 0.18 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-70-2  | Calcium   | 11.3 | 500   | 500    | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3  | Chromium  | 1.2  | 1,000 | 1,000  | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4  | Cobalt    | 0.3  | 10    | 10     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8  | Copper    | 0.92 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-89-6  | Iron      | 7.5  | 200   | 200    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1  | Lead      | 0.0  | 10    | 10     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-95-4  | Magnesium | 9.6  | 1,000 | 1,000  | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-96-5  | Manganese | 0.3  | 20    | 20     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0  | Nickel    | 3.9  | 20    | 20     | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-09-7  | Potassium | 65.7 | 500   | 500    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2  | Selenium  | 0.13 | 5     | 5      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4  | Silver    | 0.43 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-23-5  | Sodium    | 11.4 | 500   | 500    | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0  | Thallium  | 0.00 | 2     | 2      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2  | Vanadium  | 0.27 | 3     | 3      | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6  | Zinc      | 1.4  | 20    | 20     | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET****TOTAL METALS**

Page 1 of 1

Sample ID: LMW-8-1114  
SAMPLE

Lab Sample ID: ZL38F

LIMS ID: 14-25251

Matrix: Water

Data Release Authorized: *[Signature]*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>63,200</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>12,400</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>33,700</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>510</b>    |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>2,120</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>11,500</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET****TOTAL METALS**

Page 1 of 1

**Sample ID: LMW-5-1114**  
**SAMPLE**

Lab Sample ID: ZL38G

LIMS ID: 14-25252

Matrix: Water

Data Release Authorized: *XLM*

Reported: 12/19/14

QC Report No: ZL38-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/19/14

Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>93,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7439-89-6        | Iron             | 7.5  | 200   | 200           | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>52,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>230</b>    |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>2,770</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>16,400</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET****TOTAL METALS**

Page 1 of 1

**Sample ID: LMW-9-1114**  
**SAMPLE**

 Lab Sample ID: ZL38H *helly*  
 LIMS ID: 14-25253  
 Matrix: Water  
 Data Release Authorized:  
 Reported: 12/19/14

 QC Report No: ZL38-Golder Associates  
 Project: Landsburg  
 923-1000-002.R273  
 Date Sampled: 11/19/14  
 Date Received: 11/19/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>81,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>1,510</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>44,300</b> |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>160</b>    |   |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>2,490</b>  |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>15,500</b> |   |
| 200.8     | 11/21/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/21/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

**INORGANICS ANALYSIS DATA SHEET**  
**Total Mercury by Method SW7470A**

Data Release Authorized:   
Reported: 11/21/14  
Date Received: 11/19/14  
Page 1 of 1

QC Report No: ZL39-Golder Associates  
Project: Landsburg  
923-1000-002.R273

| Client/<br>ARI ID             | Date<br>Sampled | Matrix | Prep Date<br>Anal Date | RL   | Result |
|-------------------------------|-----------------|--------|------------------------|------|--------|
| LMW-10-1114<br>ZL39A 14-25256 | 11/18/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-2-1114<br>ZL39B 14-25257  | 11/18/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-4-1114<br>ZL39C 14-25258  | 11/18/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-3-1114<br>ZL39D 14-25259  | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| EB-1114<br>ZL39E 14-25260     | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-8-1114<br>ZL39F 14-25261  | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-5-1114<br>ZL39G 14-25262  | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| LMW-9-1114<br>ZL39H 14-25263  | 11/19/14        | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |
| MB-112114<br>Method Blank     | NA              | Water  | 11/21/14<br>11/21/14   | 20.0 | 20.0 U |

**Reported in ng/L**

RL-Analytical reporting limit  
U-Undetected at reported detection limit

# Chain of Custody Record & Laboratory Analysis Request

ARI Assigned Number: 7165 Turn-around Requested: Standard

ARI Client Company: Golden Phone: 425-883-0772

Client Contact: D. Morell, J. Lamberts

Client Project Name: Landsburg

Client Project #: Q231000002.R273 Samplers: Rijdecki, Lamberts

Page: 1 of 1

Date: 11/20/2014 Ice Present? Y

No. of Coolers: 3 Cooler Temps: 46.38  
39



Analytical Resources, Incorporated  
Analytical Chemists and Consultants  
4611 South 134th Place, Suite 100  
Tukwila, WA 98168  
206-695-6200 206-695-6201 (fax)

| Sample ID                                                                                                                                               | Date                                                                                                                                                   | Time                                                                                                                                             | Matrix                                                                       | No. Containers                                                           | Analysis Requested |              |              |              |              |              |              |               |      |               | Notes/Comments |  |                |
|---------------------------------------------------------------------------------------------------------------------------------------------------------|--------------------------------------------------------------------------------------------------------------------------------------------------------|--------------------------------------------------------------------------------------------------------------------------------------------------|------------------------------------------------------------------------------|--------------------------------------------------------------------------|--------------------|--------------|--------------|--------------|--------------|--------------|--------------|---------------|------|---------------|----------------|--|----------------|
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          | VOCs client list   | PCBs (LL)    | Pesticides   | SVD CS 8270  | client list  | TPH-HCID     | TAML         | -Total Metals | TAML | -Diss. Metals |                |  |                |
| Trip Blank - 11/20/14                                                                                                                                   | 11/20/14                                                                                                                                               | -                                                                                                                                                | W                                                                            | 6                                                                        | X                  |              |              |              |              |              |              |               |      |               |                |  |                |
| LMW-11-1114                                                                                                                                             |                                                                                                                                                        | 1045                                                                                                                                             | W                                                                            | 17                                                                       | X                  | X            | X            | X            | X            | X            | X            | HOLD          |      |               |                |  |                |
| LMW-6-1114                                                                                                                                              |                                                                                                                                                        | 1240                                                                                                                                             | W                                                                            | 17                                                                       | X                  | X            | X            | X            | X            | X            | X            |               |      |               |                |  |                |
| <del>LMW-7-1114</del>                                                                                                                                   |                                                                                                                                                        | -                                                                                                                                                | W                                                                            | <del>17</del>                                                            | <del>X</del>       | <del>X</del> | <del>X</del> | <del>X</del> | <del>X</del> | <del>X</del> | <del>X</del> |               |      |               |                |  | Did not sample |
| <del>LMW-7-1114-D</del>                                                                                                                                 |                                                                                                                                                        | -                                                                                                                                                | W                                                                            | <del>17</del>                                                            | <del>X</del>       | <del>X</del> | <del>X</del> | <del>X</del> | <del>X</del> | <del>X</del> | <del>X</del> |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
|                                                                                                                                                         |                                                                                                                                                        |                                                                                                                                                  |                                                                              |                                                                          |                    |              |              |              |              |              |              |               |      |               |                |  |                |
| Comments/Special Instructions<br>- Ecology EIM EDD<br>**CLIENT SPECIFIC RLS<br>and ANALYTE LIST**<br>Pls cc j.lamberts@golder.com<br>dmorell@golder.com | Relinquished by:<br>(Signature) <u>[Signature]</u><br>Printed Name: <u>J. Lamberts</u><br>Company: <u>Golden</u><br>Date & Time: <u>11/20/14 17:48</u> | Received by:<br>(Signature) <u>[Signature]</u><br>Printed Name: <u>A. Vargardsen</u><br>Company: <u>ARI</u><br>Date & Time: <u>11/20/14 1748</u> | Relinquished by:<br>(Signature)<br>Printed Name:<br>Company:<br>Date & Time: | Received by:<br>(Signature)<br>Printed Name:<br>Company:<br>Date & Time: |                    |              |              |              |              |              |              |               |      |               |                |  |                |

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, notwithstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

Validated by J. Lamberts 1/26/2015

2165:00002

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-11-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL65A

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25426

Project: Landsburg

Matrix: Water

9231000002.R273

Data Release Authorized: 

Date Sampled: 11/20/14

Reported: 12/03/14

Date Received: 11/20/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 12/01/14 14:09

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result      |
|-------------|---------------------------------------|------|------|-------------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U    |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U     |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U    |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U    |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U     |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U     |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U    |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U    |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U    |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U    |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U    |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U    |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U    |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U     |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U    |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U    |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U UJ |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U    |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U    |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U    |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U    |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U    |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U    |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U    |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U    |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U UJ |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U    |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U     |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U     |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U    |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U    |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U    |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U    |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U    |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U    |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U    |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U    |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U    |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U    |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U    |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U    |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U    |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U     |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U    |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U     |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U    |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U    |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U    |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U    |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U    |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-11-1114

Page 2 of 2

SAMPLE

Lab Sample ID: ZL65A

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25426

Project: Landsburg

Matrix: Water

9231000002.R273

Date Analyzed: 12/01/14 14:09

| CAS Number | Analyte                     | DL   | LOQ  | Result      |
|------------|-----------------------------|------|------|-------------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U     |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U    |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U    |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U    |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U    |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U    |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U    |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U    |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U    |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U    |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U    |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U    |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U    |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U    |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U    |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U    |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U    |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U    |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U UJ |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U    |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 102%  |
| d8-Toluene             | 104%  |
| Bromofluorobenzene     | 92.7% |
| d4-1,2-Dichlorobenzene | 103%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-6-1114

Page 1 of 2

SAMPLE

Lab Sample ID: ZL65B

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25427

Project: Landsburg

Matrix: Water

9231000002.R273

Data Release Authorized: 

Date Sampled: 11/20/14

Reported: 12/03/14

Date Received: 11/20/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 12/01/14 14:36

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result      |
|-------------|---------------------------------------|------|------|-------------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U    |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U     |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U    |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U    |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | < 1.0 U     |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U     |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U    |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U    |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U    |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U    |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U    |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U    |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U    |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U     |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U    |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U    |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U UJ |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U    |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U    |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U    |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U    |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U    |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U    |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U    |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U    |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U UJ |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U    |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U     |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U     |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U    |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U    |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U    |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U    |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U    |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U    |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U    |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U    |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U    |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U    |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U    |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U    |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U    |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U     |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U    |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U     |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U    |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U    |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U    |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U    |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U    |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: LMW-6-1114

Page 2 of 2

SAMPLE

Lab Sample ID: ZL65B

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25427

Project: Landsburg

Matrix: Water

9231000002.R273

Date Analyzed: 12/01/14 14:36

| CAS Number | Analyte                     | DL   | LOQ  | Result             |
|------------|-----------------------------|------|------|--------------------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U            |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U           |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U           |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U           |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U           |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U           |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U           |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U           |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U           |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U           |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U           |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U           |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U           |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U           |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U           |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U           |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U           |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U           |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U <b>UJ</b> |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U           |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 106%  |
| d8-Toluene             | 101%  |
| Bromofluorobenzene     | 95.1% |
| d4-1,2-Dichlorobenzene | 106%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: TRIP BLANK-112014

Page 1 of 2

SAMPLE

Lab Sample ID: ZL65C

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25428

Project: Landsburg

Matrix: Water

9231000002.R273

Data Release Authorized: 

Date Sampled: 11/20/14

Reported: 12/03/14

Date Received: 11/20/14

Instrument/Analyst: NT2/LH

Sample Amount: 10.0 mL

Date Analyzed: 12/01/14 15:03

Purge Volume: 10.0 mL

| CAS Number  | Analyte                               | DL   | LOQ  | Result              |
|-------------|---------------------------------------|------|------|---------------------|
| 74-87-3     | Chloromethane                         | 0.09 | 0.50 | < 0.50 U            |
| 74-83-9     | Bromomethane                          | 0.25 | 1.0  | < 1.0 U             |
| 75-01-4     | Vinyl Chloride                        | 0.06 | 0.10 | < 0.10 U            |
| 75-00-3     | Chloroethane                          | 0.09 | 0.20 | < 0.20 U            |
| 75-09-2     | Methylene Chloride                    | 0.48 | 1.0  | 2.6 <del>B</del> J+ |
| 67-64-1     | Acetone                               | 2.1  | 5.0  | < 5.0 U             |
| 75-15-0     | Carbon Disulfide                      | 0.04 | 0.20 | < 0.20 U            |
| 75-35-4     | 1,1-Dichloroethene                    | 0.05 | 0.20 | < 0.20 U            |
| 75-34-3     | 1,1-Dichloroethane                    | 0.05 | 0.20 | < 0.20 U            |
| 156-60-5    | trans-1,2-Dichloroethene              | 0.05 | 0.20 | < 0.20 U            |
| 156-59-2    | cis-1,2-Dichloroethene                | 0.04 | 0.20 | < 0.20 U            |
| 67-66-3     | Chloroform                            | 0.03 | 0.20 | < 0.20 U            |
| 107-06-2    | 1,2-Dichloroethane                    | 0.07 | 0.20 | < 0.20 U            |
| 78-93-3     | 2-Butanone                            | 0.81 | 5.0  | < 5.0 U             |
| 71-55-6     | 1,1,1-Trichloroethane                 | 0.04 | 0.20 | < 0.20 U            |
| 56-23-5     | Carbon Tetrachloride                  | 0.04 | 0.20 | < 0.20 U            |
| 108-05-4    | Vinyl Acetate                         | 0.07 | 0.20 | < 0.20 U UJ         |
| 75-27-4     | Bromodichloromethane                  | 0.05 | 0.20 | < 0.20 U            |
| 78-87-5     | 1,2-Dichloropropane                   | 0.04 | 0.20 | < 0.20 U            |
| 10061-01-5  | cis-1,3-Dichloropropene               | 0.06 | 0.20 | < 0.20 U            |
| 79-01-6     | Trichloroethene                       | 0.05 | 0.20 | < 0.20 U            |
| 124-48-1    | Dibromochloromethane                  | 0.05 | 0.20 | < 0.20 U            |
| 79-00-5     | 1,1,2-Trichloroethane                 | 0.13 | 0.20 | < 0.20 U            |
| 71-43-2     | Benzene                               | 0.03 | 0.20 | < 0.20 U            |
| 10061-02-6  | trans-1,3-Dichloropropene             | 0.08 | 0.20 | < 0.20 U            |
| 110-75-8    | 2-Chloroethylvinylether               | 0.25 | 0.50 | < 0.50 U UJ         |
| 75-25-2     | Bromoform                             | 0.06 | 0.20 | < 0.20 U            |
| 108-10-1    | 4-Methyl-2-Pentanone (MIBK)           | 0.97 | 2.5  | < 2.5 U             |
| 591-78-6    | 2-Hexanone                            | 0.90 | 5.0  | < 5.0 U             |
| 127-18-4    | Tetrachloroethene                     | 0.05 | 0.20 | < 0.20 U            |
| 79-34-5     | 1,1,2,2-Tetrachloroethane             | 0.06 | 0.10 | < 0.10 U            |
| 108-88-3    | Toluene                               | 0.04 | 0.20 | < 0.20 U            |
| 108-90-7    | Chlorobenzene                         | 0.02 | 0.20 | < 0.20 U            |
| 100-41-4    | Ethylbenzene                          | 0.04 | 0.20 | < 0.20 U            |
| 100-42-5    | Styrene                               | 0.05 | 0.20 | < 0.20 U            |
| 75-69-4     | Trichlorofluoromethane                | 0.04 | 0.20 | < 0.20 U            |
| 76-13-1     | 1,1,2-Trichloro-1,2,2-trifluoroethane | 0.04 | 0.20 | < 0.20 U            |
| 179601-23-1 | m,p-Xylene                            | 0.05 | 0.40 | < 0.40 U            |
| 95-47-6     | o-Xylene                              | 0.03 | 0.20 | < 0.20 U            |
| 95-50-1     | 1,2-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U            |
| 541-73-1    | 1,3-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U            |
| 106-46-7    | 1,4-Dichlorobenzene                   | 0.04 | 0.20 | < 0.20 U            |
| 107-02-8    | Acrolein                              | 2.5  | 2.5  | < 2.5 U             |
| 74-88-4     | Iodomethane                           | 0.23 | 0.50 | < 0.50 U            |
| 107-13-1    | Acrylonitrile                         | 0.60 | 1.0  | < 1.0 U             |
| 563-58-6    | 1,1-Dichloropropene                   | 0.03 | 0.10 | < 0.10 U            |
| 74-95-3     | Dibromomethane                        | 0.14 | 0.20 | < 0.20 U            |
| 630-20-6    | 1,1,1,2-Tetrachloroethane             | 0.04 | 0.20 | < 0.20 U            |
| 96-12-8     | 1,2-Dibromo-3-chloropropane           | 0.04 | 0.50 | < 0.50 U            |
| 96-18-4     | 1,2,3-Trichloropropane                | 0.13 | 0.20 | < 0.20 U            |

## ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge &amp; Trap GC/MS-Method SW8260C

Sample ID: TRIP BLANK-112014

Page 2 of 2

SAMPLE

Lab Sample ID: ZL65C

QC Report No: ZL65-Golder Associates

LIMS ID: 14-25428

Project: Landsburg

Matrix: Water

9231000002.R273

Date Analyzed: 12/01/14 15:03

| CAS Number | Analyte                     | DL   | LOQ  | Result      |
|------------|-----------------------------|------|------|-------------|
| 110-57-6   | trans-1,4-Dichloro-2-butene | 0.32 | 1.0  | < 1.0 U     |
| 108-67-8   | 1,3,5-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U    |
| 95-63-6    | 1,2,4-Trimethylbenzene      | 0.02 | 0.20 | < 0.20 U    |
| 87-68-3    | Hexachlorobutadiene         | 0.07 | 0.20 | < 0.20 U    |
| 106-93-4   | 1,2-Dibromoethane           | 0.07 | 0.10 | < 0.10 U    |
| 74-97-5    | Bromochloromethane          | 0.06 | 0.20 | < 0.20 U    |
| 594-20-7   | 2,2-Dichloropropane         | 0.05 | 0.10 | < 0.10 U    |
| 142-28-9   | 1,3-Dichloropropane         | 0.06 | 0.10 | < 0.10 U    |
| 98-82-8    | Isopropylbenzene            | 0.02 | 0.20 | < 0.20 U    |
| 103-65-1   | n-Propylbenzene             | 0.02 | 0.20 | < 0.20 U    |
| 108-86-1   | Bromobenzene                | 0.06 | 0.20 | < 0.20 U    |
| 95-49-8    | 2-Chlorotoluene             | 0.02 | 0.10 | < 0.10 U    |
| 106-43-4   | 4-Chlorotoluene             | 0.02 | 0.20 | < 0.20 U    |
| 98-06-6    | tert-Butylbenzene           | 0.03 | 0.20 | < 0.20 U    |
| 135-98-8   | sec-Butylbenzene            | 0.02 | 0.20 | < 0.20 U    |
| 99-87-6    | 4-Isopropyltoluene          | 0.03 | 0.10 | < 0.10 U    |
| 104-51-8   | n-Butylbenzene              | 0.02 | 0.20 | < 0.20 U    |
| 120-82-1   | 1,2,4-Trichlorobenzene      | 0.11 | 0.50 | < 0.50 U    |
| 91-20-3    | Naphthalene                 | 0.12 | 0.50 | < 0.50 U UJ |
| 87-61-6    | 1,2,3-Trichlorobenzene      | 0.11 | 0.20 | < 0.20 U    |

Reported in µg/L (ppb)

## Volatile Surrogate Recovery

|                        |       |
|------------------------|-------|
| d4-1,2-Dichloroethane  | 108%  |
| d8-Toluene             | 101%  |
| Bromofluorobenzene     | 93.0% |
| d4-1,2-Dichlorobenzene | 105%  |

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

EPA SW-846 indicates that vinyl chloride and styrene may degrade in the presence of acid preservative.

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-11-1114  
SAMPLE

Lab Sample ID: ZL65A  
LIMS ID: 14-25426  
Matrix: Water  
Data Release Authorized: *MW*  
Reported: 12/04/14

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273  
Date Sampled: 11/20/14  
Date Received: 11/20/14

Date Extracted: 11/27/14  
Date Analyzed: 12/03/14 00:53  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-11-1114  
SAMPLE

Lab Sample ID: ZL65A  
LIMS ID: 14-25426  
Matrix: Water  
Date Analyzed: 12/03/14 00:53

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 63.2% | 2-Fluorobiphenyl       | 65.6% |
| d14-p-Terphenyl      | 68.8% | d4-1,2-Dichlorobenzene | 64.8% |
| d5-Phenol            | 68.5% | 2-Fluorophenol         | 65.6% |
| 2,4,6-Tribromophenol | 59.7% | d4-2-Chlorophenol      | 66.7% |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 2

Sample ID: LMW-6-1114  
SAMPLE

Lab Sample ID: ZL65B  
LIMS ID: 14-25427  
Matrix: Water  
Data Release Authorized: *mm*  
Reported: 12/04/14

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273  
Date Sampled: 11/20/14  
Date Received: 11/20/14

Date Extracted: 11/27/14  
Date Analyzed: 12/03/14 01:27  
Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL  
Final Extract Volume: 0.50 mL  
Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 2 of 2

Sample ID: LMW-6-1114  
SAMPLE

Lab Sample ID: ZL65B  
LIMS ID: 14-25427  
Matrix: Water  
Date Analyzed: 12/03/14 01:27

QC Report No: ZL65-Golder Associates  
Project: Landsburg  
9231000002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

Reported in µg/L (ppb)

#### Semivolatile Surrogate Recovery

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 60.8% | 2-Fluorobiphenyl       | 57.2% |
| d14-p-Terphenyl      | 64.4% | d4-1,2-Dichlorobenzene | 62.0% |
| d5-Phenol            | 71.5% | 2-Fluorophenol         | 68.5% |
| 2,4,6-Tribromophenol | 56.3% | d4-2-Chlorophenol      | 69.9% |

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Extraction Method: SW3510C

Page 1 of 1

Sample ID: LMW-11-1114

SAMPLE

Lab Sample ID: ZL65A

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: 

Reported: 12/17/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

Date Extracted: 11/27/14

Date Analyzed: 12/16/14 19:54

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

**Pest/PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 80.5% |
| Tetrachlorometaxylene | 88.0% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

Pesticides/PCB by GC/ECD Method SW8081B

Extraction Method: SW3510C

Page 1 of 1

Sample ID: LMW-6-1114

SAMPLE

Lab Sample ID: ZL65B

LIMS ID: 14-25427

Matrix: Water

Data Release Authorized: 

Reported: 12/17/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

Date Extracted: 11/27/14

Date Analyzed: 12/16/14 20:11

Instrument/Analyst: ECD6/YZ

GPC Cleanup: No

Sulfur Cleanup: No

Florisil Cleanup: No

Sample Amount: 500 mL

Final Extract Volume: 5.0 mL

Dilution Factor: 1.00

Silica Gel: No

| CAS Number | Analyte             | DL     | LOQ   | Result    |
|------------|---------------------|--------|-------|-----------|
| 319-84-6   | alpha-BHC           | 0.0085 | 0.050 | < 0.050 U |
| 319-85-7   | beta-BHC            | 0.0098 | 0.050 | < 0.050 U |
| 319-86-8   | delta-BHC           | 0.0087 | 0.050 | < 0.050 U |
| 58-89-9    | gamma-BHC (Lindane) | 0.016  | 0.050 | < 0.050 U |
| 76-44-8    | Heptachlor          | 0.011  | 0.050 | < 0.050 U |
| 309-00-2   | Aldrin              | 0.010  | 0.050 | < 0.050 U |
| 1024-57-3  | Heptachlor Epoxide  | 0.0079 | 0.050 | < 0.050 U |
| 959-98-8   | Endosulfan I        | 0.0089 | 0.050 | < 0.050 U |
| 60-57-1    | Dieldrin            | 0.017  | 0.10  | < 0.10 U  |
| 72-55-9    | 4,4'-DDE            | 0.018  | 0.10  | < 0.10 U  |
| 72-20-8    | Endrin              | 0.017  | 0.10  | < 0.10 U  |
| 33213-65-9 | Endosulfan II       | 0.014  | 0.10  | < 0.10 U  |
| 72-54-8    | 4,4'-DDD            | 0.019  | 0.10  | < 0.10 U  |
| 1031-07-8  | Endosulfan Sulfate  | 0.024  | 0.10  | < 0.10 U  |
| 50-29-3    | 4,4'-DDT            | 0.017  | 0.10  | < 0.10 U  |
| 72-43-5    | Methoxychlor        | 0.074  | 0.50  | < 0.50 U  |
| 53494-70-5 | Endrin Ketone       | 0.015  | 0.10  | < 0.10 U  |
| 7421-93-4  | Endrin Aldehyde     | 0.016  | 0.10  | < 0.10 U  |
| 5103-74-2  | trans-Chlordane     | 0.0082 | 0.050 | < 0.050 U |
| 5103-71-9  | cis-Chlordane       | 0.0082 | 0.050 | < 0.050 U |
| 8001-35-2  | Toxaphene           | 0.22   | 5.0   | < 5.0 U   |

Reported in µg/L (ppb)

## Pest/PCB Surrogate Recovery

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 86.0% |
| Tetrachlorometaxylene | 93.5% |

# This analyte (CAS registry No. 5103-74-2) is named trans-Chlordane in EPA Method 8081B(Feb 2007). It has also been named beta-Chlordane.

\$ This analyte (CAS registry No. 5103-71-9) is named cis-Chlordane in EPA Method 8081B(Feb 2007). It has also been named alpha-Chlordane.

## ORGANICS ANALYSIS DATA SHEET

NWTPH-HCID Method by GC/FID

Extraction Method: SW3510C

Page 1 of 1

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Matrix: Water

Data Release Authorized: *MMW*

Reported: 11/25/14

| ARI ID                | Sample ID                 | Extraction Date | Analysis Date | DL  | Range       | Result   |
|-----------------------|---------------------------|-----------------|---------------|-----|-------------|----------|
| MB-112214<br>14-25426 | Method Blank              | 11/22/14        | 11/24/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 92.2%    |
| ZL65A<br>14-25426     | LMW-11-1114<br>HC ID: --- | 11/22/14        | 11/24/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 94.0%    |
| ZL65B<br>14-25427     | LMW-6-1114<br>HC ID: ---  | 11/22/14        | 11/24/14      | 1.0 | Gas         | < 0.25 U |
|                       |                           |                 |               |     | Diesel      | < 0.50 U |
|                       |                           |                 |               |     | Oil         | < 0.50 U |
|                       |                           |                 |               |     | o-Terphenyl | 97.0%    |

Reported in mg/L (ppm)

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

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

Oil value based on the total peaks in the range from C24 to C38.

HC ID: DRO/RRO indicates results of organics or additional hydrocarbons in ranges are not identifiable.

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-11-1114  
SAMPLE

Lab Sample ID: ZL65A

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

Date Extracted: 11/27/14

Date Analyzed: 12/09/14 01:45

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: Yes

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 72.8% |
| Tetrachlorometaxylene | 52.8% |

ORGANICS ANALYSIS DATA SHEET  
PCB by GC/ECD Method SW8082A  
Extraction Method: SW3510C  
Page 1 of 1

Sample ID: LMW-6-1114  
SAMPLE

Lab Sample ID: ZL65B

LIMS ID: 14-25427

Matrix: Water

Data Release Authorized: 

Reported: 12/10/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

Date Extracted: 11/27/14

Date Analyzed: 12/09/14 02:05

Instrument/Analyst: ECD5/JGR

GPC Cleanup: No

Sulfur Cleanup: Yes

Sample Amount: 1000 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

Silica Gel: No

Acid Cleanup: Yes

| CAS Number | Analyte      | DL     | LOQ   | Result    |
|------------|--------------|--------|-------|-----------|
| 12674-11-2 | Aroclor 1016 | 0.0025 | 0.010 | < 0.010 U |
| 53469-21-9 | Aroclor 1242 | 0.0025 | 0.010 | < 0.010 U |
| 12672-29-6 | Aroclor 1248 | 0.0025 | 0.010 | < 0.010 U |
| 11097-69-1 | Aroclor 1254 | 0.0025 | 0.010 | < 0.010 U |
| 11096-82-5 | Aroclor 1260 | 0.0028 | 0.010 | < 0.010 U |
| 11104-28-2 | Aroclor 1221 | 0.0025 | 0.010 | < 0.010 U |
| 11141-16-5 | Aroclor 1232 | 0.0025 | 0.010 | < 0.010 U |

Reported in µg/L (ppb)

**PCB Surrogate Recovery**

|                       |       |
|-----------------------|-------|
| Decachlorobiphenyl    | 58.8% |
| Tetrachlorometaxylene | 48.2% |

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LMW-11-1114  
SAMPLE

Lab Sample ID: ZL65A

LIMS ID: 14-25426

Matrix: Water

Data Release Authorized: *[Signature]*

Reported: 12/19/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | <b>7440-38-2</b> | <b>Arsenic</b>   | 0.05 | 3     | <b>7</b>      |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>58,200</b> |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>1,410</b>  |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>26,300</b> |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>170</b>    |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>2,100</b>  |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>28,800</b> |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

## INORGANICS ANALYSIS DATA SHEET

## TOTAL METALS

Page 1 of 1

Sample ID: LMW-6-1114  
SAMPLE

Lab Sample ID: ZL65B

LIMS ID: 14-25427

Matrix: Water

Data Release Authorized: *8/2/14*

Reported: 12/19/14

QC Report No: ZL65-Golder Associates

Project: Landsburg

9231000002.R273

Date Sampled: 11/20/14

Date Received: 11/20/14

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number       | Analyte          | DL   | LOQ   | Result        | Q |
|-----------|-----------|-----------------|---------------|------------------|------------------|------|-------|---------------|---|
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7429-90-5        | Aluminum         | 7.6  | 1,000 | 1,000         | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-36-0        | Antimony         | 0.01 | 3     | 3             | U |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-38-2        | Arsenic          | 0.05 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-39-3        | Barium           | 1.3  | 500   | 500           | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-41-7        | Beryllium        | 0.16 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-43-9        | Cadmium          | 0.18 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-70-2</b> | <b>Calcium</b>   | 11.3 | 500   | <b>27,300</b> |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-47-3        | Chromium         | 1.2  | 1,000 | 1,000         | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-48-4        | Cobalt           | 0.3  | 10    | 10            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-50-8        | Copper           | 0.92 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-89-6</b> | <b>Iron</b>      | 7.5  | 200   | <b>2,410</b>  |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7439-92-1        | Lead             | 0.0  | 10    | 10            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-95-4</b> | <b>Magnesium</b> | 9.6  | 1,000 | <b>14,000</b> |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7439-96-5</b> | <b>Manganese</b> | 0.3  | 20    | <b>30</b>     |   |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-02-0        | Nickel           | 3.9  | 20    | 20            | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-09-7</b> | <b>Potassium</b> | 65.7 | 500   | <b>720</b>    |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7782-49-2        | Selenium         | 0.13 | 5     | 5             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-22-4        | Silver           | 0.43 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | <b>7440-23-5</b> | <b>Sodium</b>    | 11.4 | 500   | <b>7,070</b>  |   |
| 200.8     | 11/24/14  | 200.8           | 11/25/14      | 7440-28-0        | Thallium         | 0.00 | 2     | 2             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-62-2        | Vanadium         | 0.27 | 3     | 3             | U |
| 3010A     | 11/24/14  | 6010C           | 11/25/14      | 7440-66-6        | Zinc             | 1.4  | 20    | 20            | U |

Reported in ug/L (ppb).

U-Analyte undetected at given LOQ

LOQ-Reporting Limit

INORGANICS ANALYSIS DATA SHEET  
Total Mercury by Method SW7470A

Data Release Authorized:  
Reported: 12/01/14  
Date Received: 11/20/14  
Page 1 of 1

A handwritten signature in black ink, appearing to be 'JR' or similar, written over the 'Data Release Authorized' text.

QC Report No: ZL66-Golder Associates  
Project: Landsburg  
9231000002.R273

| Client/<br>ARI ID             | Date<br>Sampled | Matrix | Prep Date<br>Anal Date | RL   | Result |
|-------------------------------|-----------------|--------|------------------------|------|--------|
| LMW-11-1114<br>ZL66A 14-25429 | 11/20/14        | Water  | 11/24/14<br>11/28/14   | 20.0 | 20.0 U |
| LMW-6-1114<br>ZL66B 14-25430  | 11/20/14        | Water  | 11/24/14<br>11/28/14   | 20.0 | 20.0 U |
| MB-112414<br>Method Blank     | NA              | Water  | 11/24/14<br>11/28/14   | 20.0 | 20.0 U |

## Reported in ng/L

RL-Analytical reporting limit  
U-Undetected at reported detection limit

# Chain of Custody Record & Laboratory Analysis Request

|                                                      |                                           |                                                         |
|------------------------------------------------------|-------------------------------------------|---------------------------------------------------------|
| ARI Assigned Number:<br><b>2138-7565</b>             | Turn-around Requested:<br><b>Standard</b> | Page: <b>1</b> of <b>1</b>                              |
| ARI Client Company:<br><b>Goldner Associates</b>     | Phone:<br><b>425 883-0777</b>             | Date:<br><b>11/18/2014</b>                              |
| Client Contact:<br><b>Doug Morell, Jill Lamberts</b> | No. of Coolers:<br><b>7</b>               | Ice Present? <b>y</b><br>Cooler Temps: <b>0.4 - 4.7</b> |
| Client Project Name:<br><b>Landsburg</b>             |                                           |                                                         |
| Client Project #:<br><b>923-1000-002-R273</b>        | Samplers:<br><b>Rydecki, Lamberts</b>     |                                                         |



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants  
4611 South 134th Place, Suite 100  
Tukwila, WA 98168  
206-695-6200 206-695-6201 (fax)  
www.arilabs.com

| Sample ID         | Date     | Time | Matrix | No. Containers | Analysis Requested  |                               |            |                           |          |                        |                      |  | Notes/Comments |  |
|-------------------|----------|------|--------|----------------|---------------------|-------------------------------|------------|---------------------------|----------|------------------------|----------------------|--|----------------|--|
|                   |          |      |        |                | VOCs<br>Client list | PCBs (CLL)                    | Pesticides | SVOCs 8270<br>Client list | TPH-HCID | TAM L<br>-Total Metals | TAM L<br>Diss Metals |  |                |  |
| Trip Blank-111814 | 11/18/14 | -    | W      | 9-6            | X                   |                               |            |                           |          |                        |                      |  |                |  |
| LMW-10-1114       |          | 945  | W      | 17             | X                   | X                             | X          | X                         | X        | X                      | HOLD                 |  |                |  |
| LMW-2-1114        |          | 1140 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| LMW-4-1114        |          | 1310 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| Trip Blank-111914 | 11/19/14 | -    | W      | 6              | X                   | HOLD unless there are detects |            |                           |          |                        |                      |  |                |  |
| LMW-3-1114        |          | 0920 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      | HOLD                 |  |                |  |
| EB-1114           |          | 0900 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| LMW-8-1114        |          | 1005 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| LMW-5-1114        |          | 1130 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |
| LMW-9-1114        |          | 1315 | W      | 17             | X                   | X                             | X          | X                         | X        | X                      |                      |  |                |  |

|                                                                                                                                                                            |                                                    |                                                |                                 |                             |
|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----------------------------------------------------|------------------------------------------------|---------------------------------|-----------------------------|
| Comments/Special Instructions<br><b>-Ecology ETM EDD</b><br><br><b>** CLIENT SPECIFIC RLS and ANALYTE LIST **</b><br><br>PLS cc jlamberts@golder.com<br>dmorell@golder.com | Relinquished by:<br>(Signature) <i>[Signature]</i> | Received by:<br>(Signature) <i>[Signature]</i> | Relinquished by:<br>(Signature) | Received by:<br>(Signature) |
|                                                                                                                                                                            | Printed Name:<br><b>J. Lamberts</b>                | Printed Name:<br><b>Taylor Street</b>          | Printed Name:                   | Printed Name:               |
|                                                                                                                                                                            | Company:<br><b>Goldner</b>                         | Company:<br><b>ARL</b>                         | Company:                        | Company:                    |
|                                                                                                                                                                            | Date & Time:<br><b>11/19/14 1442</b>               | Date & Time:<br><b>11/19/14 1442</b>           | Date & Time:                    | Date & Time:                |

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

Validated by J. Lamberts 1/26/2015

2138-000002  
2569-000002

**ORGANICS ANALYSIS DATA SHEET****Semivolatiles by SW8270D GC/MS****Extraction Method: SW3520C**

Page 1 of 2

**Sample ID: LMW-4-1114****SAMPLE**

Lab Sample ID: ZS69A

LIMS ID: 15-518

Matrix: Water

Data Release Authorized:

Reported: 01/22/15

QC Report No: ZS69-Golder Associates

Project: Landsburg

923-1000-002.R273

Date Sampled: 11/18/14

Date Received: 11/19/14

Date Extracted: 01/15/15

Date Analyzed: 01/21/15 17:13

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                      | DL   | LOQ | Result  |
|------------|------------------------------|------|-----|---------|
| 108-95-2   | Phenol                       | 0.27 | 1.0 | < 1.0 U |
| 111-44-4   | Bis-(2-Chloroethyl) Ether    | 0.25 | 1.0 | < 1.0 U |
| 95-57-8    | 2-Chlorophenol               | 0.22 | 1.0 | < 1.0 U |
| 541-73-1   | 1,3-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 106-46-7   | 1,4-Dichlorobenzene          | 0.27 | 1.0 | < 1.0 U |
| 100-51-6   | Benzyl Alcohol               | 0.55 | 2.0 | < 2.0 U |
| 95-50-1    | 1,2-Dichlorobenzene          | 0.25 | 1.0 | < 1.0 U |
| 95-48-7    | 2-Methylphenol               | 0.21 | 1.0 | < 1.0 U |
| 108-60-1   | 2,2'-Oxybis(1-Chloropropane) | 0.24 | 1.0 | < 1.0 U |
| 106-44-5   | 4-Methylphenol               | 0.47 | 2.0 | < 2.0 U |
| 621-64-7   | N-Nitroso-Di-N-Propylamine   | 0.27 | 1.0 | < 1.0 U |
| 67-72-1    | Hexachloroethane             | 0.30 | 2.0 | < 2.0 U |
| 98-95-3    | Nitrobenzene                 | 0.25 | 1.0 | < 1.0 U |
| 78-59-1    | Isophorone                   | 0.42 | 1.0 | < 1.0 U |
| 88-75-5    | 2-Nitrophenol                | 0.26 | 3.0 | < 3.0 U |
| 105-67-9   | 2,4-Dimethylphenol           | 1.1  | 3.0 | < 3.0 U |
| 65-85-0    | Benzoic Acid                 | 3.9  | 20  | < 20 U  |
| 111-91-1   | bis(2-Chloroethoxy) Methane  | 0.24 | 1.0 | < 1.0 U |
| 120-83-2   | 2,4-Dichlorophenol           | 1.1  | 3.0 | < 3.0 U |
| 120-82-1   | 1,2,4-Trichlorobenzene       | 0.25 | 1.0 | < 1.0 U |
| 91-20-3    | Naphthalene                  | 0.25 | 1.0 | < 1.0 U |
| 106-47-8   | 4-Chloroaniline              | 1.7  | 5.0 | < 5.0 U |
| 87-68-3    | Hexachlorobutadiene          | 0.34 | 3.0 | < 3.0 U |
| 59-50-7    | 4-Chloro-3-methylphenol      | 1.1  | 3.0 | < 3.0 U |
| 91-57-6    | 2-Methylnaphthalene          | 0.30 | 1.0 | < 1.0 U |
| 77-47-4    | Hexachlorocyclopentadiene    | 1.1  | 5.0 | < 5.0 U |
| 88-06-2    | 2,4,6-Trichlorophenol        | 1.0  | 3.0 | < 3.0 U |
| 95-95-4    | 2,4,5-Trichlorophenol        | 1.1  | 5.0 | < 5.0 U |
| 91-58-7    | 2-Chloronaphthalene          | 0.25 | 1.0 | < 1.0 U |
| 88-74-4    | 2-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 131-11-3   | Dimethylphthalate            | 0.26 | 1.0 | < 1.0 U |
| 208-96-8   | Acenaphthylene               | 0.27 | 1.0 | < 1.0 U |
| 99-09-2    | 3-Nitroaniline               | 1.5  | 3.0 | < 3.0 U |
| 83-32-9    | Acenaphthene                 | 0.25 | 1.0 | < 1.0 U |
| 51-28-5    | 2,4-Dinitrophenol            | 3.4  | 20  | < 20 U  |
| 100-02-7   | 4-Nitrophenol                | 1.8  | 10  | < 10 U  |
| 132-64-9   | Dibenzofuran                 | 0.31 | 1.0 | < 1.0 U |
| 606-20-2   | 2,6-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |
| 121-14-2   | 2,4-Dinitrotoluene           | 1.1  | 3.0 | < 3.0 U |

**ORGANICS ANALYSIS DATA SHEET**  
**Semivolatiles by SW8270D GC/MS**  
**Extraction Method: SW3520C**  
 Page 2 of 2

**Sample ID: LMW-4-1114**  
**SAMPLE**

Lab Sample ID: ZS69A  
 LIMS ID: 15-518  
 Matrix: Water  
 Date Analyzed: 01/21/15 17:13

QC Report No: ZS69-Golder Associates  
 Project: Landsburg  
 923-1000-002.R273

| CAS Number | Analyte                    | DL   | LOQ | Result  |
|------------|----------------------------|------|-----|---------|
| 84-66-2    | Diethylphthalate           | 0.27 | 1.0 | < 1.0 U |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 0.27 | 1.0 | < 1.0 U |
| 86-73-7    | Fluorene                   | 0.29 | 1.0 | < 1.0 U |
| 100-01-6   | 4-Nitroaniline             | 2.0  | 3.0 | < 3.0 U |
| 534-52-1   | 4,6-Dinitro-2-Methylphenol | 3.6  | 10  | < 10 U  |
| 86-30-6    | N-Nitrosodiphenylamine     | 0.30 | 1.0 | < 1.0 U |
| 101-55-3   | 4-Bromophenyl-phenylether  | 0.24 | 1.0 | < 1.0 U |
| 118-74-1   | Hexachlorobenzene          | 0.28 | 1.0 | < 1.0 U |
| 87-86-5    | Pentachlorophenol          | 1.9  | 10  | < 10 U  |
| 85-01-8    | Phenanthrene               | 0.32 | 1.0 | < 1.0 U |
| 86-74-8    | Carbazole                  | 0.31 | 1.0 | < 1.0 U |
| 120-12-7   | Anthracene                 | 0.26 | 1.0 | < 1.0 U |
| 84-74-2    | Di-n-Butylphthalate        | 0.29 | 1.0 | < 1.0 U |
| 206-44-0   | Fluoranthene               | 0.30 | 1.0 | < 1.0 U |
| 129-00-0   | Pyrene                     | 0.28 | 1.0 | < 1.0 U |
| 85-68-7    | Butylbenzylphthalate       | 0.30 | 1.0 | < 1.0 U |
| 91-94-1    | 3,3'-Dichlorobenzidine     | 1.8  | 5.0 | < 5.0 U |
| 56-55-3    | Benzo(a)anthracene         | 0.29 | 1.0 | < 1.0 U |
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1  | 3.0 | < 3.0 U |
| 218-01-9   | Chrysene                   | 0.32 | 1.0 | < 1.0 U |
| 117-84-0   | Di-n-Octyl phthalate       | 0.27 | 1.0 | < 1.0 U |
| 205-99-2   | Benzo(b)fluoranthene       | 0.32 | 1.0 | < 1.0 U |
| 207-08-9   | Benzo(k)fluoranthene       | 0.34 | 1.0 | < 1.0 U |
| 50-32-8    | Benzo(a)pyrene             | 0.30 | 1.0 | < 1.0 U |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.36 | 1.0 | < 1.0 U |
| 53-70-3    | Dibenz(a,h)anthracene      | 0.39 | 1.0 | < 1.0 U |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.39 | 1.0 | < 1.0 U |
| 108-39-4   | 3&4-Methylphenol           | 0.80 | 2.0 | < 2.0 U |
| 90-12-0    | 1-Methylnaphthalene        | 0.26 | 1.0 | < 1.0 U |
| TOTBFA     | Total Benzofluoranthenes   | 0.80 | 2.0 | < 2.0 U |

UJ

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 62.0% | 2-Fluorobiphenyl       | 57.2% |
| d14-p-Terphenyl      | 67.6% | d4-1,2-Dichlorobenzene | 56.8% |
| d5-Phenol            | 83.5% | 2-Fluorophenol         | 77.9% |
| 2,4,6-Tribromophenol | 85.1% | d4-2-Chlorophenol      | 82.7% |

# Chain of Custody Record & Laboratory Analysis Request

|                                                                                                                                                                                   |  |                                                                                                                     |             |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                 |                |                                                                                                                                                                                                                                                                  |  |                             |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|--|---------------------------------------------------------------------------------------------------------------------|-------------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----------------|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|--|-----------------------------|--|--|--|----------------|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--------------------------------------------------|--|
| ARI Assigned Number:<br><b>2012</b>                                                                                                                                               |  | Turn-around Requested:<br><b>Standard</b>                                                                           |             | Page: <b>1</b> of <b>1</b>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      |                |  <b>Analytical Resources, Incorporated</b><br>Analytical Chemists and Consultants<br>4611 South 134th Place, Suite 100<br>Tukwila, WA 98168<br>206-695-6200 206-695-6201 (fax) |  |                             |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
| ARI Client Company:<br><b>Golder Associates</b>                                                                                                                                   |  | Phone:<br><b>425-883-0777</b>                                                                                       |             | Date:<br><b>2/12/15</b>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |                |                                                                                                                                                                                                                                                                  |  | Ice Present? <b>Y</b>       |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
| Client Contact:<br><b>Doug Morell, Jill Lamberts</b>                                                                                                                              |  |                                                                                                                     |             | No. of Coolers: <b>1</b>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                        |                |                                                                                                                                                                                                                                                                  |  | Cooler Temps: <b>5.7</b>    |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
| Client Project Name:<br><b>Landsburg</b>                                                                                                                                          |  |                                                                                                                     |             | Analysis Requested                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |                |                                                                                                                                                                                                                                                                  |  |                             |  |  |  | Notes/Comments |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
| Client Project #:<br><b>923-1000-002-R273</b>                                                                                                                                     |  | Samplers:<br><b>Lamberts, Hunt</b>                                                                                  |             | <div style="display: flex;"> <div style="writing-mode: vertical-rl; transform: rotate(180deg); border: 1px solid black; padding: 2px;">             BEHP by 8230 D           </div> <table border="1" style="width: 100%; height: 100%;"> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> <tr><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr> </table> </div> |                |                                                                                                                                                                                                                                                                  |  |                             |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  | Pls analyze under existing MSA btwn Golder + ARI |  |
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| Sample ID                                                                                                                                                                         |  | Date                                                                                                                | Time        | Matrix                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          | No. Containers |                                                                                                                                                                                                                                                                  |  |                             |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
| <b>LMW-4-0215</b>                                                                                                                                                                 |  | <b>2/12/15</b>                                                                                                      | <b>1450</b> | <b>W</b>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                        | <b>4</b>       |                                                                                                                                                                                                                                                                  |  |                             |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
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| Comments/Special Instructions<br><b>- Ecology Eim EDD</b><br><b>**Client specific RLS and analyze first**</b><br><b>Pls CC j.lamberts@golder.com</b><br><b>dmorell@golder.com</b> |  | Relinquished by:<br>(Signature)  |             | Received by:<br>(Signature)                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                 |                | Relinquished by:<br>(Signature)                                                                                                                                                                                                                                  |  | Received by:<br>(Signature) |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
|                                                                                                                                                                                   |  | Printed Name:<br><b>J. Lamberts</b>                                                                                 |             | Printed Name:<br><b>A. Volgardsen</b>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |                | Printed Name:                                                                                                                                                                                                                                                    |  | Printed Name:               |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
|                                                                                                                                                                                   |  | Company:<br><b>Golder</b>                                                                                           |             | Company:<br><b>ARI</b>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          |                | Company:                                                                                                                                                                                                                                                         |  | Company:                    |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |
|                                                                                                                                                                                   |  | Date & Time:<br><b>2/12/15 1607</b>                                                                                 |             | Date & Time:<br><b>2/12/15 1607</b>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             |                | Date & Time:                                                                                                                                                                                                                                                     |  | Date & Time:                |  |  |  |                |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |                                                  |  |

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the Invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

Validated by J. Lamberts 2/23/2015

2012: 000002

ORGANICS ANALYSIS DATA SHEET  
Semivolatiles by SW8270D GC/MS  
Extraction Method: SW3520C  
Page 1 of 1

Sample ID: LMW-4-0215  
SAMPLE

Lab Sample ID: ZW42A

LIMS ID: 15-2811

Matrix: Water

Data Release Authorized: *JB*

Reported: 02/18/15

QC Report No: ZW42-Golder Associates

Project: Landsburg

923-1000-002-R273

Date Sampled: 02/12/15

Date Received: 02/12/15

Date Extracted: 02/16/15

Date Analyzed: 02/18/15 12:04

Instrument/Analyst: NT6/JZ

Sample Amount: 500 mL

Final Extract Volume: 0.50 mL

Dilution Factor: 1.00

| CAS Number | Analyte                    | DL  | LOQ | Result  |
|------------|----------------------------|-----|-----|---------|
| 117-81-7   | bis(2-Ethylhexyl)phthalate | 2.1 | 3.0 | < 3.0 U |

Reported in µg/L (ppb)

**Semivolatile Surrogate Recovery**

|                      |       |                        |       |
|----------------------|-------|------------------------|-------|
| d5-Nitrobenzene      | 64.0% | 2-Fluorobiphenyl       | 64.0% |
| d14-p-Terphenyl      | 80.0% | d4-1,2-Dichlorobenzene | 59.6% |
| d5-Phenol            | 60.8% | 2-Fluorophenol         | 61.1% |
| 2,4,6-Tribromophenol | 69.6% | d4-2-Chlorophenol      | 65.3% |

**Date:** January 27, 2015  
**To:** Project File

**Project No.:** 923-1000. R209  
**Company:** Golder, Redmond

**From:** T. Stapp, Senior Chemist

**cc:** D. Morell, J. Lamberts

**Email:**

**RE: TECHNICAL REVIEW OF LABORATORY RAW DATA DELIVERABLE FOR SEMI-VOLATILE ANALYSIS; SDG #ZL38 & ZL39 / NOVEMBER, 2014 SAMPLE COLLECTION**

The following describes a summary review of analytical responses for semi-volatile analyses (EPA method 8270C) associated with well water samples collected at the Landsburg Mine site in Black Diamond, WA, on November 18, 2014 by Golder personnel. The data is reported by Analytical Resource Associates, Inc. associated with sample delivery group (SDG) numbers ZL38 and ZL39.

## 1.0 FINDINGS

SDG #ZL38 reports bis-2-ethylhexyl-phthalate (BEHP) at 12 ug/L (ppb), above the method reporting limit (MRL) of 3.0 ug/L for the semi-volatile compounds in sample LMW-4-1114. No other detectable semi-volatile compounds were reported above the method detection limit (MDLs) for any samples from the eight sample set submitted on December 4, 2014. Therefore, raw data was requested of the laboratory including extraction/ sequence records, mass response data, gas chromatographic/ mass spectrometry ion scans, and supporting quality control (QC) records to consider information associated with the BEHP detection.

### 1.1 REVIEW

A summary review was performed on the raw data to support the previous SDG deliverable and data validation report generated on the original data (Golder, 2015). Metrics for elements of data validation were found to meet compliance limits and exhibit acceptable instrument response including, continuing calibration, the calibration tune criteria, surrogate and internal standard recovery, blank spike recovery and method blanks. Consistency of sample handling and extraction was confirmed and all sample chromatograms were reviewed.

The November 28, 2014 BEHP detection noted for sample LMW-4-1114 is clearly visible on the associated chromatogram, however BEHP is also detected in significant response (above the baseline but below the MDL) for samples LMW-10-1114 and LMW-3-1114 in the same analytical set. Sample LMW-2-1114, although present in the original analytical run sequence (see below), was not represented with a chromatogram on this date due to a QC deficiency that required re-extraction and re-analysis on December 2, 2014. However, a BEHP response is also visible on the chromatogram from this date.

c:\1\stapp\1\ tms-pro\jd\m's\landsburg\2014\_memo\techmemo\_b2ehp cont\am.docx

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Since these samples bracket sample LMW-4-1114, there is evidence that a chronic condition of contamination exists for BEHP.

### ***1.1.1 Semi-volatile analytical sequence for November 28, 2014:***

- LMW-10-1114 data file #...b/11281413.D
- LMW-2-1114 data file #...b/11281414.D
- LMW-4-1114 data file #...b/11281415.D
- LMW-3-1114 data file #...b/11281416.D
- EB-1114 data file #...b/11281417.D
- LMW-8-1114 data file #...b/11281418.D
- LMW-9-1114 data file #...b/11281419.D
- LMW-5-1114 data file #...b/11281420.D

## **2.0 CLOSING**

Laboratory protocol precludes reporting BEHP detects below a certain response, and the sample report sheets therefore indicate a non-detect for all other samples in the client sample set. However, it is clear that the chromatograms show visible instrument response at the retention time expected for BEHP for nearly all of the samples in the sample set, including the method blank. It appears an instrument process error could be responsible for the BEHP contaminant found in all associated samples.

The chromatograms for each of the associated samples including the method blank in the client sample set, exhibit a signal response above the respective baselines but below the MDL for BEHP at the expected retention time (chromatograms attached). Therefore, although there is non-detect status for all other samples and QC blanks, there is evidence that an instrument process error has allowed the BEHP contaminant to be present. An instrument process error could include carry-over contamination from the injection needle, and/or column bleed. Although the laboratory is diligent in tracking and eliminating instrument process error, there is the possibility of preparation error from extraction and handling of the individual samples as well. Phthalates, as a chemical class, are ubiquitous in many products used to extract and handle samples and sample solvent extracts. Therefore, of the many target sources available to create contamination, it is difficult to pin-point the exact cause. Golder feels the BEHP contaminant that was detectable in sample LMW-4-1114, will likely disappear from subsequent sample collections for this well.

## **3.0 REFERENCE**

2015, Golder. Landsburg Mine Site Interim Groundwater Monitoring Report – November 2014. Recipient: Mr. Bill Kombol, Palmer Coking Coal Company, 31407 Highway 169, Black Diamond, WA 98010. January 26, 2015.

Data File: /chem2/nt6.i/20141128.b/11281402.d

Date : 28-NOV-2014 11:04

Client ID: ZK73MBW1

Sample Info: ZK73MBW1,

Volume Injected (uL): 1.0

Column phase: ZB-5msi

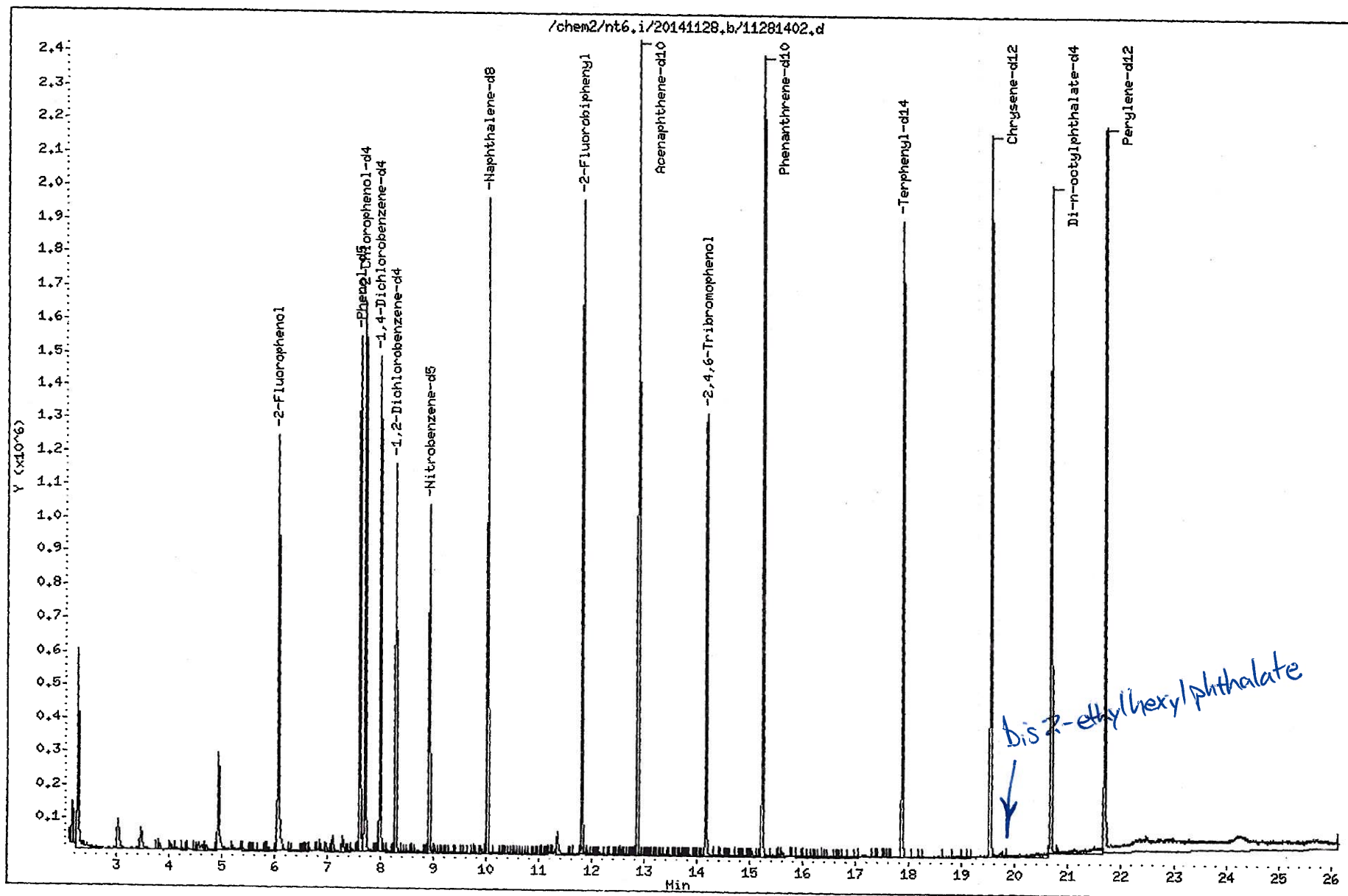
# METHOD BLANK

Instrument: nt6.i

Operator: JZ

Column diameter: 0.32

Page 6



Date : 28-NOV-2014 17:17

Client ID: LMW-10-1114

Sample Info: ZL38A

Volume Injected (uL): 1.0

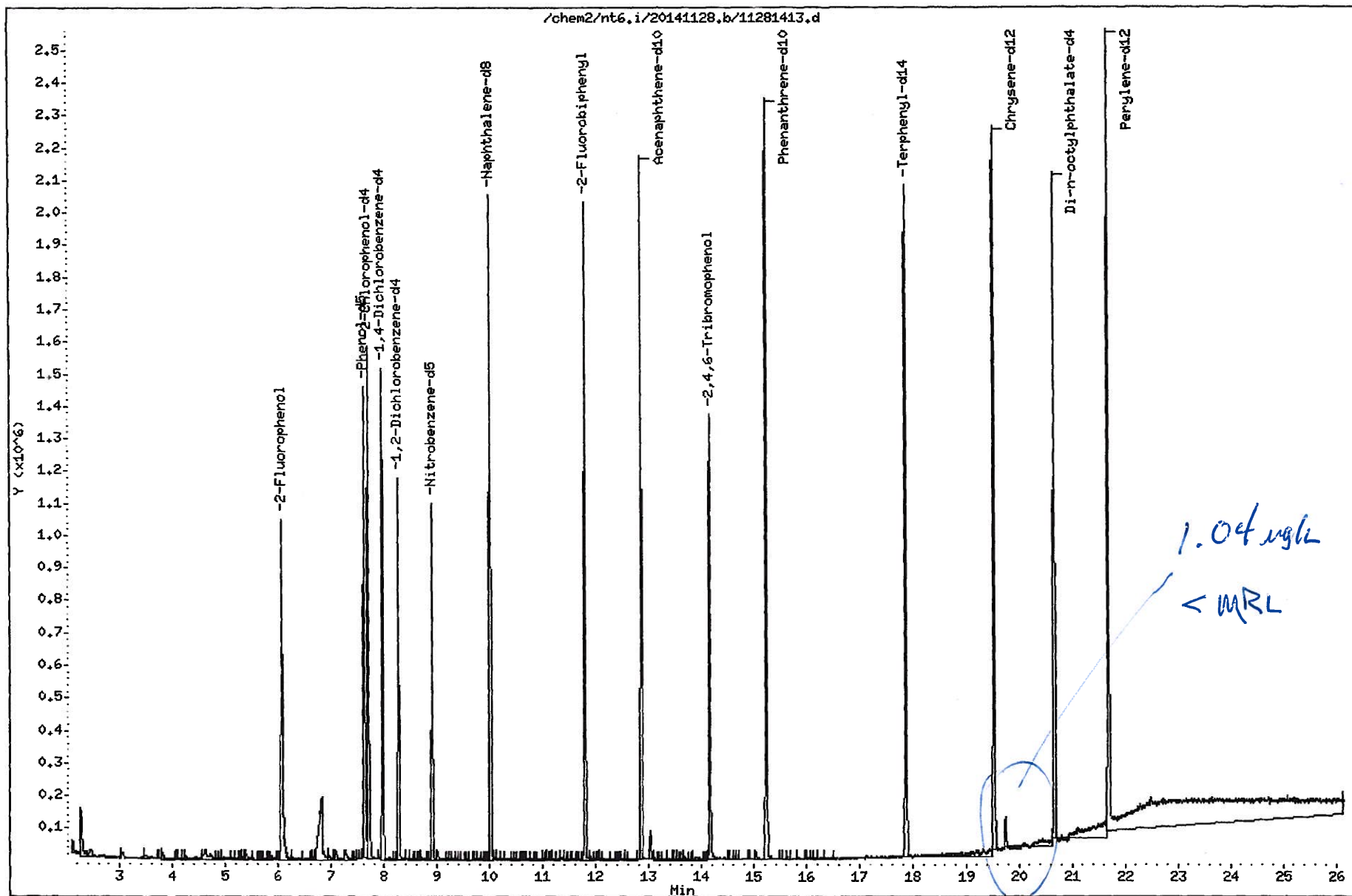
Column phase: ZB-5msi

Instrument: nt6.i

Operator: JZ

Column diameter: 0.32

LMW-10-1114



ND @ MRL

Date : 28-NOV-2014 18:25

Client ID: LMW-4-1114

Sample Info: ZL38C

Volume Injected (uL): 1.0

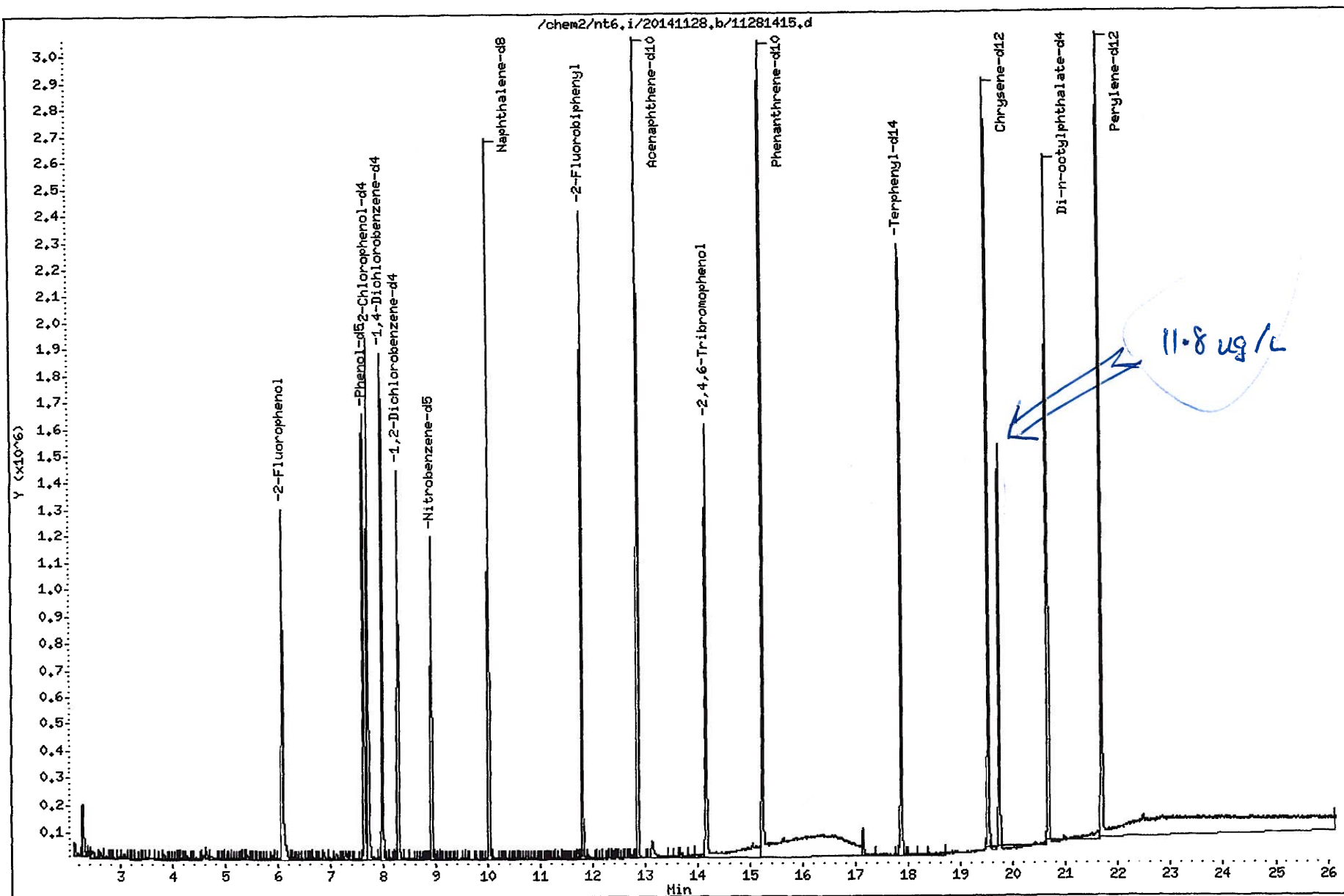
Column phase: ZB-5msi

Instrument: nt6.i

Operator: JZ

Column diameter: 0.32

LMW-4-1114



11.8 ug/L

Date : 28-NOV-2014 18:59

Client ID: LMW-3-1114

Sample Info: ZL38D

Volume Injected (uL): 1.0

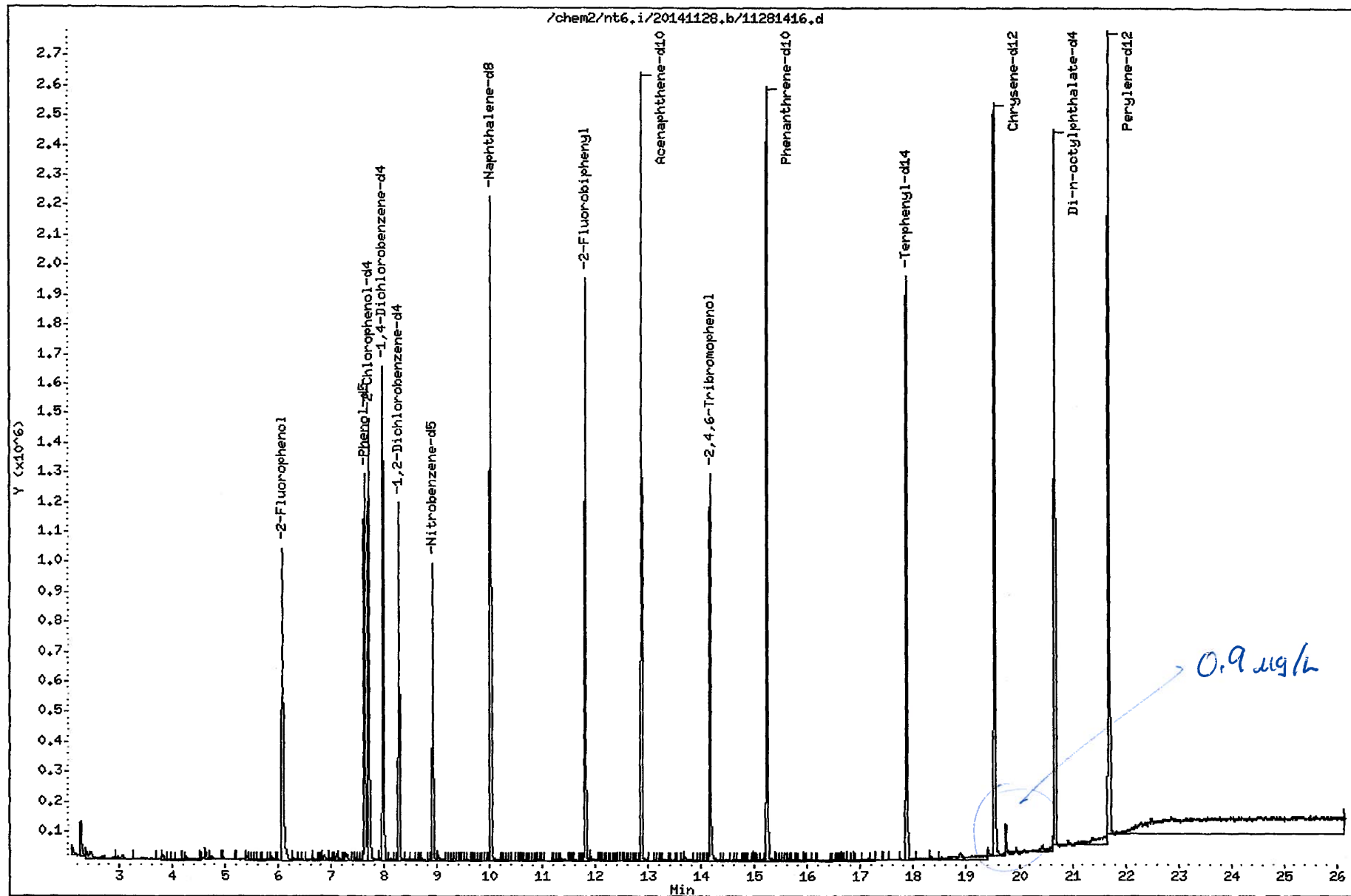
Column phase: ZB-5msi

Instrument: nt6.i

Operator: JZ

Column diameter: 0.32

LMW-3-1114



ND @ MRL

Data File: /chem2/nt6.i/20141128.b/11281417.d

Date : 28-NOV-2014 19:33

Client ID: EB-1114

Sample Info: ZL38E

Volume Injected (uL): 1.0

Column phase: ZB-5msi

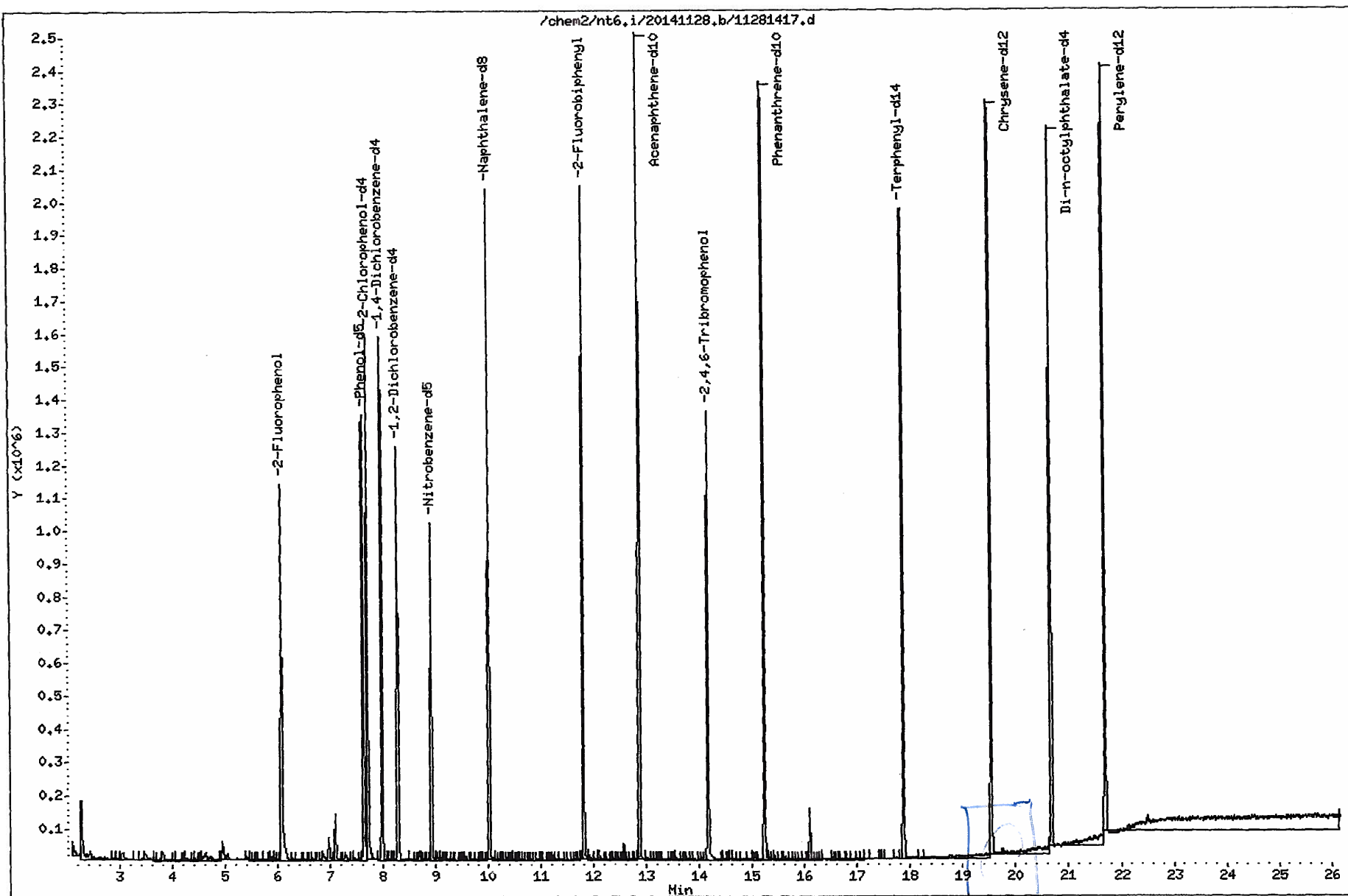
Instrument: nt6.i

Operator: JZ

Column diameter: 0.32

EQUIPMENT  
BLANK

Page 6



Date : 28-NOV-2014 20:06

Client ID: LMW-8-1114

Sample Info: ZL38F

Volume Injected (uL): 1.0

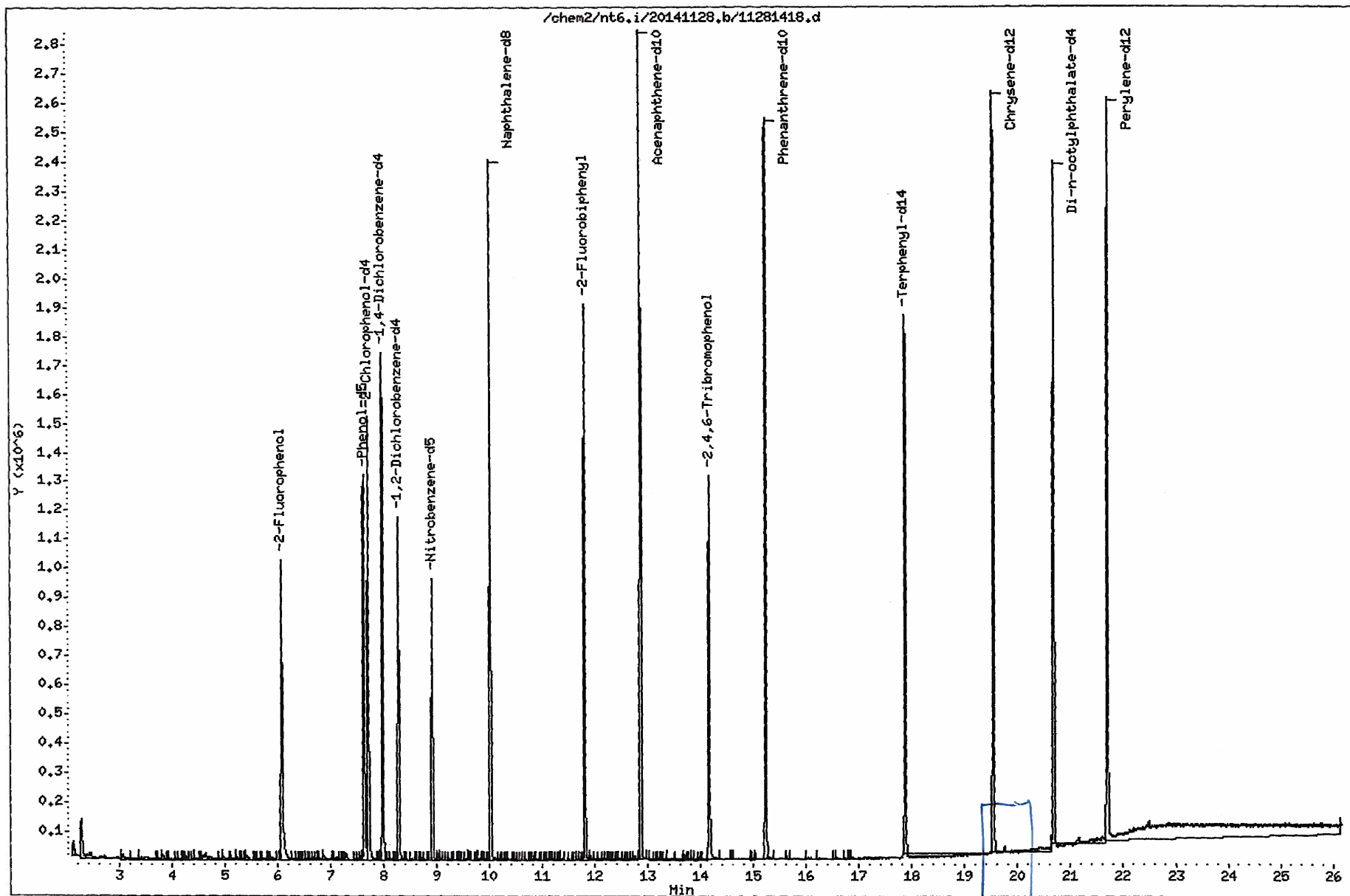
Column phase: ZB-5msi

Instrument: nt6.i

Operator: JZ

Column diameter: 0.32

LMW-8-1114



Date : 28-NOV-2014 21:14

Client ID: LMW-9-1114

Sample Info: ZL38H

Volume Injected (uL): 1.0

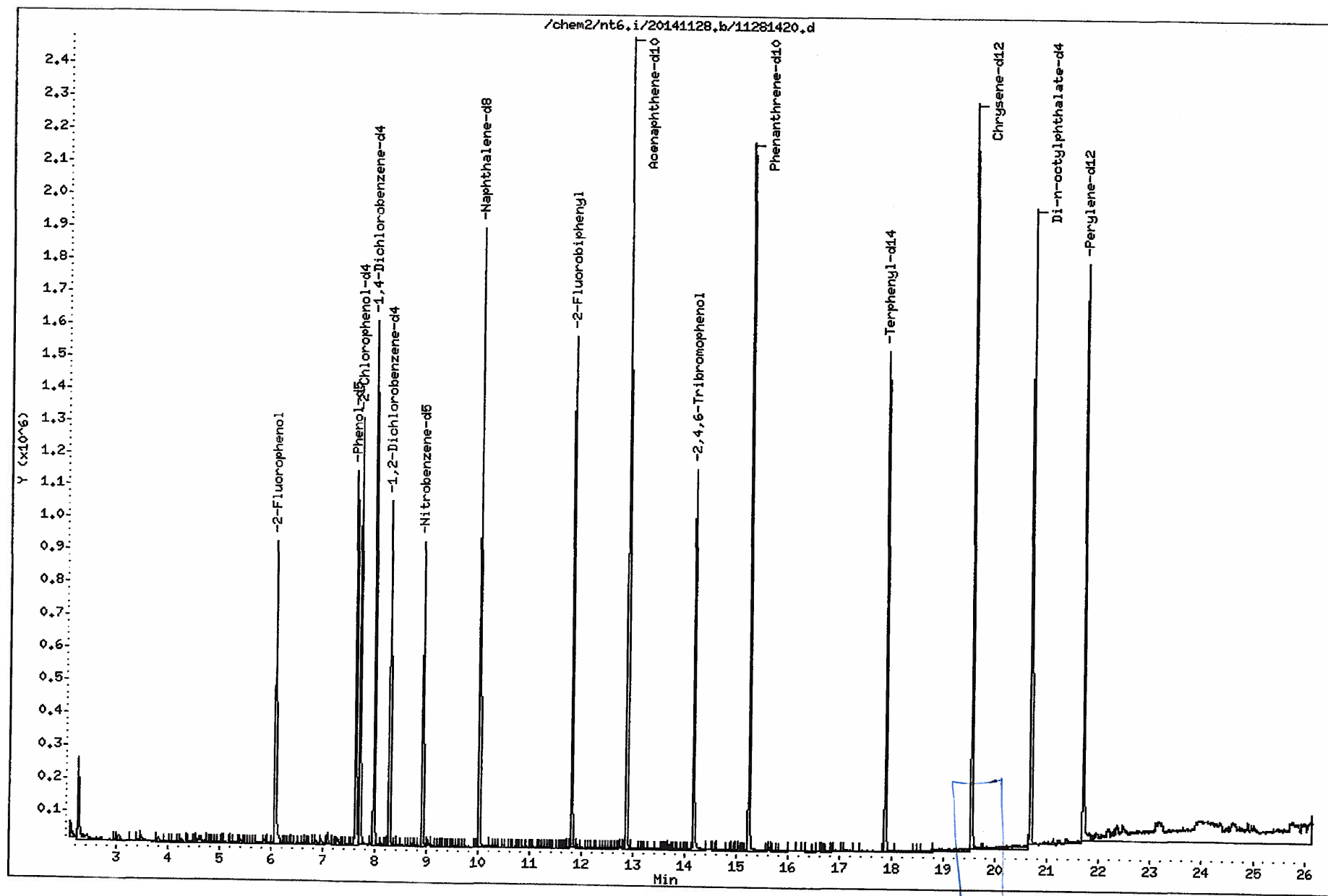
Column phase: ZB-5msi

Instrument: nt6.i

Operator: JZ

Column diameter: 0.32

LMW-9-1114



LMW-5-1114

