

| Appendix | Lab | Lab ID | Sample Date | Description | CAM Form Included | Lab Presumptive Certainty? | QC Performance Standards Met | CAM COMPLIANCE | ESM QAQC doc | Result? |
|----------|-------|----------|-------------|--------------------------|-------------------|----------------------------|------------------------------|-------------------|--------------|---|
| D | GWA | 84427 | 06/01/05 | Soil - Test Pits/PT spl | Yes | NO | No | CAM Non-Compliant | Completed | Data Not Compromised - Tank 1 data not utilized other than identification of tank contents. |
| D | GWA | 85632 | 07/13/05 | Soil - Wells | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| D | GWA | 87113 | 08/30/05 | Soil Grid Locations/TCLP | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| D | GWA | 87813 | 08/31/05 | Soil Grid Locations | Yes | Yes | Yes | CAM-Compliant | Completed | Data Not Compromised |
| D | GWA | 96205 | 06/21/06 | Asbestos | No | No | N/A | NON-CAM | Completed | Data Not Compromised |
| E | RC | 10061 | 03/23/06 | Packer Testing | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10068 | 03/23/06 | Packer Testing | Yes | Yes | Yes | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10070 | 03/24/06 | Packer Testing | Yes | Yes | Yes | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10079 | 03/27/06 | Packer Testing | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10088 | 03/28/06 | Packer Testing | Yes | Yes | Yes | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10094 | 03/29/06 | Packer Testing | Yes | Yes | Yes | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10103 | 03/30/06 | Packer Testing | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10105 | 03/31/06 | Packer Testing | Yes | Yes | Yes | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10154 | 04/06/06 | GW - wells | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10170 | 04/10/06 | GW - PZ-1 - PZ-3 | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 10171 | 04/10/06 | GW - wells | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| E | RC | 11371 | 11/15/06 | PZ-4 - PZ-7 | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| F | RC | 10426 | 05/30/06 | SW / Sed | Yes | Yes | No | CAM-Compliant | Completed | Data Not Compromised |
| G | ALPHA | L0604251 | 03/29/06 | Soil Gas | No | No | N/A | NON-CAM | Completed | Data Not Compromised |

ES&M LAB RESULTS QUALITY REVIEW

| | | | |
|--|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10061 |
| Job #: | 2004-301 | Sample Collection Date: | 3/23/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | No * | | |
| Holding time requirements met? | Yes | | |
| * Sample received at 9° C because it did not have time to cool between collection and lab receipt. | | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | N/A | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | N/A | |
| **RPD Calc: $100 * (\text{diff btwn sample \& dup}) / (\text{average of sample \& dup})$ | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| * Reporting Limits were high due to necessary dilutions. | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. The project narrative explained all non-conformances, see below: | | | |
| Describe Non-Conformances | | | |
| <ol style="list-style-type: none"> 1. VOC CCV: Vinyl Chloride showed high recovery. No impact to the data is suspected, as no Vinyl Chloride was found in the field samples. 2. A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 3. Samples: The quantitation limit for 4-isopropyltoluene due to possible carryover from a previous sample analysis. 4. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|---|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10068 |
| Job #: | 2004-301 | Sample Collection Date: | 3/23/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| N/A | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | N/A | |
| <small>**RPD Calc: 100*(diff btwn sample & dup)/(average of sample & dup)</small> | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| <small>* Reporting Limits were high due to necessary dilutions.</small> | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. | | | |
| Describe Non-Conformances | | | |
| <ol style="list-style-type: none"> 1. A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 2. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|---|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10070 |
| Job #: | 2004-301 | Sample Collection Date: | 3/24/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| N/A | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | N/A | |
| <small>**RPD Calc: 100*(diff btwn sample & dup)/(average of sample & dup)</small> | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| <small>* Reporting Limits were high due to necessary dilutions.</small> | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. | | | |
| Describe Non-Conformances | | | |
| 1. A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 2. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|--|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10079 |
| Job #: | 2004-301 | Sample Collection Date: | 3/27/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| N/A | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | N/A | |
| <small>**RPD Calc: 100*(diff btwn sample & dup)/(average of sample & dup)</small> | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| <small>* Reporting Limits were high due to necessary dilutions.</small> | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. The project narrative explained all non-conformances, see below: | | | |
| Describe Non-Conformances | | | |
| <ol style="list-style-type: none"> 1. The LCS showed a high recovery for Tetrachloroethene. This was most likely due to carryover from a previous sample analysis. The LCSD results were acceptable. 2. A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 3. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|---|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10088 |
| Job #: | 2004-301 | Sample Collection Date: | 3/28/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| N/A | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | N/A | |
| <small>**RPD Calc: 100*(diff btwn sample & dup)/(average of sample & dup)</small> | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| <small>* Reporting Limits were high due to necessary dilutions.</small> | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. | | | |
| Describe Non-Conformances | | | |
| 1. VOC: The following compounds were quantified with quadratic fit: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 2. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|---|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10094 |
| Job #: | 2004-301 | Sample Collection Date: | 3/29/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| N/A | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | N/A | |
| <small>**RPD Calc: 100*(diff btwn sample & dup)/(average of sample & dup)</small> | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| <small>* Reporting Limits were high due to necessary dilutions.</small> | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. | | | |
| Describe Non-Conformances | | | |
| 1. A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 2. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|--|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10103 |
| Job #: | 2004-301 | Sample Collection Date: | 3/27/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| N/A | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | N/A | |
| **RPD Calc: $100 * (\text{diff btwn sample \& dup}) / (\text{average of sample \& dup})$ | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| * Reporting Limits were high due to necessary dilutions. | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. The project narrative explained all non-conformances, see below: | | | |
| Describe Non-Conformances | | | |
| <ol style="list-style-type: none"> 1. The PRD between the LCS and the LCSD for Styrene was above the acceptance limit. The recoveries for both were acceptable. No data impact expected. 2. A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 3. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|--|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10105 |
| Job #: | 2004-301 | Sample Collection Date: | 3/31/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | No* | | |
| Holding time requirements met? | Yes | | |
| * Sample received at 12° C because it did not have time to cool between collection and lab receipt. | | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | N/A | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | N/A | |
| **RPD Calc: $100 * (\text{diff btwn sample \& dup}) / (\text{average of sample \& dup})$ | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | Yes | | |
| Action Limits | Yes | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. | | | |
| Describe Non-Conformances | | | |
| 1. A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|--|----------------------------|--------------------------------|---------------------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10154 |
| Job #: | 2006-056 | Sample Collection Date: | 04/06/06 - 04/07/06 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| DUP (VPH) and MS/MSD included with samples | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | Yes | |
| <small>**RPD Calc: 100*(diff btwn sample & dup)/(average of sample & dup)</small> | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| <small>* Reporting Limits were high due to necessary dilutions.</small> | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. The project narrative explained all non-conformances, see below: | | | |
| Describe Non-Conformances | | | |
| <ol style="list-style-type: none"> 1. EPH: The following samples had surrogates that did not meet the acceptance criteria: 10154-05, 07, 08, and 09. All other batch QC has acceptable recoveries. It is suspected that this may have been caused by matrix interference. The samples could not be re-extracted due to insufficient sample volume. 2. VOC 8260: LCS 10154-53 did not meet acceptance criteria for dichlorodifluoromethane. LCS 10154-51 did not meet the acceptance limits for 2,2-dichloropropane, and dichlorodifluoromethane. The RPD for styrene did not meet the acceptance criteria, however both recoveries were acceptable. LCS 10154-57 did not meet the acceptance limits for 2,2-dichloropropane, and dichlorodifluoromethane. The RPD for styrene did not meet the acceptance criteria due to a low recovery in the LCSD. These compounds noted with failures are known to be problematic in the method. 3. EPH: The fractionation check sample (LCS) for the batch of silica gel in use for these samples met the method acceptance criteria. 4. VPH Target compounds and ranges were determined by GC/MS. Ranges were determined in a similar manner as described in the MassDEP APH method of 2/2000. 5. PCB: The sample peaks most closely resembled Aroclor 1242, however there was also a similarity to Aroclor 1248. Possibly due to weathering, the Aroclor identification is difficult and not absolute. Chromatograms are provided. Samples that contained non-aroclor peaks have been noted on the report pages. Quantification is quadratic. 6. VOC: The following compounds were quantified with quadratic fit: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 7. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|---|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10170 |
| Job #: | 2006-056 | Sample Collection Date: | 4/10/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | | |
| **RPD Calc: $100 * (\text{diff btwn sample \& dup}) / (\text{average of sample \& dup})$ | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | No* | | |
| Action Limits | Yes | | |
| * Reporting Limits were high due to necessary dilutions. | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. The project narrative explained all non-conformances, see below: | | | |
| Describe Non-Conformances | | | |
| <ol style="list-style-type: none"> 1. VOC: The continuing calibration verification for naphthene did not meet acceptance criteria (69%). Samples 10170-01, 03, & 04 were analyzed in this window. All calibration check compounds (CCC) met acceptance criteria. No further action required. 2. VOC 8260: LCS 10170-51 did not meet acceptance criteria for dichlorodifluoromethane and acetone. LCS 10170-53 did not meet the acceptance limits for 2,2-dichloropropane, bromomethane, styrene, and dichlorodifluoromethane. The RPD for styrene did not meet the acceptance criteria. These compounds noted with failures are known to be problematic in the method. 3. PCB: The sample peaks most closely resembled Aroclor 1242, however there was also a similarity to Aroclor 1248. Possibly due to weathering, the Aroclor identification is difficult and not absolute. Chromatograms are provided. Samples that contained non-aroclor peaks have been noted on the report pages. Quantification is quadratic. 4. PCB: A sulfur clean up was performed on samples 10170-01 and -05. 5. VOC: The following compounds were quantified with quadratic fit: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 6. Dilutions performed during the analysis are noted on the result pages. 7. Lead only was requested by the customer. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|---|----------------------------|--------------------------------|-----------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 10171 |
| Job #: | 2006-056 | Sample Collection Date: | 4/10/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | <u>Yes</u> | | |
| Preservation requirements met? | <u>Yes</u> | | |
| Holding time requirements met? | <u>Yes</u> | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| Field Dup(s) vs. Environmental Sample: | | RPD <51%? | |
| **RPD Calc: 100*(diff btwn sample & dup)/(average of sample & dup) | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | <u>Yes</u> | | |
| Reporting Limits | <u>No*</u> | | |
| Action Limits | <u>Yes</u> | | |
| * Reporting Limits were high due to necessary dilutions. | | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. The project narrative explained all non-conformances, see below: | | | |
| Describe Non-Conformances | | | |
| <ol style="list-style-type: none"> 1. VOC: The continuing calibration verification for naphthene did not meet acceptance criteria (69%). Samples 10171-02, 03, 04, 06, 07, & 09 were analyzed in this window. All calibration check compounds (CCC) met acceptance criteria. No further action required. 2. EPH: The following samples had surrogates that did not meet the acceptance criteria: 10171-09. All other batch QC has acceptable recoveries. It is suspected that this may have been caused by matrix interference. The samples could not be re-extracted due to insufficient sample volume. 3. VOC 8260: LCS 10171-51 did not meet acceptance criteria for dichlorodifluoromethane and acetone. LCS 10171-53 did not meet the acceptance limits for 2,2-dichloropropane, bromomethane, styrene, and dichlorodifluoromethane. The RPD for styrene did not meet the acceptance criteria. These compounds noted with failures are known to be problematic in the method. 4. EPH: The fractionation check sample (LCS) for the batch of silica gel in use for these samples met the method acceptance criteria. 5. VPH Target compounds and ranges were determined by GC/MS. Ranges were determined in a similar manner as described in the MassDEP APH method of 2/2000. 6. PCB: The sample peaks most closely resembled Aroclor 1242, however there was also a similarity to Aroclor 1248. Possibly due to weathering, the Aroclor identification is difficult and not absolute. Chromatograms are provided. Samples that contained non-aroclor peaks have been noted on the report pages. Quantification is quadratic. 7. VOC: The following compounds were quantified with quadratic fit: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene. 8. Dilutions performed during the analysis are noted on the result pages. 9. RCRA metals only were requested by the customer | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

ES&M LAB RESULTS QUALITY REVIEW

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|--|----------------------------|--------------------------------|------------|
| Site: | DND Lewis Chemical | | |
| Lab: | Resource Laboratories, LLC | Lab ID: | 11371 |
| Job #: | 2006-056 | Sample Collection Date: | 11/15/2006 |
| Were sampling and analytical methods requirements met? | | | |
| Correct containers used? | Yes | | |
| Preservation requirements met? | Yes | | |
| Holding time requirements met? | Yes | | |
| Correct # of dupes, matrix spikes and matrix spike dupes, trip blanks (based on number of samples)? | | | |
| N/A | | | |
| Field Dup(s) vs. Environmental Sample: | RPD <51%? | | |
| <div style="display: flex; justify-content: space-between;"> **RPD Calc: $100 * (\text{diff btwn sample \& dup}) / (\text{average of sample \& dup})$ <div style="border-bottom: 1px solid black; width: 40%;"></div> </div> | | | |
| Were the following analytical precision and accuracy requirements met? | | | |
| Detection Limits | Yes | | |
| Reporting Limits | Yes | | |
| Action Limits | Yes | | |
| Review lab QC reports and project narrative. | | | |
| Lab Data Certification indicates that presumptive certainty has been met. The project narrative explained all non-conformances, see below: | | | |
| Describe Non-Conformances | | | |
| <ol style="list-style-type: none"> 1. VOC: LCS 0403790 did not meet acceptance criteria for bromoform and 1,2 dibromo-3-chloropropane. LCSD 0403790 did not meet acceptance criteria for bromoform and 1,2 dibromo-3-chloropropane. These compounds are known to be problematic in the method. 2. Dilutions performed during the analysis are noted on the result pages. | | | |
| Observations? | | | |
| All notes were reviewed and do not indicate compromised data. | | | |

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10061
Date Received: 3/23/06

Project: 2004-301 DND Lewis

Attached please find results for the analysis of the samples received on the date referenced above.

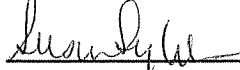
This is a reissue of report 10061, including the associated batch QC, case narrative and MCP Report Certification Form.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

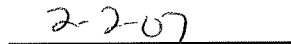
Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager



Date

Total number of pages



Resource Laboratories, LLC Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Lab Number: 10061-01
Sample Designation: ESM-03B (21-30)
Date Sampled: 3/23/06
Date Analyzed: 3/23/06
Matrix: Water
Instrument Dilution Factor: 20
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 40 | 1,1,2-trichloroethane | U | 40 |
| chloromethane | U | 40 | 1,3-dichloropropane | U | 40 |
| vinyl chloride | U | 40 | tetrachloroethene | 8700 | 40 |
| bromomethane | U | 40 | dibromochloromethane | U | 40 |
| chloroethane | U | 40 | 1,2-dibromoethane | U | 40 |
| trichlorofluoromethane | U | 40 | chlorobenzene | U | 40 |
| diethyl ether | U | 200 | 1,1,1,2-tetrachloroethane | U | 40 |
| acetone | U | 200 | ethylbenzene | 41 | 40 |
| 1,1-dichloroethene | 160 | 20 | m&p-xylenes | 93 | 40 |
| methylene chloride | U | 100 | o-xylene | U | 40 |
| carbon disulfide | U | 40 | styrene | U | 40 |
| methyl t-butyl ether (MTBE) | U | 40 | bromoform | U | 40 |
| trans-1,2-dichloroethene | U | 40 | isopropylbenzene | U | 40 |
| 1,1-dichloroethane | U | 40 | 1,1,2,2-tetrachloroethane | U | 40 |
| 2-butanone (MEK) | U | 200 | 1,2,3-trichloropropane | U | 40 |
| 2,2-dichloropropane | U | 40 | n-propylbenzene | U | 40 |
| cis-1,2-dichloroethene | 370 | 40 | bromobenzene | U | 40 |
| chloroform | U | 40 | 1,3,5-trimethylbenzene | U | 40 |
| bromochloromethane | U | 40 | 2-chlorotoluene | U | 40 |
| tetrahydrofuran (THF) | U | 200 | 4-chlorotoluene | U | 40 |
| 1,1,1-trichloroethane | 2500 | 40 | tert-butylbenzene | U | 40 |
| 1,1-dichloropropene | U | 40 | 1,2,4-trimethylbenzene | 60 | 40 |
| carbon tetrachloride | U | 40 | sec-butylbenzene | U | 40 |
| 1,2-dichloroethane | U | 40 | 1,3-dichlorobenzene | U | 40 |
| benzene | U | 40 | 4-isopropyltoluene | U | 100 |
| trichloroethene | 4200 | 40 | 1,4-dichlorobenzene | U | 40 |
| 1,2-dichloropropane | U | 40 | 1,2-dichlorobenzene | 240 | 40 |
| bromodichloromethane | U | 40 | n-butylbenzene | U | 40 |
| dibromomethane | U | 40 | 1,2-dibromo-3-chloropropane | U | 40 |
| 4-methyl-2-pentanone (MIBK) | U | 200 | 1,2,4-trichlorobenzene | U | 40 |
| cis-1,3-dichloropropene | U | 40 | hexachlorobutadiene | U | 40 |
| toluene | 180 | 40 | naphthalene | U | 100 |
| trans-1,3-dichloropropene | U | 40 | 1,2,3-trichlorobenzene | U | 40 |
| 2-hexanone | U | 200 | | | |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 98 | 78-114 | | | |
| toluene-D8 | 103 | 88-110 | | | |
| 4-bromofluorobenzene | 96 | 86-115 | | | |

U = Below quantitation limit

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10061

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 9 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

No exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

VOC CCV: Vinyl Chloride showed high recovery. No impact to the data is suspected, as no Vinyl Chloride was found in the field samples.

A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene.

Samples: The quantitation limit for 4-isopropyltoluene was raised due to possible carryover from a previous sample analysis.

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10061

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|------------|----------|---------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |

| | | |
|---|---|---------------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|---|---------------------|
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|--|---------------------|
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
|---|--|---------------------|

| | | |
|---|---|----------------------|
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No () NA |
|---|---|----------------------|

| | | |
|---|---|---------------------|
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes () No (x) |
|---|---|---------------------|

| | | |
|---|--|---------------------|
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes (x) No () |
|---|--|---------------------|

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Position: Lab DirectorPrinted Name: Susan C. SylvesterDate: 2-2-07

Lab Number: 10061-50
Sample Designation: Method Blank
Date Sampled: N/A
Date Analyzed: 3/23/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| dibromofluoromethane | 102 | 78-114 |
| toluene-D8 | 108 | 88-110 |
| 4-bromofluorobenzene | 98 | 86-115 |

U = Below quantitation limit

Lab Number: 10061-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\VOA03\MAR06\032306\W3032322.D
Date Analyzed: 3/23/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-------------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 17 | 86% | 17 | 83% | 3% |
| chloromethane | 20 | 98% | 18 | 91% | 8% |
| vinyl chloride | 23 | 115% | 21 | 106% | 8% |
| bromomethane | 16 | 79% | 16 | 82% | 4% |
| chloroethane | 20 | 100% | 19 | 97% | 3% |
| trichlorofluoromethane | 21 | 104% | 20 | 99% | 5% |
| diethylether | 21 | 103% | 21 | 104% | 1% |
| acetone | 17 | 83% | 16 | 81% | 3% |
| 1,1-dichloroethene | 21 | 105% | 20 | 98% | 7% |
| methylene chloride | 22 | 108% | 21 | 104% | 4% |
| carbon disulfide | 20 | 99% | 18 | 92% | 7% |
| methyl-t-butyl ether (MTBE) | 41 | 103% | 40 | 101% | 2% |
| trans-1,2-dichloroethene | 21 | 105% | 20 | 101% | 4% |
| isopropyl ether (DIPE) | 21 | 106% | 21 | 103% | 3% |
| ethyl-t-butyl ether (ETBE) | 20 | 100% | 20 | 102% | 2% |
| 1,1-dichloroethane | 20 | 101% | 20 | 100% | 1% |
| t-butanol (TBA) | 89 | 89% | 89 | 89% | 1% |
| 2-butanone (MEK) | 18 | 92% | 18 | 88% | 4% |
| 2,2-dichloropropane | 21 | 104% | 19 | 96% | 8% |
| cis-1,2-dichloroethene | 22 | 112% | 21 | 106% | 5% |
| chloroform | 22 | 109% | 21 | 107% | 2% |
| bromochloromethane | 22 | 109% | 22 | 108% | 2% |
| tetrahydrofuran (THF) | 19 | 96% | 19 | 96% | 0% |
| 1,1,1-trichloroethane | 21 | 103% | 20 | 102% | 1% |
| 1,1-dichloropropene | 21 | 106% | 21 | 103% | 3% |
| t-amyl-methyl ether (TAME) | 20 | 101% | 20 | 99% | 2% |
| carbon tetrachloride | 18 | 88% | 17 | 87% | 1% |
| 1,2-dichloroethane | 21 | 105% | 20 | 100% | 4% |
| benzene | 22 | 108% | 20 | 102% | 6% |
| trichloroethene | 21 | 103% | 21 | 104% | 1% |
| 1,2-dichloropropane | 21 | 103% | 20 | 100% | 3% |
| bromodichloromethane | 18 | 92% | 18 | 92% | 0% |
| dibromomethane | 22 | 110% | 22 | 108% | 2% |
| 4-methyl-2-pentanone (MIBK) | 19 | 95% | 19 | 96% | 1% |
| cis-1,3-dichloropropene | 20 | 100% | 20 | 99% | 1% |
| toluene | 22 | 108% | 20 | 101% | 7% |
| trans-1,3-dichloropropene | 18 | 92% | 18 | 92% | 0% |
| 2-hexanone | 18 | 91% | 17 | 87% | 4% |
| 1,1,2-trichloroethane | 21 | 106% | 21 | 105% | 1% |
| 1,3-dichloropropane | 21 | 106% | 21 | 103% | 3% |
| tetrachloroethene | 21 | 107% | 20 | 102% | 4% |
| dibromochloromethane | 19 | 93% | 19 | 94% | 0% |
| 1,2-dibromoethane (EDB) | 21 | 104% | 20 | 102% | 2% |
| chlorobenzene | 21 | 104% | 20 | 100% | 4% |
| 1,1,1,2-tetrachloroethane | 20 | 100% | 19 | 94% | 6% |
| ethylbenzene | 22 | 108% | 20 | 99% | 9% |
| m&p-xylenes | 44 | 109% | 41 | 103% | 6% |
| o-xylene | 22 | 108% | 21 | 103% | 5% |
| styrene | 21 | 106% | 19 | 97% | 9% |
| bromoform | 18 | 89% | 18 | 91% | 2% |
| isopropylbenzene | 23 | 113% | 22 | 108% | 5% |
| 1,1,2,2-tetrachloroethane | 19 | 95% | 18 | 92% | 3% |
| 1,2,3-trichloropropane | 19 | 94% | 18 | 91% | 3% |
| n-propylbenzene | 21 | 106% | 21 | 103% | 3% |
| bromobenzene | 21 | 104% | 20 | 102% | 3% |
| 1,3,5-trimethylbenzene | 21 | 104% | 20 | 98% | 6% |
| 2-chlorotoluene | 21 | 105% | 20 | 98% | 7% |
| 4-chlorotoluene | 20 | 102% | 19 | 96% | 6% |
| tert-butylbenzene | 19 | 97% | 19 | 95% | 2% |
| 1,2,4-trimethylbenzene | 21 | 106% | 20 | 102% | 4% |
| sec-butylbenzene | 21 | 103% | 20 | 99% | 4% |
| 1,3-dichlorobenzene | 20 | 101% | 20 | 98% | 3% |
| 4-isopropyltoluene | 22 | 110% | 22 | 109% | 1% |
| 1,4-dichlorobenzene | 19 | 97% | 19 | 96% | 1% |
| 1,2-dichlorobenzene | 21 | 104% | 20 | 99% | 6% |
| n-butylbenzene | 22 | 109% | 21 | 104% | 5% |
| 1,2-dibromo-3-chloropropane (| 19 | 93% | 19 | 93% | 0% |
| 1,2,4-trichlorobenzene | 18 | 92% | 18 | 91% | 1% |
| hexachlorobutadiene | 20 | 98% | 19 | 96% | 2% |
| naphthalene | 16 | 78% | 16 | 78% | 0% |
| 1,2,3-trichlorobenzene | 19 | 93% | 19 | 94% | 2% |
| 1,4-dioxane | 39 | 97% | 35 | 87% | 11% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 102% | 100% |
| SS toluene-D8 | 106% | 100% |
| SS 4-bromofluorobenzene | 109% | 105% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

Resource Laboratories, LLC
124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

**CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST**

ANALYSIS REQUEST

10061

Company Name:

Phone #: 508-285-9700

Company Address:

FAX #: 508-285-9957

Project Manager:

Site Location (City, State):

Invoice To:

Project ID / Name:

Protocol:

RCRA SDWA NPDES

NDDES OTHER

Lab Sample ID

Field ID

CONTAINERS

Matrix

Preservation Method

Sampling

SAMPLER

(Lab Use Only)

WATER
SOLID
OTHER

HCl
HNO₃
H₂SO₄
NaOH
MeOH
OTHER (Specify)

DATE

TIME

3/23/06 6:10 AM

☐ VOC 8260-NH List ☐ MADEP VPH ☐ MEGRO

☒ VOC 8260 ☐ VOC8015GRO ☐ VOC 624

☐ VOC 8260 BTEX, MIBE, Naphthalene only

☐ VOC 524.2 ☐ VOC 524.2 NH List

☐ TPH Fingerprint ☐ MEDRO ☐ DRO 8015 ☐ EPH

☐ 8270PAH ☐ 8270ABN ☐ 625

☐ 8082 PCB ☐ 8081 Pesticides ☐ 608

☐ O&G 1664 ☐ O&G SM5520F

☐ pH ☐ BOD ☐ Conductivity

☐ TSS ☐ TDS ☐ TS

☐ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals

☐ Total Metals-List ☐ Dissolved Metals-List

☐ Ammonia ☐ COD

☐ T-Phosphate ☐ Phenol

☐ Cyanide ☐ Sulfide

☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride

☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitibility/FP

☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC

☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)

☐ Standard Drinking Water Test ☐ Bacteria P/A

Grab (G) or Composite (C)

TAT REQUESTED

SPECIAL INSTRUCTIONS

Priority (24 hr) ☒

Expedited (48 hr) ☐

10 Business Days ☐

Other ☐

E-Mail Address

REPORTING INSTRUCTIONS

☐ FAX ☐ OTHER (specify)

☐ PDF ☐ Excel Spreadsheet

RECEIVED ON ICE ☒ YES ☐ NO

TEMPERATURE

9 °C

Lab Use Only

Quote #

PO #

Relinquished by Sampler:

Relinquished by:

Date

Time

Received by:

Date

Time

Received by Laboratory:

Date

Time

Way Bill#:

Date

Time

Date

Time

Date

Time

Date

Time

Date

Time

**CUSTODY
RECORD**

Sample per instructions from Joe Callahan / ASAP turnaround

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10068
Date Received: 3/23/06

Project: 2004-301 DND Lewis

Attached please find results for the analysis of the samples received on the date referenced above.

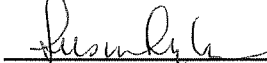
This is a reissue of report 10061, including the associated batch QC, case narrative and MCP Report Certification Form.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

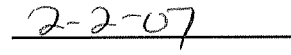
Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager



Date

Total number of pages



Resource Laboratories, LLC Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Lab Number: 10068-01
Sample Designation: ESM-03B (30-40)
Date Sampled: 3/23/06
Date Analyzed: 3/24/06
Matrix: Water
Instrument Dilution Factor: 20
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 40 | 1,1,2-trichloroethane | U | 40 |
| chloromethane | U | 40 | 1,3-dichloropropane | U | 40 |
| vinyl chloride | U | 40 | tetrachloroethene | 6600 | 40 |
| bromomethane | U | 40 | dibromochloromethane | U | 40 |
| chloroethane | U | 40 | 1,2-dibromoethane | U | 40 |
| trichlorofluoromethane | U | 40 | chlorobenzene | U | 40 |
| diethyl ether | U | 200 | 1,1,1,2-tetrachloroethane | U | 40 |
| acetone | U | 200 | ethylbenzene | 43 | 40 |
| 1,1-dichloroethene | 660 | 20 | m&p-xylenes | 110 | 40 |
| methylene chloride | 870 | 100 | o-xylene | U | 40 |
| carbon disulfide | U | 40 | styrene | U | 40 |
| methyl t-butyl ether (MTBE) | U | 40 | bromoform | U | 40 |
| trans-1,2-dichloroethene | U | 40 | isopropylbenzene | U | 40 |
| 1,1-dichloroethane | 55 | 40 | 1,1,2,2-tetrachloroethane | U | 40 |
| 2-butanone (MEK) | U | 200 | 1,2,3-trichloropropane | U | 40 |
| 2,2-dichloropropane | U | 40 | n-propylbenzene | U | 40 |
| cis-1,2-dichloroethene | 73 | 40 | bromobenzene | U | 40 |
| chloroform | U | 40 | 1,3,5-trimethylbenzene | U | 40 |
| bromochloromethane | U | 40 | 2-chlorotoluene | U | 40 |
| tetrahydrofuran (THF) | U | 200 | 4-chlorotoluene | U | 40 |
| 1,1,1-trichloroethane | 4100 | 40 | tert-butylbenzene | U | 40 |
| 1,1-dichloropropene | U | 40 | 1,2,4-trimethylbenzene | U | 40 |
| carbon tetrachloride | U | 40 | sec-butylbenzene | U | 40 |
| 1,2-dichloroethane | 210 | 40 | 1,3-dichlorobenzene | U | 40 |
| benzene | U | 40 | 4-isopropyltoluene | U | 40 |
| trichloroethene | 12000 | 40 | 1,4-dichlorobenzene | U | 40 |
| 1,2-dichloropropane | U | 40 | 1,2-dichlorobenzene | 160 | 40 |
| bromodichloromethane | U | 40 | n-butylbenzene | U | 40 |
| dibromomethane | U | 40 | 1,2-dibromo-3-chloropropane | U | 40 |
| 4-methyl-2-pentanone (MIBK) | U | 200 | 1,2,4-trichlorobenzene | U | 40 |
| cis-1,3-dichloropropene | U | 40 | hexachlorobutadiene | U | 40 |
| toluene | 170 | 40 | naphthalene | U | 100 |
| trans-1,3-dichloropropene | U | 40 | 1,2,3-trichlorobenzene | U | 40 |
| 2-hexanone | U | 200 | | | |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 98 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 100 | 86-115 | | | |

U = Below quantitation limit

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10068

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 6 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

No exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

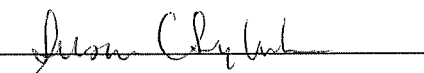
Other

A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

| MADEP MCP Analytical Method Report Certification Form | | | | | |
|--|--|----------|---------------|--------------------------|-------------|
| Laboratory Name: Resource Laboratories, LLC | | | | | Lab # 10068 |
| Project Location Hyde Park | | | Project # | MADEP RTN (if available) | |
| This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers) | | | | | |
| Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other: | | | | | |
| MCP SW-846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | | | Yes (x) No () | |
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | | | Yes (x) No () | |
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | | | Yes (x) No () | |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | | | Yes () No () NA | |
| E | Were all QC performance standards and recommendations for the specified methods achieved? | | | Yes (x) No () | |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | | | Yes (x) No () | |
| I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete. | | | | | |

Signature: 

Position: Lab Director

Printed Name: Susan C. Sylvester

Date: 2-2-07

Lab Number: 10068-50
Sample Designation: Method Blank
Date Sampled: N/A
Date Analyzed: 3/24/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 98 | 78-114 | | | |
| toluene-D8 | 103 | 88-110 | | | |
| 4-bromofluorobenzene | 100 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10068-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\VOA03\MAR06\032406\V3032424.D
Date Analyzed: 3/24/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-------------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 17 | 83% | 18 | 88% | 5% |
| chloromethane | 19 | 93% | 19 | 94% | 1% |
| vinyl chloride | 22 | 109% | 23 | 117% | 7% |
| bromomethane | 15 | 76% | 16 | 81% | 7% |
| chloroethane | 19 | 95% | 19 | 97% | 2% |
| trichlorofluoromethane | 20 | 102% | 21 | 106% | 3% |
| diethylether | 21 | 104% | 21 | 103% | 1% |
| acetone | 16 | 80% | 16 | 79% | 2% |
| 1,1-dichloroethene | 20 | 99% | 21 | 103% | 4% |
| methylene chloride | 21 | 106% | 21 | 105% | 1% |
| carbon disulfide | 19 | 95% | 19 | 96% | 1% |
| methyl-t-butyl ether (MTBE) | 40 | 100% | 40 | 99% | 0% |
| trans-1,2-dichloroethene | 20 | 100% | 21 | 106% | 6% |
| isopropyl ether (DIPE) | 21 | 105% | 20 | 102% | 3% |
| ethyl-t-butyl ether (ETBE) | 21 | 104% | 21 | 105% | 1% |
| 1,1-dichloroethane | 20 | 100% | 20 | 102% | 1% |
| t-butanol (TBA) | 88 | 88% | 86 | 86% | 2% |
| 2-butanone (MEK) | 17 | 86% | 17 | 86% | 0% |
| 2,2-dichloropropane | 19 | 96% | 20 | 98% | 3% |
| cis-1,2-dichloroethene | 21 | 105% | 21 | 105% | 0% |
| chloroform | 21 | 105% | 21 | 107% | 3% |
| bromochloromethane | 21 | 105% | 22 | 109% | 4% |
| tetrahydrofuran (THF) | 19 | 94% | 18 | 92% | 3% |
| 1,1,1-trichloroethane | 20 | 99% | 21 | 106% | 6% |
| 1,1-dichloropropene | 21 | 104% | 22 | 110% | 5% |
| t-amyl-methyl ether (TAME) | 19 | 97% | 21 | 103% | 6% |
| carbon tetrachloride | 17 | 83% | 18 | 91% | 9% |
| 1,2-dichloroethane | 21 | 103% | 20 | 102% | 1% |
| benzene | 21 | 104% | 21 | 105% | 1% |
| trichloroethene | 20 | 101% | 22 | 108% | 6% |
| 1,2-dichloropropane | 20 | 100% | 20 | 102% | 2% |
| bromodichloromethane | 18 | 90% | 19 | 95% | 5% |
| dibromomethane | 21 | 107% | 22 | 109% | 2% |
| 4-methyl-2-pentanone (MIBK) | 19 | 96% | 18 | 92% | 4% |
| cis-1,3-dichloropropene | 20 | 99% | 20 | 102% | 3% |
| toluene | 20 | 101% | 20 | 102% | 1% |
| trans-1,3-dichloropropene | 18 | 92% | 18 | 92% | 1% |
| 2-hexanone | 18 | 88% | 17 | 85% | 4% |
| 1,1,2-trichloroethane | 21 | 103% | 21 | 103% | 0% |
| 1,3-dichloropropane | 21 | 104% | 20 | 102% | 2% |
| tetrachloroethene | 21 | 103% | 21 | 107% | 4% |
| dibromochloromethane | 19 | 94% | 18 | 92% | 2% |
| 1,2-dibromoethane (EDB) | 20 | 102% | 20 | 99% | 4% |
| chlorobenzene | 20 | 100% | 20 | 101% | 2% |
| 1,1,1,2-tetrachloroethane | 19 | 95% | 19 | 97% | 2% |
| ethylbenzene | 21 | 104% | 21 | 104% | 0% |
| m&p-xylenes | 43 | 106% | 43 | 107% | 0% |
| o-xylene | 21 | 105% | 21 | 107% | 2% |
| styrene | 20 | 98% | 21 | 105% | 6% |
| bromoform | 18 | 91% | 18 | 89% | 2% |
| isopropylbenzene | 22 | 111% | 22 | 109% | 2% |
| 1,1,2,2-tetrachloroethane | 18 | 89% | 18 | 91% | 2% |
| 1,2,3-trichloropropane | 18 | 92% | 18 | 88% | 4% |
| n-propylbenzene | 21 | 105% | 21 | 107% | 2% |
| bromobenzene | 21 | 104% | 20 | 102% | 2% |
| 1,3,5-trimethylbenzene | 20 | 99% | 20 | 102% | 3% |
| 2-chlorotoluene | 19 | 96% | 20 | 102% | 7% |
| 4-chlorotoluene | 21 | 104% | 20 | 100% | 4% |
| tert-butylbenzene | 19 | 94% | 19 | 97% | 3% |
| 1,2,4-trimethylbenzene | 20 | 102% | 20 | 102% | 0% |
| sec-butylbenzene | 20 | 100% | 20 | 102% | 2% |
| 1,3-dichlorobenzene | 19 | 97% | 20 | 99% | 2% |
| 4-isopropyltoluene | 21 | 106% | 22 | 109% | 3% |
| 1,4-dichlorobenzene | 19 | 93% | 19 | 97% | 5% |
| 1,2-dichlorobenzene | 20 | 99% | 20 | 101% | 2% |
| n-butylbenzene | 21 | 103% | 22 | 111% | 8% |
| 1,2-dibromo-3-chloropropane (| 18 | 92% | 19 | 93% | 1% |
| 1,2,4-trichlorobenzene | 18 | 90% | 19 | 93% | 3% |
| hexachlorobutadiene | 19 | 97% | 20 | 100% | 3% |
| naphthalene | 15 | 76% | 15 | 76% | 0% |
| 1,2,3-trichlorobenzene | 18 | 91% | 18 | 92% | 1% |
| 1,4-dioxane | 33 | 83% | 36 | 91% | 9% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 98% | 100% |
| SS toluene-D8 | 101% | 103% |
| SS 4-bromofluorobenzene | 105% | 101% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

Resource Laboratories, LLC
124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

Company Name:

Phone #: 603-285-9760

Company Address:

FAX #: 603-285-9957

Project Manager:

Site Location (City, State):

Invoice To:

Protocol:

RCRA SDWA NPDES
MCP NHDES OTHER

Lab Sample ID

(Lab Use Only)

Field ID

CONTAINERS

WATER

SOLID

OTHER

HCl

HNO₃

H₂SO₄

ICE

MeOH

OTHER (Specify)

DATE

TIME

SAMPLER

☐ VOC-NH Petroleum Remediation Short List

☐ VOC-NH Hazardous Waste Remediation Short List

☐ VOC-NH Petroleum & Haz Waste Full List ☐ MADEP VPH ☐ MEGRO

☒ VOC 8260 ☐ VOC8015GRO ☐ VOC 624

☐ VOC 8260 BTEX, MIBE, Naphthalene only ☐ VOC 8260 + Oxygenates

☐ VOC 524.2 ☐ VOC 524.2 NH Petroleum & Haz. Waste Full List

☐ TPH 8100 ☐ MEDRO ☐ DRO 8015 ☐ EPH

☐ 8270PAH ☐ 8270ABN ☐ 625

☐ 8082 PCB ☐ 8081 Pesticides ☐ 608

☐ O&G 1664 ☐ O&G SM5520F

☐ pH ☐ BOD ☐ Conductivity

☐ TSS ☐ TDS ☐ TS

☐ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals

☐ Total Metals-list ☐ Dissolved Metals-list

☐ Ammonia ☐ COD

☐ T-Phosphate ☐ Phenol (subcontract)

☐ Cyanide ☐ Sulfide

☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride

☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitibility/FP

☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC

☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)

Grab (G) or Composite (C)

TAT REQUESTED

Priority (24 hr) ☒

Expedited (48 hr) ☐

10 Business Days ☐

Other ☐

E-Mail Address

Quote #

PO #

SPECIAL INSTRUCTIONS

REPORTING INSTRUCTIONS

☐ FAX ☐ EDD ☐ OTHER (specify)

RECEIVED ON ICE

☒ YES

☐ NO

TEMPERATURE

8.0

°C

Lab Use Only

CUSTODY RECORD

Relinquished by: *[Signature]*

Date: 3/23/06

Time: 1500

Received by:

Received by:

Date:

Time:

Relinquished by:

Date:

Time:

Received by:

Received by:

Date:

Time:

Relinquished by:

Date:

Time:

Received by:

Received by:

Date:

Time:

CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST

10068

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10070
Date Received: 3/24/06

Project: 2004-301 DND Lewis

Attached please find results for the analysis of the samples received on the date referenced above.


This is a reissue of report 10061, including the associated batch QC, case narrative and MCP Report Certification Form.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

2-2-07

Date

Total number of pages

8

Resource Laboratories, LLC Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Lab Number: 10070-01
 Sample Designation: ESM-03B (40-50)
 Date Sampled: 3/24/06
 Date Analyzed: 3/28/06
 Matrix: Water
 Instrument Dilution Factor: 200
 Analyst: LMM

VOLATILE ORGANICS
 SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 400 | trans-1,3-dichloropropene | U | 400 |
| chloromethane | U | 400 | 2-hexanone | U | 2000 |
| vinyl chloride | U | 400 | 1,1,2-trichloroethane | U | 400 |
| bromomethane | U | 400 | 1,3-dichloropropane | U | 400 |
| chloroethane | U | 400 | tetrachloroethene | 9200 | 400 |
| trichlorofluoromethane | U | 400 | dibromochloromethane | U | 400 |
| diethyl ether | U | 2000 | 1,2-dibromoethane | U | 400 |
| acetone | U | 2000 | chlorobenzene | U | 400 |
| 1,1-dichloroethene | 2500 | 200 | 1,1,1,2-tetrachloroethane | U | 400 |
| methylene chloride | 10000 | 1000 | ethylbenzene | U | 400 |
| carbon disulfide | U | 400 | m&p-xylenes | U | 400 |
| methyl t-butyl ether (MTBE) | U | 400 | o-xylene | U | 400 |
| trans-1,2-dichloroethene | U | 400 | styrene | U | 400 |
| isopropyl ether (DIPE) | U | 400 | bromoform | U | 400 |
| ethyl t-butyl ether (ETBE) | U | 400 | isopropylbenzene | U | 400 |
| 1,1-dichloroethane | U | 400 | 1,1,2,2-tetrachloroethane | U | 400 |
| t-butanol (TBA) | U | 10000 | 1,2,3-trichloropropane | U | 400 |
| 2-butanone (MEK) | U | 2000 | n-propylbenzene | U | 400 |
| 2,2-dichloropropane | U | 400 | bromobenzene | U | 400 |
| cis-1,2-dichloroethene | U | 400 | 1,3,5-trimethylbenzene | U | 400 |
| chloroform | U | 400 | 2-chlorotoluene | U | 400 |
| bromochloromethane | U | 400 | 4-chlorotoluene | U | 400 |
| tetrahydrofuran (THF) | U | 2000 | tert-butylbenzene | U | 400 |
| 1,1,1-trichloroethane | 19000 | 400 | 1,2,4-trimethylbenzene | U | 400 |
| 1,1-dichloropropene | U | 400 | sec-butylbenzene | U | 400 |
| t-amyl-methyl ether (TAME) | U | 400 | 1,3-dichlorobenzene | U | 400 |
| carbon tetrachloride | U | 400 | 4-isopropyltoluene | U | 400 |
| 1,2-dichloroethane | 1500 | 400 | 1,4-dichlorobenzene | U | 400 |
| benzene | U | 400 | 1,2-dichlorobenzene | U | 400 |
| trichloroethene | 50000 | 400 | n-butylbenzene | U | 400 |
| 1,2-dichloropropane | U | 400 | 1,2-dibromo-3-chloropropane | U | 400 |
| bromodichloromethane | U | 400 | 1,2,4-trichlorobenzene | U | 400 |
| dibromomethane | U | 400 | hexachlorobutadiene | U | 400 |
| 4-methyl-2-pentanone (MIBK) | U | 2000 | naphthalene | U | 1000 |
| cis-1,3-dichloropropene | U | 400 | 1,2,3-trichlorobenzene | U | 400 |
| toluene | 1200 | 400 | 1,4-dioxane | U | 10000 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 101 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 99 | 86-115 | | | |

U = Below quantitation limit

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10070

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 10 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

The LCS showed a high recovery for Tetrachloroethene. This was most likely due to carryover from a previous sample analysis. The LCSD results were acceptable.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10070

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|------------|----------|---------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |

| | | |
|---|---|---------------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|---|---------------------|
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
|---|---|---------------------|

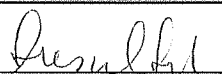
| | | |
|---|--|---------------------|
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
|---|--|---------------------|

| | | |
|---|---|----------------------|
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No () NA |
|---|---|----------------------|

| | | |
|---|---|---------------------|
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|--|---------------------|
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes (x) No () |
|---|--|---------------------|

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Position: Lab DirectorPrinted Name: Susan C. SylvesterDate: 2-2-07

Lab Number: 10070-50
 Sample Designation: Method Blank
 Date Sampled: N/A
 Date Analyzed: 3/28/06
 Matrix: Water
 Instrument Dilution Factor: 1
 Analyst: LMM

VOLATILE ORGANICS
 SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |

| SURROGATE STANDARDS | Recovery (%) | Acceptance Limits (%) |
|----------------------|--------------|-----------------------|
| dibromofluoromethane | 98 | 78-114 |
| toluene-D8 | 101 | 88-110 |
| 4-bromofluorobenzene | 101 | 86-115 |

U = Below quantitation limit

Lab Number: 10070-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\VOA03\032806\V3032824.D
Date Analyzed: 3/28/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-------------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 18 | 92% | 18 | 90% | 2% |
| chloromethane | 20 | 101% | 20 | 99% | 2% |
| vinyl chloride | 23 | 115% | 23 | 114% | 1% |
| bromomethane | 19 | 95% | 19 | 98% | 1% |
| chloroethane | 21 | 108% | 21 | 103% | 3% |
| trichlorofluoromethane | 22 | 112% | 22 | 108% | 4% |
| diethylether | 21 | 103% | 22 | 111% | 7% |
| acetone | 17 | 87% | 17 | 86% | 1% |
| 1,1-dichloroethene | 22 | 109% | 21 | 105% | 3% |
| methylene chloride | 23 | 116% | 23 | 113% | 3% |
| carbon disulfide | 20 | 101% | 20 | 98% | 4% |
| methyl-t-butyl ether (MTBE) | 41 | 103% | 41 | 103% | 0% |
| trans-1,2-dichloroethene | 22 | 110% | 21 | 105% | 4% |
| isopropyl ether (DIPE) | 21 | 107% | 22 | 111% | 4% |
| ethyl-t-butyl ether (ETBE) | 22 | 109% | 21 | 105% | 4% |
| 1,1-dichloroethane | 21 | 107% | 21 | 105% | 2% |
| t-butanol (TBA) | 85 | 85% | 87 | 87% | 3% |
| 2-butanone (MEK) | 18 | 89% | 18 | 90% | 2% |
| 2,2-dichloropropane | 20 | 99% | 19 | 94% | 5% |
| cis-1,2-dichloroethene | 23 | 116% | 23 | 114% | 2% |
| chloroform | 22 | 110% | 22 | 109% | 1% |
| bromochloromethane | 23 | 113% | 23 | 114% | 1% |
| tetrahydrofuran (THF) | 17 | 87% | 18 | 90% | 4% |
| 1,1,1-trichloroethane | 21 | 106% | 21 | 103% | 2% |
| 1,1-dichloropropene | 23 | 114% | 21 | 107% | 6% |
| t-amyl-methyl ether (TAME) | 21 | 105% | 21 | 104% | 1% |
| carbon tetrachloride | 18 | 92% | 18 | 92% | 1% |
| 1,2-dichloroethane | 21 | 107% | 21 | 103% | 4% |
| benzene | 22 | 109% | 21 | 107% | 2% |
| trichloroethene | 23 | 115% | 22 | 109% | 5% |
| 1,2-dichloropropane | 21 | 107% | 21 | 107% | 0% |
| bromodichloromethane | 19 | 95% | 20 | 98% | 2% |
| dibromomethane | 23 | 113% | 23 | 113% | 0% |
| 4-methyl-2-pentanone (MIBK) | 19 | 97% | 19 | 98% | 1% |
| cis-1,3-dichloropropene | 21 | 106% | 21 | 106% | 0% |
| toluene | 23 | 114% | 22 | 110% | 4% |
| trans-1,3-dichloropropene | 19 | 98% | 19 | 97% | 1% |
| 2-hexanone | 17 | 84% | 17 | 87% | 4% |
| 1,1,2-trichloroethane | 22 | 109% | 22 | 111% | 2% |
| 1,3-dichloropropane | 22 | 108% | 20 | 102% | 5% |
| tetrachloroethene | 29 | 146% # | 25 | 125% | 15% |
| dibromochloromethane | 19 | 97% | 20 | 98% | 2% |
| 1,2-dibromoethane (EDB) | 21 | 105% | 21 | 104% | 1% |
| chlorobenzene | 21 | 108% | 20 | 100% | 5% |
| 1,1,1,2-tetrachloroethane | 20 | 102% | 20 | 99% | 3% |
| ethylbenzene | 22 | 109% | 21 | 103% | 6% |
| m&p-xylenes | 44 | 109% | 42 | 105% | 3% |
| o-xylene | 22 | 109% | 21 | 104% | 5% |
| styrene | 20 | 100% | 20 | 102% | 3% |
| bromoform | 18 | 92% | 19 | 93% | 2% |
| isopropylbenzene | 23 | 114% | 22 | 112% | 1% |
| 1,1,2,2-tetrachloroethane | 18 | 91% | 18 | 90% | 1% |
| 1,2,3-trichloropropane | 18 | 89% | 17 | 86% | 3% |
| n-propylbenzene | 22 | 109% | 21 | 105% | 3% |
| bromobenzene | 22 | 108% | 21 | 105% | 3% |
| 1,3,5-trimethylbenzene | 20 | 102% | 20 | 99% | 3% |
| 2-chlorotoluene | 21 | 103% | 20 | 99% | 4% |
| 4-chlorotoluene | 21 | 104% | 20 | 99% | 5% |
| tert-butylbenzene | 20 | 99% | 19 | 94% | 6% |
| 1,2,4-trimethylbenzene | 21 | 108% | 20 | 100% | 6% |
| sec-butylbenzene | 20 | 102% | 20 | 98% | 4% |
| 1,3-dichlorobenzene | 20 | 102% | 20 | 100% | 2% |
| 4-isopropyltoluene | 22 | 111% | 21 | 107% | 4% |
| 1,4-dichlorobenzene | 20 | 100% | 19 | 97% | 3% |
| 1,2-dichlorobenzene | 21 | 106% | 20 | 100% | 5% |
| n-butylbenzene | 22 | 110% | 21 | 103% | 6% |
| 1,2-dibromo-3-chloropropane (| 18 | 88% | 18 | 92% | 4% |
| 1,2,4-trichlorobenzene | 18 | 92% | 18 | 89% | 3% |
| hexachlorobutadiene | 20 | 100% | 19 | 94% | 6% |
| naphthalene | 15 | 75% | 15 | 78% | 1% |
| 1,2,3-trichlorobenzene | 19 | 93% | 18 | 89% | 4% |
| 1,4-dioxane | 34 | 85% | 34 | 86% | 0% |

SURROGATE STANDARDS

| | | |
|---------------------------|------|------|
| SS dibromodifluoromethane | 98% | 103% |
| SS toluene-D8 | 105% | 105% |
| SS 4-bromodifluorobenzene | 105% | 107% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

= This analyte showed recovery outside the acceptance limits.



Resource Laboratories, LLC
124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST

ANALYSIS REQUEST

10070

PAGE ____ OF ____

Company Name:

Phone #: 508 285 9700

Company Address:

FAX #: 508 285 9757

Project Manager:

Site Location (City, State):

184 West Main St. North MA
Hyde Park, MA

Project ID / Name: DUD Lewis 2004-301

Invoice To:

Protocol: RCRA SDWA NPDES
MCE NHDES OTHER

Lab Sample ID (Lab Use Only)

Field ID

CONTAINERS

MATRIX

PRESERVATION METHOD

SAMPLING

WATER

SOLID

OTHER

HCl

HNO₃

H₂SO₄

ICE

MeOH

OTHER (Specify)

DATE

TIME

SAMPLER

- ☐ VOC-NH Petroleum Remediation Short List
☐ VOC-NH Hazardous Waste Remediation Short List
☐ VOC-NH Petroleum & Haz Waste Full List ☐ MADEP VPH ☐ MEGRO
☒ VOC 8260 ☐ VOC8015GRO ☐ VOC 624
☐ VOC 8260 BTEX, MIBE, Naphthalene only ☐ VOC 8260 + Oxygenates
☐ VOC 524.2 ☐ VOC 524.2 NH Petroleum & Haz. Waste Full List
☐ TPH 8100 ☐ MEDRO ☐ DRO 8015 ☐ EPH
☐ 8270PAH ☐ 8270ABN ☐ 625
☐ 8082 PCB ☐ 8081 Pesticides ☐ 608
☐ O&G 1664 ☐ O&G SM5520F
☐ pH ☐ BOD ☐ Conductivity
☐ TSS ☐ TDS ☐ TS
☐ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals
☐ Total Metals-list ☐ Dissolved Metals-list
☐ Ammonia ☐ COD
☐ T-Phosphate ☐ Phenol (subcontract)
☐ Cyanide ☐ Sulfide
☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride
☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitibility/FP
☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC
☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)

Grab (G) or Composite (C)

TAT REQUESTED

Priority (24 hr) ☒

Expected (48 hr) ☐

10 Business Days ☐

Other ☐

E-Mail Address

Quote #

PO #

SPECIAL INSTRUCTIONS

per site specific agreement, TAT-ASAP

REPORTING INSTRUCTIONS

☐ FAX ☐ EDD ☒ OTHER (specify)

call to callahan

RECEIVED ON ICE

TEMPERATURE

YES ☒ NO ☐

10 °C

Lab Use Only

CUSTODY RECORD

Relinquished by Sampler:

Date

Time

Received by:

Received by:

Received by Laboratory:

Date

Time

Relinquished by:

Date

Time

Received by:

Received by:

Received by Laboratory:

Date

Time

Relinquished by:

Date

Time

Received by:

Received by:

Received by Laboratory:

Date

Time

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10079
Date Received: 3/27/06

Project: 2004-301 Lewis Chemical

Attached please find results for the analysis of the samples received on the date referenced above.

This is a reissue of report 10061, including the associated batch QC, case narrative and MCP Report Certification Form.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

2-2-07
Date

Total number of pages

8

Resource Laboratories, LLC Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Lab Number: 10079-01
Sample Designation: ESM-05B (20-30)
Date Sampled: 3/27/06
Date Analyzed: 3/28/06
Matrix: Water
Instrument Dilution Factor: 200
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 400 | trans-1,3-dichloropropene | U | 400 |
| chloromethane | U | 400 | 2-hexanone | U | 2000 |
| vinyl chloride | U | 400 | 1,1,2-trichloroethane | U | 400 |
| bromomethane | U | 400 | 1,3-dichloropropane | U | 400 |
| chloroethane | U | 400 | tetrachloroethene | 23000 | 400 |
| trichlorofluoromethane | U | 400 | dibromochloromethane | U | 400 |
| diethyl ether | U | 2000 | 1,2-dibromoethane | U | 400 |
| acetone | U | 2000 | chlorobenzene | U | 400 |
| 1,1-dichloroethene | 2900 | 200 | 1,1,1,2-tetrachloroethane | U | 400 |
| methylene chloride | 2000 | 1000 | ethylbenzene | U | 400 |
| carbon disulfide | U | 400 | m&p-xylenes | 750 | 400 |
| methyl t-butyl ether (MTBE) | U | 400 | o-xylene | U | 400 |
| trans-1,2-dichloroethene | U | 400 | styrene | U | 400 |
| isopropyl ether (DIPE) | U | 400 | bromoform | U | 400 |
| ethyl t-butyl ether (ETBE) | U | 400 | isopropylbenzene | U | 400 |
| 1,1-dichloroethane | U | 400 | 1,1,2,2-tetrachloroethane | U | 400 |
| t-butanol (TBA) | U | 10000 | 1,2,3-trichloropropane | U | 400 |
| 2-butanone (MEK) | U | 2000 | n-propylbenzene | U | 400 |
| 2,2-dichloropropane | U | 400 | bromobenzene | U | 400 |
| cis-1,2-dichloroethene | 460 | 400 | 1,3,5-trimethylbenzene | U | 400 |
| chloroform | U | 400 | 2-chlorotoluene | U | 400 |
| bromochloromethane | U | 400 | 4-chlorotoluene | U | 400 |
| tetrahydrofuran (THF) | U | 2000 | tert-butylbenzene | U | 400 |
| 1,1,1-trichloroethane | 38000 | 400 | 1,2,4-trimethylbenzene | U | 400 |
| 1,1-dichloropropene | U | 400 | sec-butylbenzene | U | 400 |
| t-amyl-methyl ether (TAME) | U | 400 | 1,3-dichlorobenzene | U | 400 |
| carbon tetrachloride | U | 400 | 4-isopropyltoluene | U | 400 |
| 1,2-dichloroethane | U | 400 | 1,4-dichlorobenzene | U | 400 |
| benzene | U | 400 | 1,2-dichlorobenzene | 2000 | 400 |
| trichloroethene | 54000 | 400 | n-butylbenzene | U | 400 |
| 1,2-dichloropropane | U | 400 | 1,2-dibromo-3-chloropropane | U | 400 |
| bromodichloromethane | U | 400 | 1,2,4-trichlorobenzene | U | 400 |
| dibromomethane | U | 400 | hexachlorobutadiene | U | 400 |
| 4-methyl-2-pentanone (MIBK) | U | 2000 | naphthalene | U | 1000 |
| cis-1,3-dichloropropene | U | 400 | 1,2,3-trichlorobenzene | U | 400 |
| toluene | 1500 | 400 | 1,4-dioxane | U | 10000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 100 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 105 | 86-115 | | | |

U = Below quantitation limit

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10079

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 2 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

The LCS showed a high recovery for Tetrachloroethene. This was most likely due to carryover from a previous sample analysis. The LCSD results were acceptable.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10079

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|------------|----------|---------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |

| | | |
|---|---|------------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
|---|---|------------------|

| | | |
|---|---|------------------|
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
|---|---|------------------|

| | | |
|---|--|------------------|
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
|---|--|------------------|

| | | |
|---|---|-------------------|
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No () NA |
|---|---|-------------------|

| | | |
|---|---|------------------|
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes () No (x) |
|---|---|------------------|

| | | |
|---|--|------------------|
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes (x) No () |
|---|--|------------------|

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Susan C. SylvesterPosition: Lab DirectorPrinted Name: Susan C. SylvesterDate: 2-2-07

Lab Number: 10079-50
 Sample Designation: Method Blank
 Date Sampled: N/A
 Date Analyzed: 3/28/06
 Matrix: Water
 Instrument Dilution Factor: 1
 Analyst: LMM

VOLATILE ORGANICS
 SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 98 | 78-114 | | | |
| toluene-D8 | 101 | 88-110 | | | |
| 4-bromofluorobenzene | 101 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10079-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\VOA03\032806\W3032824.D
Date Analyzed: 3/28/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-------------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 18 | 92% | 18 | 90% | 2% |
| chloromethane | 20 | 101% | 20 | 99% | 2% |
| vinyl chloride | 23 | 115% | 23 | 114% | 1% |
| bromomethane | 19 | 95% | 19 | 98% | 1% |
| chloroethane | 21 | 106% | 21 | 103% | 3% |
| trichlorofluoromethane | 22 | 112% | 22 | 108% | 4% |
| diethylether | 21 | 103% | 22 | 111% | 7% |
| acetone | 17 | 87% | 17 | 88% | 1% |
| 1,1-dichloroethene | 22 | 109% | 21 | 105% | 3% |
| methylene chloride | 23 | 116% | 23 | 113% | 3% |
| carbon disulfide | 20 | 101% | 20 | 98% | 4% |
| methyl-t-butyl ether (MTBE) | 41 | 103% | 41 | 103% | 0% |
| trans-1,2-dichloroethene | 22 | 110% | 21 | 105% | 4% |
| isopropyl ether (DIPE) | 21 | 107% | 22 | 111% | 4% |
| ethyl-t-butyl ether (ETBE) | 22 | 108% | 21 | 105% | 4% |
| 1,1-dichloroethane | 21 | 107% | 21 | 105% | 2% |
| t-butanol (TBA) | 85 | 85% | 87 | 87% | 3% |
| 2-butanone (MEK) | 18 | 89% | 18 | 90% | 2% |
| 2,2-dichloropropane | 20 | 98% | 19 | 94% | 5% |
| cis-1,2-dichloroethene | 23 | 116% | 23 | 114% | 2% |
| chloroform | 22 | 110% | 22 | 109% | 1% |
| bromochloromethane | 23 | 113% | 23 | 114% | 1% |
| tetrahydrofuran (THF) | 17 | 87% | 18 | 90% | 4% |
| 1,1,1-trichloroethane | 21 | 106% | 21 | 103% | 2% |
| 1,1-dichloropropene | 23 | 114% | 21 | 107% | 6% |
| t-amyl-methyl ether (TAME) | 21 | 105% | 21 | 104% | 1% |
| carbon tetrachloride | 18 | 92% | 18 | 92% | 1% |
| 1,2-dichloroethane | 21 | 107% | 21 | 103% | 4% |
| benzene | 22 | 109% | 21 | 107% | 2% |
| trichloroethene | 23 | 115% | 22 | 109% | 5% |
| 1,2-dichloropropane | 21 | 107% | 21 | 107% | 0% |
| bromodichloromethane | 19 | 95% | 20 | 98% | 2% |
| dibromomethane | 23 | 113% | 23 | 113% | 0% |
| 4-methyl-2-pentanone (MIBK) | 19 | 97% | 19 | 98% | 1% |
| cis-1,3-dichloropropene | 21 | 108% | 21 | 108% | 0% |
| toluene | 23 | 114% | 22 | 110% | 4% |
| trans-1,3-dichloropropene | 19 | 98% | 19 | 97% | 1% |
| 2-hexanone | 17 | 84% | 17 | 87% | 4% |
| 1,1,2-trichloroethane | 22 | 109% | 22 | 111% | 2% |
| 1,3-dichloropropane | 22 | 108% | 20 | 102% | 5% |
| tetrachloroethene | 29 | 146% # | 25 | 125% | 15% |
| dibromochloromethane | 19 | 97% | 20 | 98% | 2% |
| 1,2-dibromoethane (EDB) | 21 | 105% | 21 | 104% | 1% |
| chlorobenzene | 21 | 106% | 20 | 100% | 5% |
| 1,1,1,2-tetrachloroethane | 20 | 102% | 20 | 99% | 3% |
| ethylbenzene | 22 | 109% | 21 | 103% | 6% |
| m&p-xylenes | 44 | 109% | 42 | 105% | 3% |
| o-xylene | 22 | 109% | 21 | 104% | 5% |
| styrene | 20 | 100% | 20 | 102% | 3% |
| bromoform | 18 | 92% | 19 | 93% | 2% |
| isopropylbenzene | 23 | 114% | 22 | 112% | 1% |
| 1,1,2,2-tetrachloroethane | 18 | 91% | 18 | 90% | 1% |
| 1,2,3-trichloropropane | 18 | 89% | 17 | 86% | 3% |
| n-propylbenzene | 22 | 109% | 21 | 105% | 3% |
| bromobenzene | 22 | 108% | 21 | 105% | 3% |
| 1,3,5-trimethylbenzene | 20 | 102% | 20 | 99% | 3% |
| 2-chlorotoluene | 21 | 103% | 20 | 99% | 4% |
| 4-chlorotoluene | 21 | 104% | 20 | 99% | 5% |
| tert-butylbenzene | 20 | 99% | 19 | 94% | 6% |
| 1,2,4-trimethylbenzene | 21 | 108% | 20 | 100% | 6% |
| sec-butylbenzene | 20 | 102% | 20 | 98% | 4% |
| 1,3-dichlorobenzene | 20 | 102% | 20 | 100% | 2% |
| 4-isopropyltoluene | 22 | 111% | 21 | 107% | 4% |
| 1,4-dichlorobenzene | 20 | 100% | 19 | 97% | 3% |
| 1,2-dichlorobenzene | 21 | 106% | 20 | 100% | 5% |
| n-butylbenzene | 22 | 110% | 21 | 103% | 8% |
| 1,2-dibromo-3-chloropropane (| 18 | 88% | 18 | 92% | 4% |
| 1,2,4-trichlorobenzene | 18 | 92% | 18 | 89% | 3% |
| hexachlorobutadiene | 20 | 100% | 19 | 94% | 6% |
| naphthalene | 15 | 75% | 15 | 76% | 1% |
| 1,2,3-trichlorobenzene | 19 | 93% | 18 | 89% | 4% |
| 1,4-dioxane | 34 | 85% | 34 | 86% | 0% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 98% | 103% |
| SS toluene-D8 | 105% | 105% |
| SS 4-bromofluorobenzene | 105% | 107% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

= This analyte showed recovery outside the acceptance limits.

RL

Resource Laboratories, LLC
124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

CHAIN-OF-CUSTODY REC 10079
AND ANALYSIS REQUEST

ANALYSIS REQUEST

Company Name:

Phone #: (603) 285-9700

Environmental Strategies & Mgmt
Company Address:
184 West Main St. North MA

FAX #: (603) 285-9957
Site Location (City, State):
North Haverhill, MA

Project Manager:

Project ID / Name:

be Callahan

2004-201 / Lewis Chemical

Invoice To:

Protocol: RCRA SDWA NPDES
MGP NHDES OTHER

ESM

Lab

Sample ID

Field ID

CONTAINERS

Matrix

Preservation Method

Sampling

SAMPLER

(Lab Use Only)

10079-01

ESM 05 B (20-35)

2

WATER
SOLID
OTHER

HCl
HNO₃
H₂SO₄
NaOH
MeOH
OTHER (Specify)

DATE

TIME

3/27/06 1300 AF

☐ VOC 8260-NH List ☐ MADEP VPH ☐ MEGRO

☒ VOC 8260 ☐ VOC 8015GRO ☐ VOC 624

☐ VOC 8260 BTEX, MIB, Naphthalene only

☐ VOC 524.2 ☐ VOC 524.2 NH List

☐ TPH Fingerprint ☐ MEDRO ☐ DRO 8015 ☐ EPH

☐ 8270PAH ☐ 8270ABN ☐ 625

☐ 8082 PCB ☐ 8081 Pesticides ☐ 608

☐ O&G 1664 ☐ O&G SM5520F

☐ pH ☐ BOD ☐ Conductivity

☐ TSS ☐ TDS ☐ TS

☐ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals

☐ Total Metals-list ☐ Dissolved Metals-list

☐ Ammonia ☐ COD

☐ T-Phosphate ☐ Phenol

☐ Cyanide ☐ Sulfide

☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride

☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitability/FP

☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC

☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)

☐ Standard Drinking Water Test ☐ Bacteria P/A

Grab (G) or Composite (C)

TEST REQUESTED

Priority (24 hr) ☒

Expedited (48 hr) ☐

10 Business Days ☐

Other ☐

E-Mail Address

Quote #

PO #

SPECIAL INSTRUCTIONS

REPORTING INSTRUCTIONS

☒ FAX ☐ OTHER (specify)

☐ PDF ☐ Excel Spreadsheet

CUSTODY RECORD

Relinquished by Sampler:

Relinquished by:

Date

Date

Time

Time

Received by:

Received by:

RECEIVED ON ICE

TEMPERATURE

YES

NO

Date

Date

Time

Time

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Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10088
Date Received: 3/28/06

Project: 2004-301 Lewis Chemical

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

7-14-06

Date

Total number of pages

10

Resource Laboratories, LLC Certifications

New Hampshire NH902
Maine NH903

Connecticut PH-0146
Massachusetts M-NH902

Lab Number: 10088-01
Sample Designation: ESM-05B (30-40)
Date Sampled: 3/28/06
Date Analyzed: 3/29/06
Matrix: Water
Instrument Dilution Factor: 200
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 400 | trans-1,3-dichloropropene | U | 400 |
| chloromethane | U | 400 | 2-hexanone | U | 2000 |
| vinyl chloride | U | 400 | 1,1,2-trichloroethane | U | 400 |
| bromomethane | U | 400 | 1,3-dichloropropane | U | 400 |
| chloroethane | U | 400 | tetrachloroethene | 12000 | 400 |
| trichlorofluoromethane | U | 400 | dibromochloromethane | U | 400 |
| diethyl ether | U | 2000 | 1,2-dibromoethane | U | 400 |
| acetone | U | 2000 | chlorobenzene | U | 400 |
| 1,1-dichloroethene | 1700 | 200 | 1,1,1,2-tetrachloroethane | U | 400 |
| methylene chloride | 3800 | 1000 | ethylbenzene | U | 400 |
| carbon disulfide | U | 400 | m&p-xylenes | 420 | 400 |
| methyl t-butyl ether (MTBE) | U | 400 | o-xylene | U | 400 |
| trans-1,2-dichloroethene | U | 400 | styrene | U | 400 |
| isopropyl ether (DIPE) | U | 400 | bromoform | U | 400 |
| ethyl t-butyl ether (ETBE) | U | 400 | isopropylbenzene | U | 400 |
| 1,1-dichloroethane | U | 400 | 1,1,2,2-tetrachloroethane | U | 400 |
| t-butanol (TBA) | U | 10000 | 1,2,3-trichloropropane | U | 400 |
| 2-butanone (MEK) | U | 2000 | n-propylbenzene | U | 400 |
| 2,2-dichloropropane | U | 400 | bromobenzene | U | 400 |
| cis-1,2-dichloroethene | U | 400 | 1,3,5-trimethylbenzene | U | 400 |
| chloroform | U | 400 | 2-chlorotoluene | U | 400 |
| bromochloromethane | U | 400 | 4-chlorotoluene | U | 400 |
| tetrahydrofuran (THF) | U | 2000 | tert-butylbenzene | U | 400 |
| 1,1,1-trichloroethane | 29000 | 400 | 1,2,4-trimethylbenzene | U | 400 |
| 1,1-dichloropropene | U | 400 | sec-butylbenzene | U | 400 |
| t-amyl-methyl ether (TAME) | U | 400 | 1,3-dichlorobenzene | U | 400 |
| carbon tetrachloride | U | 400 | 4-isopropyltoluene | U | 400 |
| 1,2-dichloroethane | U | 400 | 1,4-dichlorobenzene | U | 400 |
| benzene | U | 400 | 1,2-dichlorobenzene | 960 | 400 |
| trichloroethene | 31000 | 400 | n-butylbenzene | U | 400 |
| 1,2-dichloropropane | U | 400 | 1,2-dibromo-3-chloropropane | U | 400 |
| bromodichloromethane | U | 400 | 1,2,4-trichlorobenzene | U | 400 |
| dibromomethane | U | 400 | hexachlorobutadiene | U | 400 |
| 4-methyl-2-pentanone (MIBK) | U | 2000 | naphthalene | U | 1000 |
| cis-1,3-dichloropropene | U | 400 | 1,2,3-trichlorobenzene | U | 400 |
| toluene | 910 | 400 | 1,4-dioxane | U | 10000 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dlbromofluoromethane | 102 | 78-114 | | | |
| toluene-D8 | 109 | 88-110 | | | |
| 4-bromofluorobenzene | 102 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10088-02
Sample Designation: ESM-05B (40-50)
Date Sampled: 3/28/06
Date Analyzed: 3/29/06
Matrix: Water
Instrument Dilution Factor: 100
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 200 | trans-1,3-dichloropropene | U | 200 |
| chloromethane | U | 200 | 2-hexanone | U | 1000 |
| vinyl chloride | U | 200 | 1,1,2-trichloroethane | U | 200 |
| bromomethane | U | 200 | 1,3-dichloropropane | U | 200 |
| chloroethane | U | 200 | tetrachloroethene | 3200 | 200 |
| trichlorofluoromethane | U | 200 | dibromochloromethane | U | 200 |
| diethyl ether | U | 1000 | 1,2-dibromoethane | U | 200 |
| acetone | U | 1000 | chlorobenzene | U | 200 |
| 1,1-dichloroethene | 590 | 100 | 1,1,1,2-tetrachloroethane | U | 200 |
| methylene chloride | 4100 | 500 | ethylbenzene | U | 200 |
| carbon disulfide | U | 200 | m&p-xylenes | U | 200 |
| methyl t-butyl ether (MTBE) | U | 200 | o-xylene | U | 200 |
| trans-1,2-dichloroethene | U | 200 | styrene | U | 200 |
| isopropyl ether (DIPE) | U | 200 | bromoform | U | 200 |
| ethyl t-butyl ether (ETBE) | U | 200 | isopropylbenzene | U | 200 |
| 1,1-dichloroethane | U | 200 | 1,1,2,2-tetrachloroethane | U | 200 |
| t-butanol (TBA) | U | 5000 | 1,2,3-trichloropropane | U | 200 |
| 2-butanone (MEK) | U | 1000 | n-propylbenzene | U | 200 |
| 2,2-dichloropropane | U | 200 | bromobenzene | U | 200 |
| cis-1,2-dichloroethene | U | 200 | 1,3,5-trimethylbenzene | U | 200 |
| chloroform | U | 200 | 2-chlorotoluene | U | 200 |
| bromochloromethane | U | 200 | 4-chlorotoluene | U | 200 |
| tetrahydrofuran (THF) | U | 1000 | tert-butylbenzene | U | 200 |
| 1,1,1-trichloroethane | 9000 | 200 | 1,2,4-trimethylbenzene | U | 200 |
| 1,1-dichloropropene | U | 200 | sec-butylbenzene | U | 200 |
| t-amyl-methyl ether (TAME) | U | 200 | 1,3-dichlorobenzene | U | 200 |
| carbon tetrachloride | U | 200 | 4-isopropyltoluene | U | 200 |
| 1,2-dichloroethane | 510 | 200 | 1,4-dichlorobenzene | U | 200 |
| benzene | U | 200 | 1,2-dichlorobenzene | 420 | 200 |
| trichloroethene | 14000 | 200 | n-butylbenzene | U | 200 |
| 1,2-dichloropropane | U | 200 | 1,2-dibromo-3-chloropropane | U | 200 |
| bromodichloromethane | U | 200 | 1,2,4-trichlorobenzene | U | 200 |
| dibromomethane | U | 200 | hexachlorobutadiene | U | 200 |
| 4-methyl-2-pentanone (MIBK) | U | 1000 | naphthalene | U | 500 |
| cis-1,3-dichloropropene | U | 200 | 1,2,3-trichlorobenzene | U | 200 |
| toluene | 310 | 200 | 1,4-dioxane | U | 5000 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| dibromofluoromethane | 101 | 78-114 |
| toluene-D8 | 105 | 88-110 |
| 4-bromofluorobenzene | 96 | 86-115 |

U = Below quantitation limit

Lab Number: 10088-03
 Sample Designation: ESM-05B (50-60)
 Date Sampled: 3/28/06
 Date Analyzed: 3/29/06
 Matrix: Water
 Instrument Dilution Factor: 100
 Analyst: LMM

VOLATILE ORGANICS
 SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 200 | trans-1,3-dichloropropene | U | 200 |
| chloromethane | U | 200 | 2-hexanone | U | 1000 |
| vinyl chloride | U | 200 | 1,1,2-trichloroethane | U | 200 |
| bromomethane | U | 200 | 1,3-dichloropropane | U | 200 |
| chloroethane | U | 200 | tetrachloroethene | 5200 | 200 |
| trichlorofluoromethane | U | 200 | dibromochloromethane | U | 200 |
| diethyl ether | U | 1000 | 1,2-dibromoethane | U | 200 |
| acetone | U | 1000 | chlorobenzene | U | 200 |
| 1,1-dichloroethene | 440 | 100 | 1,1,1,2-tetrachloroethane | U | 200 |
| methylene chloride | 2500 | 500 | ethylbenzene | U | 200 |
| carbon disulfide | U | 200 | m&p-xylenes | U | 200 |
| methyl t-butyl ether (MTBE) | U | 200 | o-xylene | U | 200 |
| trans-1,2-dichloroethene | U | 200 | styrene | U | 200 |
| isopropyl ether (DIPE) | U | 200 | bromoform | U | 200 |
| ethyl t-butyl ether (ETBE) | U | 200 | isopropylbenzene | U | 200 |
| 1,1-dichloroethane | U | 200 | 1,1,2,2-tetrachloroethane | U | 200 |
| t-butanol (TBA) | U | 5000 | 1,2,3-trichloropropane | U | 200 |
| 2-butanone (MEK) | U | 1000 | n-propylbenzene | U | 200 |
| 2,2-dichloropropane | U | 200 | bromobenzene | U | 200 |
| cis-1,2-dichloroethene | U | 200 | 1,3,5-trimethylbenzene | U | 200 |
| chloroform | U | 200 | 2-chlorotoluene | U | 200 |
| bromochloromethane | U | 200 | 4-chlorotoluene | U | 200 |
| tetrahydrofuran (THF) | U | 1000 | tert-butylbenzene | U | 200 |
| 1,1,1-trichloroethane | 9200 | 200 | 1,2,4-trimethylbenzene | U | 200 |
| 1,1-dichloropropene | U | 200 | sec-butylbenzene | U | 200 |
| t-amyl-methyl ether (TAME) | U | 200 | 1,3-dichlorobenzene | U | 200 |
| carbon tetrachloride | U | 200 | 4-isopropyltoluene | U | 200 |
| 1,2-dichloroethane | 300 | 200 | 1,4-dichlorobenzene | U | 200 |
| benzene | U | 200 | 1,2-dichlorobenzene | U | 200 |
| trichloroethene | 11000 | 200 | n-butylbenzene | U | 200 |
| 1,2-dichloropropane | U | 200 | 1,2-dibromo-3-chloropropane | U | 200 |
| bromodichloromethane | U | 200 | 1,2,4-trichlorobenzene | U | 200 |
| dibromomethane | U | 200 | hexachlorobutadiene | U | 200 |
| 4-methyl-2-pentanone (MIBK) | U | 1000 | naphthalene | U | 500 |
| cis-1,3-dichloropropene | U | 200 | 1,2,3-trichlorobenzene | U | 200 |
| toluene | 610 | 200 | 1,4-dioxane | U | 5000 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 100 | 78-114 | | | |
| toluene-D8 | 103 | 88-110 | | | |
| 4-bromofluorobenzene | 98 | 86-115 | | | |

U = Below quantitation limit

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10088

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 6 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

No exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested by the customer.

Other

VOC: The following compounds were quantified with quadratic fit:
dichlorodifluoromethane, chloromethane, vinyl chloride, bromomethane, acetone,
dibromochloromethane, isopropylbenzene, tert-butylbenzene, and hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10088

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|------------|----------|---------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |

| | | |
|---|--|----------------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No () NA |
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes (x) No () |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes (x) No () |

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Position: Lab DirectorPrinted Name: Susan C. SylvesterDate: 4-14-06

Lab Number: 10088-51
 Sample Designation: Method Blank
 Date Sampled: N/A
 Date Analyzed: 3/29/06
 Matrix: Water
 Instrument Dilution Factor: 1
 Analyst: LMM

VOLATILE ORGANICS
 SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dlbromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cls-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dlbromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 103 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 103 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10088-52
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\VOA03\032906\V3032924.D
Date Analyzed: 3/29/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-----------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 16 | 79% | 18 | 81% | 2% |
| chloromethane | 18 | 91% | 18 | 91% | 1% |
| vinyl chloride | 21 | 104% | 21 | 107% | 2% |
| bromomethane | 17 | 86% | 18 | 90% | 4% |
| chloroethane | 19 | 97% | 19 | 97% | 0% |
| trichlorofluoromethane | 20 | 99% | 20 | 101% | 2% |
| diethyl ether | 21 | 106% | 20 | 100% | 6% |
| acetone | 18 | 88% | 17 | 84% | 4% |
| 1,1-dichloroethene | 20 | 101% | 20 | 102% | 1% |
| methylene chloride | 21 | 107% | 21 | 107% | 0% |
| carbon disulfide | 18 | 92% | 19 | 95% | 3% |
| methyl-t-butyl ether (MTBE) | 41 | 104% | 41 | 103% | 1% |
| trans-1,2-dichloroethene | 20 | 100% | 20 | 102% | 2% |
| isopropyl ether (DIPE) | 21 | 106% | 21 | 106% | 1% |
| ethyl-t-butyl ether (ETBE) | 22 | 109% | 21 | 106% | 2% |
| 1,1-dichloroethane | 20 | 101% | 20 | 101% | 1% |
| t-butanol (TBA) | 92 | 92% | 90 | 90% | 2% |
| 2-butanone (MEK) | 19 | 93% | 17 | 87% | 6% |
| 2,2-dichloropropane | 18 | 89% | 18 | 90% | 0% |
| cis-1,2-dichloroethene | 22 | 109% | 22 | 108% | 2% |
| chloroform | 21 | 106% | 21 | 106% | 0% |
| bromochloromethane | 22 | 112% | 22 | 110% | 2% |
| tetrahydrofuran (THF) | 18 | 92% | 18 | 89% | 3% |
| 1,1,1-trichloroethane | 19 | 95% | 19 | 96% | 1% |
| 1,1-dichloropropene | 22 | 108% | 21 | 106% | 1% |
| t-amyl-methyl ether (TAME) | 21 | 106% | 21 | 103% | 2% |
| carbon tetrachloride | 15 | 77% | 17 | 83% | 7% |
| 1,2-dichloroethane | 20 | 101% | 20 | 100% | 1% |
| benzene | 21 | 105% | 21 | 107% | 2% |
| trichloroethene | 22 | 109% | 20 | 102% | 7% |
| 1,2-dichloropropane | 21 | 107% | 20 | 100% | 7% |
| bromodichloromethane | 17 | 84% | 17 | 86% | 2% |
| dibromomethane | 22 | 108% | 21 | 105% | 2% |
| 4-methyl-2-pentanone (MIBK) | 20 | 101% | 19 | 96% | 5% |
| cis-1,3-dichloropropene | 19 | 97% | 20 | 99% | 2% |
| toluene | 22 | 108% | 22 | 108% | 0% |
| trans-1,3-dichloropropene | 18 | 89% | 18 | 89% | 0% |
| 2-hexanone | 18 | 92% | 17 | 87% | 6% |
| 1,1,2-trichloroethane | 22 | 108% | 22 | 108% | 0% |
| 1,3-dichloropropane | 20 | 102% | 20 | 99% | 3% |
| tetrachloroethene | 21 | 105% | 20 | 102% | 2% |
| dibromochloromethane | 16 | 78% | 17 | 83% | 6% |
| 1,2-dibromoethane (EDB) | 20 | 98% | 20 | 100% | 2% |
| chlorobenzene | 20 | 99% | 19 | 97% | 2% |
| 1,1,1,2-tetrachloroethane | 17 | 86% | 18 | 91% | 5% |
| ethylbenzene | 20 | 102% | 21 | 103% | 0% |
| m&p-xylenes | 41 | 102% | 42 | 104% | 2% |
| o-xylene | 21 | 103% | 21 | 105% | 2% |
| styrene | 19 | 97% | 21 | 104% | 7% |
| bromoform | 15 | 74% | 16 | 79% | 7% |
| isopropylbenzene | 22 | 108% | 22 | 109% | 1% |
| 1,1,2,2-tetrachloroethane | 17 | 87% | 18 | 90% | 3% |
| 1,2,3-trichloropropane | 18 | 89% | 18 | 92% | 4% |
| n-propylbenzene | 20 | 99% | 21 | 105% | 5% |
| bromobenzene | 20 | 100% | 21 | 105% | 4% |
| 1,3,5-trimethylbenzene | 19 | 97% | 20 | 101% | 4% |
| 2-chlorotoluene | 19 | 93% | 20 | 98% | 5% |
| 4-chlorotoluene | 20 | 99% | 20 | 102% | 4% |
| tert-butylbenzene | 17 | 87% | 19 | 93% | 6% |
| 1,2,4-trimethylbenzene | 19 | 94% | 21 | 103% | 9% |
| sec-butylbenzene | 19 | 94% | 20 | 98% | 3% |
| 1,3-dichlorobenzene | 19 | 96% | 20 | 99% | 3% |
| 4-isopropyltoluene | 20 | 102% | 22 | 109% | 6% |
| 1,4-dichlorobenzene | 19 | 93% | 19 | 97% | 4% |
| 1,2-dichlorobenzene | 19 | 96% | 20 | 100% | 4% |
| n-butylbenzene | 21 | 106% | 22 | 111% | 5% |
| 1,2-dibromo-3-chloropropane | 16 | 80% | 18 | 88% | 9% |
| 1,2,4-trichlorobenzene | 20 | 102% | 21 | 107% | 5% |
| hexachlorobutadiene | 17 | 84% | 19 | 93% | 10% |
| naphthalene | 20 | 101% | 20 | 102% | 1% |
| 1,2,3-trichlorobenzene | 20 | 99% | 21 | 105% | 6% |
| 1,4-dioxene | 36 | 91% | 34 | 85% | 7% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 95% | 97% |
| SS toluene-D8 | 106% | 103% |
| SS 4-bromofluorobenzene | 103% | 104% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

Resource Laboratories, LLC

124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

10088

PAGE 1 OF 1

Company Name: **ES&M** Phone #: **508 285 9700**

Company Address: **184 West Main St. Norton** FAX #: **508 285 9757**

Project Manager: **Joe Callahan** Site Location (City, State): **Norton, MA**

Invoice To: **ES&M** Project ID / Name: **Leads Clean 2004-301**

Protocol: ☒ RCRA ☐ SDWA ☐ NPDES
☒ MCP ☐ NHDES ☐ OTHER

| Lab Sample ID | Field ID | # CONTAINERS | Matrix | | | Preservation Method | | | | | | Sampling | | SAMPLER |
|----------------|---------------|--------------|--------|-------|-------|---------------------|------------------|--------------------------------|------|------|-----------------|----------|------|---------|
| (Lab Use Only) | | | WATER | SOLID | OTHER | HCl | HNO ₃ | H ₂ SO ₄ | NaOH | MeOH | OTHER (Specify) | DATE | TIME | |
| 10088-01 | ESM05B(3x40) | 2 | ✓ | | | ✓ | | | | | | 3/28/06 | 0800 | AK |
| 02 | ESM05B(40-50) | 2 | ✓ | | | ✓ | | | | | | 3/28/06 | 1200 | AK |
| 03 | ESM05B(50-60) | 2 | ✓ | | | ✓ | | | | | | 3/28/06 | 1500 | AK |

- ☐ VOC 8260-NH List ☐ MADEP VPH ☐ MEGRO
☒ VOC 8260 ☐ VOC 80156RO ☐ VOC 624
☐ VOC 8260 BTEX, NMBE, Naphthalene only
☐ VOC 524.2 ☐ VOC 524.2 NH List
☐ TPH Fingerprint ☐ MEDRO ☐ DRO 8015 ☐ EPH
☐ 8270PAH ☐ 8270ABN ☐ 625
☐ 8082 PCB ☐ 8081 Pesticides ☐ 608
☐ O&G 1664 ☐ O&G SM5520F
☐ pH ☐ BOD ☐ Conductivity
☐ TSS ☐ TDS ☐ TS
☐ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals
☐ Total Metals-list ☐ Dissolved Metals-list
☐ Ammonia ☐ COD
☐ T-Phosphate ☐ Phenol
☐ Cyanide ☐ Sulfide
☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride
☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitibility/FP
☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC
☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)
☐ Standard Drinking Water Test ☐ Bacteria P/A

Grab (G) or Composite (C)

TAT REQUESTED

Priority (24 hr) ☐
Expedited (48 hr) ☐
10 Business Days ☐
Other ☒ Quote # _____
E-Mail Address _____
PO # _____

SPECIAL INSTRUCTIONS

REPORTING INSTRUCTIONS
☒ FAX ☐ OTHER (specify) _____
☐ PDF ☐ Excel Spreadsheet

RECEIVED ON ICE ☒ YES ☐ NO
 TEMPERATURE 6 °C

Lab Use Only

CUSTODY RECORD

| | | | | | |
|--------------------------|---------|------|-------------------------|---------|------|
| Relinquished by Sampler: | Date | Time | Received by: | Date | Time |
| <i>Joe Callahan</i> | 3/28/06 | 1515 | <i>Joe Callahan</i> | 3/28/06 | 1515 |
| Relinquished by: | Date | Time | Received by: | Date | Time |
| | | | | | |
| Relinquished by: | Date | Time | Received by Laboratory: | Date | Time |
| | | | | | |

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10094
Date Received: 3/29/06

Project: 2004-301 Lewis Chemical

Attached please find results for the analysis of the samples received on the date referenced above.

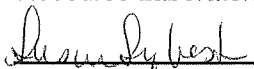
This is a reissue of report 10061, including the associated batch QC, case narrative and MCP Report Certification Form.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

2-2-07

Date

Total number of pages

Resource Laboratories, LLC Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Lab Number: 10094-01
 Sample Designation: ESM-05B (60-70)
 Date Sampled: 3/29/06
 Date Analyzed: 3/30/06
 Matrix: Water
 Instrument Dilution Factor: 20
 Analyst: LMM

VOLATILE ORGANICS
 SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 40 | trans-1,3-dichloropropene | U | 40 |
| chloromethane | U | 40 | 2-hexanone | U | 200 |
| vinyl chloride | U | 40 | 1,1,2-trichloroethane | U | 40 |
| bromomethane | U | 40 | 1,3-dichloropropane | U | 40 |
| chloroethane | U | 40 | tetrachloroethene | 3100 | 40 |
| trichlorofluoromethane | U | 40 | dibromochloromethane | U | 40 |
| diethyl ether | U | 200 | 1,2-dibromoethane | U | 40 |
| acetone | U | 200 | chlorobenzene | U | 40 |
| 1,1-dichloroethene | 220 | 20 | 1,1,1,2-tetrachloroethane | U | 40 |
| methylene chloride | 1000 | 100 | ethylbenzene | U | 40 |
| carbon disulfide | U | 40 | m&p-xylenes | 89 | 40 |
| methyl t-butyl ether (MTBE) | U | 40 | o-xylene | U | 40 |
| trans-1,2-dichloroethene | U | 40 | styrene | U | 40 |
| isopropyl ether (DIPE) | U | 40 | bromoform | U | 40 |
| ethyl t-butyl ether (ETBE) | U | 40 | isopropylbenzene | U | 40 |
| 1,1-dichloroethane | U | 40 | 1,1,2,2-tetrachloroethane | U | 40 |
| t-butanol (TBA) | U | 1000 | 1,2,3-trichloropropane | U | 40 |
| 2-butanone (MEK) | U | 200 | n-propylbenzene | U | 40 |
| 2,2-dichloropropane | U | 40 | bromobenzene | U | 40 |
| cis-1,2-dichloroethene | U | 40 | 1,3,5-trimethylbenzene | U | 40 |
| chloroform | U | 40 | 2-chlorotoluene | U | 40 |
| bromochloromethane | U | 40 | 4-chlorotoluene | U | 40 |
| tetrahydrofuran (THF) | U | 200 | tert-butylbenzene | U | 40 |
| 1,1,1-trichloroethane | 6800 | 40 | 1,2,4-trimethylbenzene | U | 40 |
| 1,1-dichloropropene | U | 40 | sec-butylbenzene | U | 40 |
| t-amyl-methyl ether (TAME) | U | 40 | 1,3-dichlorobenzene | U | 40 |
| carbon tetrachloride | U | 40 | 4-isopropyltoluene | U | 40 |
| 1,2-dichloroethane | 120 | 40 | 1,4-dichlorobenzene | U | 40 |
| benzene | U | 40 | 1,2-dichlorobenzene | 41 | 40 |
| trichloroethene | 7100 | 40 | n-butylbenzene | U | 40 |
| 1,2-dichloropropane | U | 40 | 1,2-dibromo-3-chloropropane | U | 40 |
| bromodichloromethane | U | 40 | 1,2,4-trichlorobenzene | U | 40 |
| dibromomethane | U | 40 | hexachlorobutadiene | U | 40 |
| 4-methyl-2-pentanone (MIBK) | U | 200 | naphthalene | U | 100 |
| cis-1,3-dichloropropene | U | 40 | 1,2,3-trichlorobenzene | U | 40 |
| toluene | 650 | 40 | 1,4-dioxane | U | 1000 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 95 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 105 | 86-115 | | | |

U = Below quantitation limit

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10094

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 6 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

No exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10094

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|----------|----------|---------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |

| | | |
|---|---|----------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
|---|---|----------------|

| | | |
|---|---|----------------|
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
|---|---|----------------|


| | | |
|---|--|----------------|
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
|---|--|----------------|

| | | |
|---|---|-------------------|
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No () NA |
|---|---|-------------------|

| | | |
|---|---|----------------|
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes (x) No () |
|---|---|----------------|

| | | |
|---|--|----------------|
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes (x) No () |
|---|--|----------------|

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Position: Lab DirectorPrinted Name: Susan C. SylvesterDate: 2-2-07

Lab Number: 10094-50
Sample Designation: Method Blank
Date Sampled: N/A
Date Analyzed: 3/30/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| dibromofluoromethane | 101 | 78-114 |
| toluene-D8 | 104 | 88-110 |
| 4-bromofluorobenzene | 103 | 86-115 |

U = Below quantitation limit

Lab Number: 10094-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\VOA03\033006\V3033024.D
Date Analyzed: 3/30/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-----------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 17 | 84% | 16 | 82% | 2% |
| chloromethane | 19 | 93% | 18 | 92% | 0% |
| vinyl chloride | 22 | 111% | 21 | 106% | 4% |
| bromomethane | 18 | 90% | 18 | 89% | 1% |
| chloroethane | 19 | 96% | 19 | 93% | 3% |
| trichlorofluoromethane | 20 | 102% | 20 | 101% | 1% |
| diethylether | 20 | 102% | 21 | 103% | 1% |
| acetone | 17 | 83% | 16 | 82% | 1% |
| 1,1-dichloroethene | 20 | 100% | 20 | 99% | 1% |
| methylene chloride | 22 | 108% | 21 | 107% | 0% |
| carbon disulfide | 19 | 97% | 19 | 84% | 2% |
| methyl-t-butyl ether (MTBE) | 39 | 99% | 39 | 98% | 1% |
| trans-1,2-dichloroethene | 21 | 104% | 20 | 101% | 3% |
| isopropyl ether (DIPE) | 21 | 104% | 21 | 105% | 1% |
| ethyl-t-butyl ether (ETBE) | 21 | 104% | 21 | 103% | 2% |
| 1,1-dichloroethane | 20 | 99% | 19 | 97% | 1% |
| t-butanol (TBA) | 86 | 86% | 89 | 89% | 3% |
| 2-butanone (MEK) | 18 | 89% | 18 | 90% | 0% |
| 2,2-dichloropropane | 19 | 94% | 19 | 94% | 1% |
| cis-1,2-dichloroethene | 21 | 107% | 22 | 108% | 1% |
| chloroform | 21 | 106% | 21 | 103% | 3% |
| bromochloromethane | 22 | 108% | 21 | 107% | 1% |
| tetrahydrofuran (THF) | 17 | 87% | 17 | 87% | 0% |
| 1,1,1-trichloroethane | 20 | 101% | 20 | 100% | 1% |
| 1,1-dichloropropene | 21 | 106% | 21 | 107% | 1% |
| t-amyl-methyl ether (TAME) | 20 | 101% | 20 | 101% | 0% |
| carbon tetrachloride | 17 | 85% | 17 | 87% | 2% |
| 1,2-dichloroethane | 20 | 98% | 20 | 100% | 2% |
| benzene | 21 | 103% | 21 | 106% | 2% |
| trichloroethene | 21 | 105% | 21 | 106% | 0% |
| 1,2-dichloropropane | 21 | 104% | 21 | 103% | 1% |
| bromodichloromethane | 18 | 88% | 18 | 91% | 4% |
| dibromomethane | 21 | 106% | 21 | 107% | 0% |
| 4-methyl-2-pentanone (MIBK) | 20 | 98% | 19 | 97% | 2% |
| cis-1,3-dichloropropene | 20 | 102% | 20 | 99% | 3% |
| toluene | 21 | 105% | 21 | 107% | 2% |
| trans-1,3-dichloropropene | 18 | 88% | 18 | 92% | 5% |
| 2-hexanone | 17 | 86% | 17 | 87% | 1% |
| 1,1,2-trichloroethane | 21 | 107% | 21 | 105% | 2% |
| 1,3-dichloropropane | 19 | 96% | 20 | 99% | 3% |
| tetrachloroethene | 20 | 100% | 21 | 105% | 5% |
| dibromochloromethane | 17 | 86% | 18 | 88% | 2% |
| 1,2-dibromoethane (EDB) | 19 | 97% | 20 | 99% | 3% |
| chlorobenzene | 19 | 96% | 20 | 100% | 4% |
| 1,1,1,2-tetrachloroethane | 18 | 92% | 19 | 94% | 2% |
| ethylbenzene | 20 | 99% | 20 | 101% | 2% |
| m&p-xylenes | 41 | 102% | 40 | 101% | 1% |
| o-xylene | 20 | 100% | 21 | 103% | 3% |
| styrene | 19 | 94% | 20 | 102% | 9% |
| bromoform | 17 | 84% | 17 | 86% | 3% |
| isopropylbenzene | 21 | 106% | 22 | 109% | 3% |
| 1,1,2,2-tetrachloroethane | 17 | 85% | 17 | 85% | 0% |
| 1,2,3-trichloropropane | 17 | 84% | 17 | 86% | 2% |
| n-propylbenzene | 20 | 99% | 20 | 98% | 1% |
| bromobenzene | 20 | 98% | 20 | 98% | 0% |
| 1,3,5-trimethylbenzene | 19 | 94% | 19 | 95% | 1% |
| 2-chlorotoluene | 19 | 95% | 18 | 90% | 5% |
| 4-chlorotoluene | 19 | 93% | 19 | 97% | 4% |
| tert-butylbenzene | 18 | 88% | 18 | 89% | 1% |
| 1,2,4-trimethylbenzene | 19 | 95% | 19 | 96% | 1% |
| sec-butylbenzene | 19 | 93% | 19 | 96% | 3% |
| 1,3-dichlorobenzene | 19 | 94% | 19 | 94% | 0% |
| 4-isopropyltoluene | 20 | 101% | 21 | 103% | 3% |
| 1,4-dichlorobenzene | 18 | 89% | 18 | 91% | 3% |
| 1,2-dichlorobenzene | 18 | 92% | 19 | 97% | 5% |
| n-butylbenzene | 21 | 103% | 21 | 107% | 3% |
| 1,2-dibromo-3-chloropropane | 16 | 82% | 17 | 87% | 6% |
| 1,2,4-trichlorobenzene | 18 | 92% | 19 | 97% | 4% |
| hexachlorobutadiene | 17 | 83% | 18 | 91% | 10% |
| naphthalene | 16 | 80% | 17 | 83% | 4% |
| 1,2,3-trichlorobenzene | 19 | 93% | 19 | 94% | 1% |
| 1,4-dioxane | 35 | 87% | 37 | 92% | 6% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 100% | 99% |
| SS toluene-D8 | 106% | 104% |
| SS 4-bromofluorobenzene | 103% | 104% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

RI

Resource Laboratories, LLC
124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST

ANALYSIS REQUEST

10094

PAGE 1 OF 1

Company Name:

Phone #: 508 285 9700

Company Address:

FAX #: 508 285 9457

Project Manager:

Site Location (City, State):

Environmental Strategies, LLC
184 West Main St North MA
Joe Callahan

Hyde Park MA
Project ID / Name:
2004-301 / Lewis Chemical

Invoice To:

Protocol:

ESM

RCRA

SDWA

NPDES

NHDES

OTHER

Lab Sample ID

(Lab Use Only)

Field ID

CONTAINERS

MATRIX

PRESERVATION METHOD

SAMPLING

SAMPLER

10094-01 ESM 053 (60-76)

2

WATER

SOLID

OTHER

HCl

HNO₃

H₂SO₄

NaOH

MeOH

OTHER (Specify)

DATE

TIME

3/29/06 11:30

AF

☐ VOC 8260-NH List ☐ MADEP VPH ☐ MEGRO

☒ VOC 8260 ☐ VOC 8015GRO ☐ VOC 624

☐ VOC 8260 BTEX, MIBE, Naphthalene only

☐ VOC 524.2 ☐ VOC 524.2 NH List

☐ TPH Fingerprint ☐ MEDRO ☐ DRO 8015 ☐ EPH

☐ 8270PAH ☐ 8270ABN ☐ 625

☐ 8082 PCB ☐ 8081 Pesticides ☐ 608

☐ O&G 1664 ☐ O&G SM5520F

☐ pH ☐ BOD ☐ Conductivity

☐ TSS ☐ TDS ☐ TS

☐ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals

☐ Total Metals-list ☐ Dissolved Metals-list

☐ Ammonia ☐ COD

☐ T-Phosphate ☐ Phenol

☐ Cyanide ☐ Sulfide

☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride

☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitibility/FP

☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC

☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)

☐ Standard Drinking Water Test ☐ Bacteria P/A

Grab (G) or Composite (C)

TAT REQUESTED

Priority (24 hr) ☐

Expedited (48 hr) ☐

10 Business Days ☐

Other ☐

ASAP

E-Mail Address

Quote #

PO #

SPECIAL INSTRUCTIONS

REPORTING INSTRUCTIONS

☐ FAX ☐ OTHER (specify)

☐ PDF ☐ Excel Spreadsheet

CUSTODY RECORD

Relinquished by Sampler:

Relinquished by:

Date

Time

Received by:

Received by:

Received by Laboratory:

Way Bill#:

Date

Time

Date

Time

Date

Time

Call Joe Callahan w/ results

RECEIVED ON ICE ☒ YES ☐ NO

TEMPERATURE

60 °C

Lab Use Only

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10103
Date Received: 3/30/06

Project: 2004-301 Lewis Chemical

Attached please find results for the analysis of the samples received on the date referenced above.

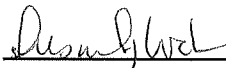
This is a reissue of report 10061, including the associated batch QC, case narrative and MCP Report Certification Form.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

2-2-07

Date

Total number of pages

9

Resource Laboratories, LLC Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Lab Number: 10103-01
Sample Designation: ESM-03B (50-60)
Date Sampled: 3/30/06
Date Analyzed: 4/1/06
Matrix: Water
Instrument Dilution Factor: 500
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 1000 | trans-1,3-dichloropropene | U | 1000 |
| chloromethane | U | 1000 | 2-hexanone | U | 5000 |
| vinyl chloride | U | 1000 | 1,1,2-trichloroethane | U | 1000 |
| bromomethane | U | 1000 | 1,3-dichloropropane | U | 1000 |
| chloroethane | U | 1000 | tetrachloroethene | 13000 | 1000 |
| trichlorofluoromethane | U | 1000 | dibromochloromethane | U | 1000 |
| diethyl ether | U | 5000 | 1,2-dibromoethane | U | 1000 |
| acetone | U | 5000 | chlorobenzene | U | 1000 |
| 1,1-dichloroethene | 3500 | 500 | 1,1,1,2-tetrachloroethane | U | 1000 |
| methylene chloride | 12000 | 3000 | ethylbenzene | U | 1000 |
| carbon disulfide | U | 1000 | m&p-xylenes | U | 1000 |
| methyl t-butyl ether (MTBE) | U | 1000 | o-xylene | U | 1000 |
| trans-1,2-dichloroethene | U | 1000 | styrene | U | 1000 |
| isopropyl ether (DIPE) | U | 1000 | bromoform | U | 1000 |
| ethyl t-butyl ether (ETBE) | U | 1000 | isopropylbenzene | U | 1000 |
| 1,1-dichloroethane | U | 1000 | 1,1,2,2-tetrachloroethane | U | 1000 |
| t-butanol (TBA) | U | 30000 | 1,2,3-trichloropropane | U | 1000 |
| 2-butanone (MEK) | U | 5000 | n-propylbenzene | U | 1000 |
| 2,2-dichloropropane | U | 1000 | bromobenzene | U | 1000 |
| cis-1,2-dichloroethene | U | 1000 | 1,3,5-trimethylbenzene | U | 1000 |
| chloroform | U | 1000 | 2-chlorotoluene | U | 1000 |
| bromochloromethane | U | 1000 | 4-chlorotoluene | U | 1000 |
| tetrahydrofuran (THF) | U | 5000 | tert-butylbenzene | U | 1000 |
| 1,1,1-trichloroethane | 24000 | 1000 | 1,2,4-trimethylbenzene | U | 1000 |
| 1,1-dichloropropene | U | 1000 | sec-butylbenzene | U | 1000 |
| t-amyl-methyl ether (TAME) | U | 1000 | 1,3-dichlorobenzene | U | 1000 |
| carbon tetrachloride | U | 1000 | 4-isopropyltoluene | U | 1000 |
| 1,2-dichloroethane | 1700 | 1000 | 1,4-dichlorobenzene | U | 1000 |
| benzene | U | 1000 | 1,2-dichlorobenzene | U | 1000 |
| trichloroethene | 58000 | 1000 | n-butylbenzene | U | 1000 |
| 1,2-dichloropropane | U | 1000 | 1,2-dibromo-3-chloropropane | U | 1000 |
| bromodichloromethane | U | 1000 | 1,2,4-trichlorobenzene | U | 1000 |
| dibromomethane | U | 1000 | hexachlorobutadiene | U | 1000 |
| 4-methyl-2-pentanone (MIBK) | U | 5000 | naphthalene | U | 3000 |
| cis-1,3-dichloropropene | U | 1000 | 1,2,3-trichlorobenzene | U | 1000 |
| toluene | 2000 | 1000 | 1,4-dioxane | U | 30000 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 101 | 78-114 | | | |
| toluene-D8 | 108 | 88-110 | | | |
| 4-bromofluorobenzene | 102 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10103-02
Sample Designation: ESM-03B (60-70)
Date Sampled: 3/30/06
Date Analyzed: 4/1/06
Matrix: Water
Instrument Dilution Factor: 500
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 1000 | trans-1,3-dichloropropene | U | 1000 |
| chloromethane | U | 1000 | 2-hexanone | U | 5000 |
| vinyl chloride | U | 1000 | 1,1,2-trichloroethane | U | 1000 |
| bromomethane | U | 1000 | 1,3-dichloropropane | U | 1000 |
| chloroethane | U | 1000 | tetrachloroethene | 20000 | 1000 |
| trichlorofluoromethane | U | 1000 | dibromochloromethane | U | 1000 |
| diethyl ether | U | 5000 | 1,2-dibromoethane | U | 1000 |
| acetone | U | 5000 | chlorobenzene | U | 1000 |
| 1,1-dichloroethene | 4900 | 500 | 1,1,1,2-tetrachloroethane | U | 1000 |
| methylene chloride | 29000 | 3000 | ethylbenzene | U | 1000 |
| carbon disulfide | U | 1000 | m&p-xylenes | U | 1000 |
| methyl t-butyl ether (MTBE) | U | 1000 | o-xylene | U | 1000 |
| trans-1,2-dichloroethene | U | 1000 | styrene | U | 1000 |
| isopropyl ether (DIPE) | U | 1000 | bromoform | U | 1000 |
| ethyl t-butyl ether (ETBE) | U | 1000 | isopropylbenzene | U | 1000 |
| 1,1-dichloroethane | 1000 | 1000 | 1,1,2,2-tetrachloroethane | U | 1000 |
| t-butanol (TBA) | U | 30000 | 1,2,3-trichloropropane | U | 1000 |
| 2-butanone (MEK) | U | 5000 | n-propylbenzene | U | 1000 |
| 2,2-dichloropropane | U | 1000 | bromobenzene | U | 1000 |
| cis-1,2-dichloroethene | U | 1000 | 1,3,5-trimethylbenzene | U | 1000 |
| chloroform | U | 1000 | 2-chlorotoluene | U | 1000 |
| bromochloromethane | U | 1000 | 4-chlorotoluene | U | 1000 |
| tetrahydrofuran (THF) | U | 5000 | tert-butylbenzene | U | 1000 |
| 1,1,1-trichloroethane | 49000 | 1000 | 1,2,4-trimethylbenzene | U | 1000 |
| 1,1-dichloropropene | U | 1000 | sec-butylbenzene | U | 1000 |
| t-amyl-methyl ether (TAME) | U | 1000 | 1,3-dichlorobenzene | U | 1000 |
| carbon tetrachloride | U | 1000 | 4-isopropyltoluene | U | 1000 |
| 1,2-dichloroethane | 3900 | 1000 | 1,4-dichlorobenzene | U | 1000 |
| benzene | U | 1000 | 1,2-dichlorobenzene | U | 1000 |
| trichloroethene | 100000 | 1000 | n-butylbenzene | U | 1000 |
| 1,2-dichloropropane | U | 1000 | 1,2-dibromo-3-chloropropane | U | 1000 |
| bromodichloromethane | U | 1000 | 1,2,4-trichlorobenzene | U | 1000 |
| dibromomethane | U | 1000 | hexachlorobutadiene | U | 1000 |
| 4-methyl-2-pentanone (MIBK) | U | 5000 | naphthalene | U | 3000 |
| cis-1,3-dichloropropene | U | 1000 | 1,2,3-trichlorobenzene | U | 1000 |
| toluene | 6600 | 1000 | 1,4-dioxane | U | 30000 |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 100 | 78-114 | | | |
| toluene-D8 | 105 | 88-110 | | | |
| 4-bromofluorobenzene | 100 | 86-115 | | | |

U = Below quantitation limit

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10103

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 10 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

The RPD between the LCS and LCSD for Styrene was above the acceptance limit. The recoveries for both were acceptable. No data impact expected.

No exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10103

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|------------|----------|---------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |

| | | |
|---|---|---------------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|---|---------------------|
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|--|---------------------|
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
|---|--|---------------------|

| | | |
|---|---|----------------------|
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No () NA |
|---|---|----------------------|

| | | |
|---|---|---------------------|
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes () No (x) |
|---|---|---------------------|

| | | |
|---|--|---------------------|
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes (x) No () |
|---|--|---------------------|

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Position: Lab DirectorPrinted Name: Susan C. SylvesterDate: 2-2-07

Lab Number: 10103-50
Sample Designation: Method Blank
Date Sampled: N/A
Date Analyzed: 3/31/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dlbromofluoromethane | 101 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 98 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10103-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\VOA03\033106\W3033134.D
Date Analyzed: 4/1/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-------------------------------|--------------|------------|--------------|------------|-------|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 16 | 81% | 16 | 80% | 1% |
| chloromethane | 18 | 91% | 18 | 90% | 1% |
| vinyl chloride | 20 | 100% | 21 | 105% | 5% |
| bromomethane | 17 | 83% | 18 | 88% | 5% |
| chloroethane | 19 | 96% | 19 | 94% | 3% |
| trichlorofluoromethane | 20 | 100% | 21 | 103% | 3% |
| diethylether | 21 | 103% | 20 | 102% | 1% |
| acetone | 17 | 87% | 17 | 87% | 0% |
| 1,1-dichloroethene | 19 | 96% | 20 | 102% | 6% |
| methylene chloride | 21 | 106% | 22 | 108% | 2% |
| carbon disulfide | 19 | 93% | 19 | 94% | 1% |
| methyl-t-butyl ether (MTBE) | 40 | 100% | 41 | 102% | 2% |
| trans-1,2-dichloroethene | 21 | 104% | 21 | 104% | 0% |
| isopropyl ether (DIPE) | 21 | 106% | 22 | 111% | 5% |
| ethyl-t-butyl ether (ETBE) | 21 | 103% | 21 | 106% | 3% |
| 1,1-dichloroethane | 19 | 97% | 20 | 102% | 5% |
| t-butanol (TBA) | 84 | 84% | 89 | 89% | 6% |
| 2-butanone (MEK) | 18 | 89% | 19 | 93% | 5% |
| 2,2-dichloropropane | 14 | 70% | 14 | 72% | 3% |
| cis-1,2-dichloroethene | 22 | 108% | 22 | 110% | 2% |
| chloroform | 21 | 104% | 22 | 108% | 3% |
| bromochloromethane | 21 | 107% | 23 | 113% | 5% |
| tetrahydrofuran (THF) | 18 | 90% | 19 | 94% | 5% |
| 1,1,1-trichloroethane | 20 | 99% | 21 | 103% | 4% |
| 1,1-dichloropropene | 21 | 105% | 20 | 102% | 3% |
| t-amyl-methyl ether (TAME) | 20 | 101% | 20 | 102% | 1% |
| carbon tetrachloride | 17 | 85% | 18 | 89% | 4% |
| 1,2-dichloroethane | 20 | 99% | 21 | 103% | 4% |
| benzene | 21 | 107% | 22 | 108% | 1% |
| trichloroethene | 22 | 108% | 22 | 108% | 0% |
| 1,2-dichloropropane | 21 | 104% | 20 | 102% | 2% |
| bromodichloromethane | 18 | 90% | 18 | 91% | 1% |
| dibromomethane | 22 | 108% | 22 | 111% | 2% |
| 4-methyl-2-pentanone (MIBK) | 18 | 92% | 20 | 99% | 7% |
| cis-1,3-dichloropropene | 19 | 96% | 20 | 99% | 3% |
| toluene | 22 | 108% | 21 | 106% | 2% |
| trans-1,3-dichloropropene | 17 | 85% | 18 | 88% | 3% |
| 2-hexanone | 16 | 82% | 17 | 87% | 6% |
| 1,1,2-trichloroethane | 21 | 103% | 21 | 104% | 1% |
| 1,3-dichloropropane | 19 | 97% | 21 | 104% | 7% |
| tetrachloroethene | 20 | 101% | 22 | 109% | 8% |
| dibromochloromethane | 18 | 89% | 19 | 96% | 8% |
| 1,2-dibromoethane (EDB) | 19 | 97% | 20 | 100% | 3% |
| chlorobenzene | 19 | 96% | 20 | 99% | 3% |
| 1,1,1,2-tetrachloroethane | 18 | 91% | 20 | 98% | 7% |
| ethylbenzene | 19 | 97% | 21 | 103% | 6% |
| m&p-xylenes | 40 | 101% | 41 | 104% | 3% |
| o-xylene | 20 | 100% | 20 | 102% | 3% |
| styrene | 15 | 74% | 20 | 101% | 31% + |
| bromoform | 17 | 87% | 19 | 93% | 7% |
| isopropylbenzene | 21 | 103% | 22 | 111% | 8% |
| 1,1,2,2-tetrachloroethane | 17 | 85% | 17 | 87% | 2% |
| 1,2,3-trichloropropane | 17 | 85% | 17 | 83% | 2% |
| n-propylbenzene | 19 | 97% | 20 | 98% | 1% |
| bromobenzene | 20 | 98% | 19 | 97% | 1% |
| 1,3,5-trimethylbenzene | 19 | 93% | 19 | 96% | 3% |
| 2-chlorotoluene | 18 | 92% | 19 | 94% | 2% |
| 4-chlorotoluene | 18 | 90% | 19 | 93% | 3% |
| tert-butylbenzene | 17 | 85% | 18 | 89% | 5% |
| 1,2,4-trimethylbenzene | 19 | 93% | 19 | 97% | 4% |
| sec-butylbenzene | 17 | 87% | 19 | 94% | 7% |
| 1,3-dichlorobenzene | 19 | 93% | 19 | 93% | 0% |
| 4-isopropyltoluene | 20 | 98% | 20 | 102% | 4% |
| 1,4-dichlorobenzene | 18 | 90% | 18 | 90% | 0% |
| 1,2-dichlorobenzene | 19 | 95% | 19 | 97% | 3% |
| n-butylbenzene | 19 | 97% | 20 | 101% | 4% |
| 1,2-dibromo-3-chloropropane (| 17 | 83% | 17 | 85% | 3% |
| 1,2,4-trichlorobenzene | 17 | 87% | 18 | 88% | 0% |
| hexachlorobutadiene | 17 | 83% | 17 | 86% | 4% |
| naphthalene | 14 | 71% | 15 | 77% | 7% |
| 1,2,3-trichlorobenzene | 17 | 87% | 18 | 91% | 4% |
| 1,4-dioxane | 35 | 88% | 36 | 89% | 2% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 99% | 101% |
| SS toluene-D8 | 106% | 108% |
| SS 4-bromofluorobenzene | 103% | 109% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

+ The RPD between LCS and LCS duplicate was above the acceptance limit. The recoveries for both were acceptable.

CHAIN-OF-CUSTODY RECORD 10103
AND ANALYSIS REQUEST
ANALYSIS REQUEST

Phone # (508) 285-4760

FAX #: (571) 285-9757

Hyde Park, WA

2004-301 / Lewis CHITARRAS

| | | | | |
|--|-----------|-----|-------|-------|
| | Protocol: | MCP | NHDES | OTHER |
|--|-----------|-----|-------|-------|

[illegible]

SPECIAL INSTRUCTIONS

REPORTING INSTRUCTIONS

☐ FAX ☐ OTHER (specify) _____☐ PDF ☐ Excel Spreadsheet

Date _____

| | |
|--|---------|
| | Date |
| | 3/30/06 |

| | |
|--|-------------|
| | U P C |
|--|-------------|

| | |
|--|------|
| | Date |
|--|------|

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

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10105
Date Received: 3/31/06

Project: 2004-301 Lewis Chemical

Attached please find results for the analysis of the samples received on the date referenced above.

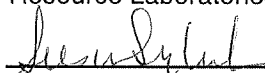
This is a reissue of report 10061, including the associated batch QC, case narrative and MCP Report Certification Form.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

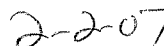
Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager



Date

Total number of pages _____

Resource Laboratories, LLC Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Lab Number: 10105-01
Sample Designation: ESM-08B(40-50)
Date Sampled: 3/31/06
Date Analyzed: 4/3/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | 250 | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | 2 | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | 2 | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | 5 | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | 40 | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | 6 | 2 |
| trichloroethene | 350 | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | 24 | 2 | 1,4-dioxane | U | 50 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 96 | 78-114 | | | |
| toluene-D8 | 108 | 88-110 | | | |
| 4-bromofluorobenzene | 106 | 86-115 | | | |

U = Below quantitation limit

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10105

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 12 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

No exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

A quadratic curve fit was used in the initial calibration for the following compounds: Dichlorodifluoromethane, Chloromethane, Vinyl Chloride, Bromomethane, Acetone, Dibromochloromethane, Isopropylbenzene, Tert-butylbenzene, and Hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10105

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|------------|----------|---------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |

| | | |
|---|---|---------------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|---|---------------------|
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|--|---------------------|
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
|---|--|---------------------|

| | | |
|---|---|----------------------|
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No () NA |
|---|---|----------------------|

| | | |
|---|---|---------------------|
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes (x) No () |
|---|---|---------------------|

| | | |
|---|--|---------------------|
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes (x) No () |
|---|--|---------------------|

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Susan C. SylvesterPosition: Lab DirectorPrinted Name: Susan C. SylvesterDate: 2-2-07

Lab Number: 10105-50
Sample Designation: Method Blank
Date Sampled: N/A
Date Analyzed: 4/3/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 101 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 100 | 86-115 | | | |

U = Below quantitation limit.

Lab Number: 10106-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\W\OA03\04\0306\W3040324.D
Date Analyzed: 4/3/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-----------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 18 | 88% | 17 | 86% | 2% |
| chloromethane | 20 | 100% | 19 | 94% | 5% |
| vinyl chloride | 22 | 112% | 21 | 104% | 7% |
| bromomethane | 20 | 99% | 19 | 95% | 4% |
| chloroethane | 21 | 103% | 20 | 100% | 3% |
| trichlorofluoromethane | 22 | 111% | 21 | 104% | 7% |
| diethylether | 22 | 109% | 22 | 110% | 0% |
| acetone | 18 | 88% | 18 | 89% | 1% |
| 1,1-dichloroethene | 22 | 111% | 21 | 106% | 5% |
| methylene chloride | 23 | 114% | 22 | 111% | 2% |
| carbon disulfide | 21 | 104% | 20 | 98% | 6% |
| methyl-t-butyl ether (MTBE) | 43 | 108% | 42 | 104% | 4% |
| trans-1,2-dichloroethene | 23 | 114% | 22 | 109% | 5% |
| isopropyl ether (DIPE) | 23 | 116% | 22 | 112% | 4% |
| ethyl-t-butyl ether (ETBE) | 23 | 113% | 22 | 111% | 2% |
| 1,1-dichloroethane | 22 | 108% | 21 | 106% | 2% |
| t-butanol (TBA) | 95 | 95% | 94 | 94% | 2% |
| 2-butanone (MEK) | 19 | 96% | 19 | 96% | 0% |
| 2,2-dichloropropane | 21 | 103% | 20 | 99% | 4% |
| cis-1,2-dichloroethene | 24 | 118% | 23 | 115% | 3% |
| chloroform | 23 | 117% | 23 | 115% | 1% |
| bromochloromethane | 24 | 120% | 24 | 121% | 1% |
| tetrahydrofuran (THF) | 19 | 94% | 19 | 96% | 3% |
| 1,1,1-trichloroethane | 22 | 112% | 22 | 110% | 2% |
| 1,1-dichloropropene | 23 | 117% | 23 | 113% | 3% |
| t-amyl-methyl ether (TAME) | 22 | 110% | 23 | 115% | 4% |
| carbon tetrachloride | 18 | 92% | 19 | 94% | 2% |
| 1,2-dichloroethane | 22 | 110% | 22 | 109% | 1% |
| benzene | 24 | 119% | 23 | 116% | 2% |
| trichloroethene | 23 | 114% | 23 | 115% | 1% |
| 1,2-dichloropropane | 22 | 109% | 22 | 110% | 2% |
| bromodichloromethane | 19 | 96% | 20 | 98% | 1% |
| dibromomethane | 24 | 118% | 23 | 115% | 2% |
| 4-methyl-2-pentanone (MIBK) | 21 | 103% | 21 | 103% | 0% |
| cis-1,3-dichloropropene | 22 | 109% | 22 | 110% | 0% |
| toluene | 24 | 120% | 23 | 116% | 3% |
| trans-1,3-dichloropropene | 19 | 97% | 19 | 96% | 1% |
| 2-hexanone | 19 | 93% | 19 | 95% | 3% |
| 1,1,2-trichloroethane | 23 | 114% | 23 | 115% | 1% |
| 1,3-dichloropropane | 21 | 104% | 20 | 99% | 5% |
| tetrachloroethene | 22 | 109% | 21 | 105% | 4% |
| dibromochloromethane | 18 | 89% | 18 | 92% | 3% |
| 1,2-dibromoethane (EDB) | 20 | 100% | 20 | 101% | 1% |
| chlorobenzene | 20 | 102% | 20 | 98% | 4% |
| 1,1,1,2-tetrachloroethane | 19 | 94% | 19 | 96% | 2% |
| ethylbenzene | 21 | 106% | 21 | 103% | 3% |
| m&p-xylenes | 43 | 107% | 42 | 106% | 0% |
| o-xylene | 22 | 108% | 21 | 106% | 2% |
| styrene | 21 | 103% | 19 | 97% | 6% |
| bromoform | 18 | 88% | 18 | 90% | 3% |
| isopropylbenzene | 23 | 114% | 22 | 110% | 4% |
| 1,1,2,2-tetrachloroethane | 18 | 92% | 17 | 87% | 6% |
| 1,2,3-trichloropropane | 18 | 89% | 17 | 86% | 4% |
| n-propylbenzene | 21 | 105% | 20 | 98% | 7% |
| bromobenzene | 21 | 104% | 20 | 101% | 4% |
| 1,3,5-trimethylbenzene | 19 | 97% | 18 | 92% | 5% |
| 2-chlorotoluene | 19 | 97% | 18 | 92% | 5% |
| 4-chlorotoluene | 20 | 99% | 19 | 94% | 5% |
| tert-butylbenzene | 18 | 92% | 18 | 90% | 2% |
| 1,2,4-trimethylbenzene | 20 | 101% | 20 | 99% | 2% |
| sec-butylbenzene | 20 | 101% | 18 | 91% | 10% |
| 1,3-dichlorobenzene | 20 | 99% | 19 | 95% | 4% |
| 4-isopropyltoluene | 22 | 108% | 21 | 104% | 4% |
| 1,4-dichlorobenzene | 19 | 96% | 19 | 93% | 3% |
| 1,2-dichlorobenzene | 21 | 103% | 19 | 97% | 5% |
| n-butylbenzene | 22 | 109% | 21 | 106% | 2% |
| 1,2-dibromo-3-chloropropane | 17 | 87% | 17 | 87% | 0% |
| 1,2,4-trichlorobenzene | 20 | 101% | 19 | 96% | 5% |
| hexachlorobutadiene | 19 | 94% | 19 | 93% | 1% |
| naphthalene | 17 | 85% | 17 | 85% | 0% |
| 1,2,3-trichlorobenzene | 20 | 101% | 20 | 100% | 1% |
| 1,4-dioxane | 37 | 92% | 36 | 90% | 3% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 104% | 102% |
| SS toluene-D8 | 107% | 110% |
| SS 4-bromofluorobenzene | 108% | 103% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

| Protocol: | MCB | NHDES | OTHER |
|----------------------|-----|-------|-------|
| 1. <i>Protocol</i> | | | |
| 2. <i>Protocol</i> | | | |
| 3. <i>Protocol</i> | | | |
| 4. <i>Protocol</i> | | | |
| 5. <i>Protocol</i> | | | |
| 6. <i>Protocol</i> | | | |
| 7. <i>Protocol</i> | | | |
| 8. <i>Protocol</i> | | | |
| 9. <i>Protocol</i> | | | |
| 10. <i>Protocol</i> | | | |
| 11. <i>Protocol</i> | | | |
| 12. <i>Protocol</i> | | | |
| 13. <i>Protocol</i> | | | |
| 14. <i>Protocol</i> | | | |
| 15. <i>Protocol</i> | | | |
| 16. <i>Protocol</i> | | | |
| 17. <i>Protocol</i> | | | |
| 18. <i>Protocol</i> | | | |
| 19. <i>Protocol</i> | | | |
| 20. <i>Protocol</i> | | | |
| 21. <i>Protocol</i> | | | |
| 22. <i>Protocol</i> | | | |
| 23. <i>Protocol</i> | | | |
| 24. <i>Protocol</i> | | | |
| 25. <i>Protocol</i> | | | |
| 26. <i>Protocol</i> | | | |
| 27. <i>Protocol</i> | | | |
| 28. <i>Protocol</i> | | | |
| 29. <i>Protocol</i> | | | |
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| <input type="checkbox"/> VOC 8260-NH List | <input type="checkbox"/> MADEP VPH | <input type="checkbox"/> MEGRO |
| <input checked="" type="checkbox"/> VOC 8260 | <input type="checkbox"/> VOC8015GRO | <input type="checkbox"/> VOC 624 |
| <input type="checkbox"/> VOC 8260 BTEX, MIBE, Naphthalene only | | |
| <input type="checkbox"/> VOC 524.2 | <input type="checkbox"/> VOC 524.2 NH List | |
| <input type="checkbox"/> TPH Fingerprint | <input type="checkbox"/> MEDRO | <input type="checkbox"/> DRO 8015 <input type="checkbox"/> EPH |
| <input type="checkbox"/> 8270PAH | <input type="checkbox"/> 8270ABN | <input type="checkbox"/> 625 |
| <input type="checkbox"/> 8082 PCB | <input type="checkbox"/> 8081 Pesticides | <input type="checkbox"/> 608 |
| <input type="checkbox"/> O&G 1664 | <input type="checkbox"/> O&G SM5520F | |
| <input type="checkbox"/> pH | <input type="checkbox"/> BOD | <input type="checkbox"/> Conductivity |
| <input type="checkbox"/> TSS | <input type="checkbox"/> TDS | <input type="checkbox"/> TS |
| <input type="checkbox"/> RCRA Metals | <input type="checkbox"/> Priority Pollutant Metals | <input type="checkbox"/> TAL Metals |
| <input type="checkbox"/> Total Metals-list | <input type="checkbox"/> Dissolved Metals-list | |
| <input type="checkbox"/> Ammonia | <input type="checkbox"/> COD | |
| <input type="checkbox"/> T-Phosphate | <input type="checkbox"/> Phenol | |
| <input type="checkbox"/> Cyanide | <input type="checkbox"/> Sulfide | |
| <input type="checkbox"/> Nitrate | <input type="checkbox"/> Nitrite | <input type="checkbox"/> Ortho P <input type="checkbox"/> Sulfate <input type="checkbox"/> Bromide <input type="checkbox"/> Chloride |
| <input type="checkbox"/> Corrosivity | <input type="checkbox"/> Reactive CN | <input type="checkbox"/> Reactive S- <input type="checkbox"/> Ignitibility/FP |
| <input type="checkbox"/> TCLP Metals | <input type="checkbox"/> TCLP VOC | <input type="checkbox"/> TCLP SVOC |
| <input type="checkbox"/> TCLP Pesticide | <input type="checkbox"/> TCLP Herbicides (subcontract) | |
| <input type="checkbox"/> Standard Drinking Water Test | <input type="checkbox"/> Bacteria P/A | |
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| Grab (G) or Composite (C) | | |

REPORTING INSTRUCTIONS

☐ FAX ☐ OTHER (specify) _____

☐ PDF ☐ Excel Spreadsheet

| | | | | | |
|--------------------------|---------|------|--|---------|---------|
| Relinquished by Sampler: | Date | Time | Received by: | Date | Time |
| <i>Andrew Mann</i> | 3/31/06 | 1200 | | | |
| Relinquished by: | Date | Time | Received by: | Date | Time |
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| Relinquished by: | Date | Time | Received by Laboratory Valerie Billar | Date | Time |
| | | | | 3/31/06 | 1:23 PM |

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10154
Date Received: 4/7/06

Project: 2006-056 DND Lewis

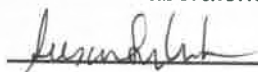
Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

4-27-06

Date

Total number of pages

71

Resource Laboratories, LLC Certifications

New Hampshire NH902
Maine NH903

Connecticut PH-0146
Massachusetts M-NH902

Lab Number: 10154-05
 Sample Designation: ESM14
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Instrument Dilution Factor: 1
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | 330 | 100 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | 330 | 100 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 2 |
| benzene | C5-C8 | U | 1 |
| toluene | C5-C8 | 11 | 2 |
| ethylbenzene | C9-C12 | 240 | 2 |
| m&p-xylenes | C9-C12 | 42 | 2 |
| o-xylene | C9-C12 | 12 | 2 |
| naphthalene | N/A | U | 5 |
| C5-C8 Aliphatics (1,2) | N/A | 320 | 100 |
| C9-C12 Aliphatics (1,3) | N/A | U | 100 |
| C9-C10 Aromatics (1) | N/A | U | 100 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 93% | |
| 2,5-dibromotoluene as aliphatic | | 95% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or Internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10154-06
 Sample Designation: DUP
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Instrument Dilution Factor: 1
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | 380 | 100 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | 390 | 100 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 2 |
| benzene | C5-C8 | U | 1 |
| toluene | C5-C8 | 14 | 2 |
| ethylbenzene | C9-C12 | 290 | 2 |
| m&p-xylenes | C9-C12 | 51 | 2 |
| o-xylene | C9-C12 | 14 | 2 |
| naphthalene | N/A | U | 5 |
| C5-C8 Aliphatics (1,2) | N/A | 360 | 100 |
| C9-C12 Aliphatics (1,3) | N/A | U | 100 |
| C9-C10 Aromatics (1) | N/A | U | 100 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 99% | |
| 2,5-dibromotoluene as aliphatic | | 102% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10154-07
 Sample Designation: ESM13
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH \leq 2
 Temperature: Received on Ice at 4 \pm 2°C
 Instrument Dilution Factor: 5
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | 4000 | 500 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | 6200 | 500 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 10 |
| benzene | C5-C8 | U | 5 |
| toluene | C5-C8 | 370 | 10 |
| ethylbenzene | C9-C12 | 1600 | 10 |
| m&p-xylenes | C9-C12 | 2200 | 10 |
| o-xylene | C9-C12 | 390 | 10 |
| naphthalene | N/A | U | 30 |
| C5-C8 Aliphatics (1,2) | N/A | 3600 | 500 |
| C9-C12 Aliphatics (1,3) | N/A | U | 500 |
| C9-C10 Aromatics (1) | N/A | 2100 | 500 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 106% | |
| 2,5-dibromotoluene as aliphatic | | 108% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10154-08
 Sample Designation: ESM-8B
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH \leq 2
 Temperature: Received on Ice at 4 \pm 2°C
 Instrument Dilution Factor: 1
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | U | 100 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | U | 100 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 2 |
| benzene | C5-C8 | U | 1 |
| toluene | C5-C8 | U | 2 |
| ethylbenzene | C9-C12 | U | 2 |
| m&p-xylenes | C9-C12 | U | 2 |
| o-xylene | C9-C12 | U | 2 |
| naphthalene | N/A | U | 5 |
| C5-C8 Aliphatics (1,2) | N/A | U | 100 |
| C9-C12 Aliphatics (1,3) | N/A | U | 100 |
| C9-C10 Aromatics (1) | N/A | U | 100 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 94% | |
| 2,5-dibromotoluene as aliphatic | | 96% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10154-09
 Sample Designation: ESM-8
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Instrument Dilution Factor: 1
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | 340 | 100 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | U | 100 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 2 |
| benzene | C5-C8 | U | 1 |
| toluene | C5-C8 | U | 2 |
| ethylbenzene | C9-C12 | U | 2 |
| m&p-xylenes | C9-C12 | U | 2 |
| o-xylene | C9-C12 | U | 2 |
| naphthalene | N/A | U | 5 |
| C5-C8 Aliphatics (1,2) | N/A | 340 | 100 |
| C9-C12 Aliphatics (1,3) | N/A | U | 100 |
| C9-C10 Aromatics (1) | N/A | U | 100 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 104% | |
| 2,5-dibromotoluene as aliphatic | | 103% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or Internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10154-10
 Sample Designation: ESM-3
 Date Sampled: 4/7/06
 Date Received: 4/7/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Instrument Dilution Factor: 100
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | 190000 | 10000 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | 14000 | 10000 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 200 |
| benzene | C5-C8 | U | 100 |
| toluene | C5-C8 | 40000 | 200 |
| ethylbenzene | C9-C12 | 1600 | 200 |
| m&p-xylenes | C9-C12 | 5300 | 200 |
| o-xylene | C9-C12 | 1400 | 200 |
| naphthalene | N/A | U | 500 |
| C5-C8 Aliphatics (1,2) | N/A | 150000 | 10000 |
| C9-C12 Aliphatics (1,3) | N/A | U | 10000 |
| C9-C10 Aromatics (1) | N/A | U | 10000 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 101% | |
| 2,5-dibromotoluene as aliphatic | | 103% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10154-01
Sample Designation: ESM11
Date Sampled: 4/6/06
Date Analyzed: 4/13/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | 2 | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | 5 | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 92 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 101 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-03
Sample Designation: ESM12
Date Sampled: 4/6/06
Date Analyzed: 4/13/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DiPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 89 | 78-114 | | | |
| toluene-D8 | 100 | 88-110 | | | |
| 4-bromofluorobenzene | 101 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-04
Sample Designation: ESM02
Date Sampled: 4/6/06
Date Analyzed: 4/13/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | 17 | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | 18 | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | 10 | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | 6 | 2 |
| trans-1,2-dichloroethene | 2 | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | 2 | 2 |
| 1,1-dichloroethane | 69 | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | 5 | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | 4 | 2 | 1,2,4-trimethylbenzene | 34 | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | 12 | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | 11 | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | 6 | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | 98 | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | 14 | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | 19 | 2 | 1,4-dioxane | U | 50 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 94 | 78-114 | | | |
| toluene-D8 | 102 | 88-110 | | | |
| 4-bromofluorobenzene | 114 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-05
Sample Designation: ESM14
Date Sampled: 4/6/06
Date Analyzed: 4/13/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | 20 | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | 8 | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | 2 | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | 270 | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | 44 | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | 13 | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | 42 | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | 15 | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | 8 | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | 12 | 2 | 1,4-dioxane | U | 50 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| dibromofluoromethane | 95 | 78-114 |
| toluene-D8 | 104 | 88-110 |
| 4-bromofluorobenzene | 101 | 86-115 |

U = Below quantitation limit

Lab Number: 10154-07
Sample Designation: ESM13
Date Sampled: 4/6/06
Date Analyzed: 4/12/06
Matrix: Water
Instrument Dilution Factor: 5
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 10 | trans-1,3-dichloropropene | U | 10 |
| chloromethane | U | 10 | 2-hexanone | U | 50 |
| vinyl chloride | 61 | 10 | 1,1,2-trichloroethane | U | 10 |
| bromomethane | U | 10 | 1,3-dichloropropane | U | 10 |
| chloroethane | U | 10 | tetrachloroethene | 150 | 10 |
| trichlorofluoromethane | U | 10 | dibromochloromethane | U | 10 |
| diethyl ether | U | 50 | 1,2-dibromoethane | U | 10 |
| acetone | U | 50 | chlorobenzene | U | 10 |
| 1,1-dichloroethene | U | 5 | 1,1,1,2-tetrachloroethane | U | 10 |
| methylene chloride | 38 | 30 | ethylbenzene | 1700 | 10 |
| carbon disulfide | U | 10 | m&p-xylenes | 2400 | 10 |
| methyl t-butyl ether (MTBE) | U | 10 | o-xylene | 370 | 10 |
| trans-1,2-dichloroethene | U | 10 | styrene | U | 10 |
| isopropyl ether (DIPE) | U | 10 | bromoform | U | 10 |
| ethyl t-butyl ether (ETBE) | U | 10 | isopropylbenzene | U | 10 |
| 1,1-dichloroethane | 25 | 10 | 1,1,2,2-tetrachloroethane | U | 10 |
| t-butanol (TBA) | U | 300 | 1,2,3-trichloropropane | U | 10 |
| 2-butanone (MEK) | U | 50 | n-propylbenzene | U | 10 |
| 2,2-dichloropropane | U | 10 | bromobenzene | U | 10 |
| cis-1,2-dichloroethene | 2200 | 10 | 1,3,5-trimethylbenzene | U | 10 |
| chloroform | U | 10 | 2-chlorotoluene | U | 10 |
| bromochloromethane | U | 10 | 4-chlorotoluene | U | 10 |
| tetrahydrofuran (THF) | U | 50 | tert-butylbenzene | U | 10 |
| 1,1,1-trichloroethane | 14 | 10 | 1,2,4-trimethylbenzene | 14 | 10 |
| 1,1-dichloropropene | U | 10 | sec-butylbenzene | 15 | 10 |
| t-amyl-methyl ether (TAME) | U | 10 | 1,3-dichlorobenzene | U | 10 |
| carbon tetrachloride | U | 10 | 4-isopropyltoluene | 130 | 10 |
| 1,2-dichloroethane | 20 | 10 | 1,4-dichlorobenzene | U | 10 |
| benzene | U | 10 | 1,2-dichlorobenzene | U | 10 |
| trichloroethene | 57 | 10 | n-butylbenzene | U | 10 |
| 1,2-dichloropropane | U | 10 | 1,2-dibromo-3-chloropropane | U | 10 |
| bromodichloromethane | U | 10 | 1,2,4-trichlorobenzene | U | 10 |
| dibromomethane | U | 10 | hexachlorobutadiene | U | 10 |
| 4-methyl-2-pentanone (MIBK) | U | 50 | naphthalene | U | 30 |
| cis-1,3-dichloropropene | U | 10 | 1,2,3-trichlorobenzene | U | 10 |
| toluene | 340 | 10 | 1,4-dioxane | U | 300 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 92 | 78-114 | | | |
| toluene-D8 | 101 | 88-110 | | | |
| 4-bromofluorobenzene | 106 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-08
 Sample Designation: ESM-8B
 Date Sampled: 4/6/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Instrument Dilution Factor: 1
 Analyst: CWC

VOLATILE ORGANICS
 SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | 6 | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | 2 | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | 7 | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 93 | 78-114 | | | |
| toluene-D8 | 103 | 88-110 | | | |
| 4-bromofluorobenzene | 105 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-09
Sample Designation: ESM-8
Date Sampled: 4/6/06
Date Analyzed: 4/13/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | 12 | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | 44 | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | 16 | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | 42 | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | 17 | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | 11 | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | 39 | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 91 | 78-114 | | | |
| toluene-DB | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 104 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-10
Sample Designation: ESM-3
Date Sampled: 4/7/06
Date Analyzed: 4/14/06
Matrix: Water
Instrument Dilution Factor: 200
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 400 | trans-1,3-dichloropropene | U | 400 |
| chloromethane | U | 400 | 2-hexanone | U | 2000 |
| vinyl chloride | U | 400 | 1,1,2-trichloroethane | U | 400 |
| bromomethane | U | 400 | 1,3-dichloropropane | U | 400 |
| chloroethane | U | 400 | tetrachloroethene | 4300 | 400 |
| trichlorofluoromethane | U | 400 | dibromochloromethane | U | 400 |
| diethyl ether | U | 2000 | 1,2-dibromoethane | U | 400 |
| acetone | U | 2000 | chlorobenzene | U | 400 |
| 1,1-dichloroethene | 310 | 200 | 1,1,1,2-tetrachloroethane | U | 400 |
| methylene chloride | U | 1000 | ethylbenzene | 1200 | 400 |
| carbon disulfide | U | 400 | m&p-xylenes | 4000 | 400 |
| methyl t-butyl ether (MTBE) | U | 400 | o-xylene | 1200 | 400 |
| trans-1,2-dichloroethene | U | 400 | styrene | U | 400 |
| isopropyl ether (DIPE) | U | 400 | bromoform | U | 400 |
| ethyl t-butyl ether (ETBE) | U | 400 | isopropylbenzene | U | 400 |
| 1,1-dichloroethane | 1600 | 400 | 1,1,1,2-tetrachloroethane | U | 400 |
| t-butanol (TBA) | U | 8000 | 1,2,3-trichloropropane | U | 400 |
| 2-butanone (MEK) | U | 2000 | n-propylbenzene | U | 400 |
| 2,2-dichloropropane | U | 400 | bromobenzene | U | 400 |
| cis-1,2-dichloroethene | 63000 | 400 | 1,3,5-trimethylbenzene | U | 400 |
| chloroform | U | 400 | 2-chlorotoluene | U | 400 |
| bromochloromethane | U | 400 | 4-chlorotoluene | U | 400 |
| tetrahydrofuran (THF) | U | 2000 | tert-butylbenzene | U | 400 |
| 1,1,1-trichloroethane | 25000 | 400 | 1,2,4-trimethylbenzene | U | 400 |
| 1,1-dichloropropene | U | 400 | sec-butylbenzene | U | 400 |
| t-amyl-methyl ether (TAME) | U | 400 | 1,3-dichlorobenzene | U | 400 |
| carbon tetrachloride | U | 400 | 4-isopropyltoluene | U | 400 |
| 1,2-dichloroethane | U | 400 | 1,4-dichlorobenzene | U | 400 |
| benzene | U | 400 | 1,2-dichlorobenzene | 850 | 400 |
| trichloroethene | 9400 | 400 | n-butylbenzene | U | 400 |
| 1,2-dichloropropane | U | 400 | 1,2-dibromo-3-chloropropane | U | 400 |
| bromodichloromethane | U | 400 | 1,2,4-trichlorobenzene | U | 400 |
| dibromomethane | U | 400 | hexachlorobutadiene | U | 400 |
| 4-methyl-2-pentanone (MIBK) | U | 2000 | naphthalene | U | 1000 |
| cis-1,3-dichloropropene | U | 400 | 1,2,3-trichlorobenzene | U | 400 |
| toluene | 36000 | 400 | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 93 | 78-114 | | | |
| toluene-D8 | 102 | 88-110 | | | |
| 4-bromofluorobenzene | 101 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-13
Sample Designation: ESM-16
Date Sampled: 4/7/06
Date Analyzed: 4/13/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | 2 | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | 41 | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | 100 | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | 16 | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | 3 | 2 |
| 1,1-dichloroethane | 14 | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | 2 | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | 3 | 2 | 1,3,5-trimethylbenzene | 7 | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | 3 | 2 | 1,2,4-trimethylbenzene | 15 | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | 110 | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | 3 | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | 110 | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 94 | 78-114 | | | |
| toluene-D8 | 106 | 88-110 | | | |
| 4-bromofluorobenzene | 109 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-14
Sample Designation: ESM-03B-D
Date Sampled: 4/7/06
Date Analyzed: 4/14/06
Matrix: Water
Instrument Dilution Factor: 200
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 400 | trans-1,3-dichloropropene | U | 400 |
| chloromethane | U | 400 | 2-hexanone | U | 2000 |
| vinyl chloride | U | 400 | 1,1,2-trichloroethane | U | 400 |
| bromomethane | U | 400 | 1,3-dichloropropane | U | 400 |
| chloroethane | U | 400 | tetrachloroethene | 14000 | 400 |
| trichlorofluoromethane | U | 400 | dibromochloromethane | U | 400 |
| diethyl ether | U | 2000 | 1,2-dibromoethane | U | 400 |
| acetone | U | 2000 | chlorobenzene | U | 400 |
| 1,1-dichloroethene | 3000 | 200 | 1,1,1,2-tetrachloroethane | U | 400 |
| methylene chloride | 12000 | 1000 | ethylbenzene | U | 400 |
| carbon disulfide | U | 400 | m&p-xylenes | U | 400 |
| methyl t-butyl ether (MTBE) | U | 400 | o-xylene | U | 400 |
| trans-1,2-dichloroethene | U | 400 | styrene | U | 400 |
| isopropyl ether (DIPE) | U | 400 | bromoforn | U | 400 |
| ethyl t-butyl ether (ETBE) | U | 400 | isopropylbenzene | U | 400 |
| 1,1-dichloroethane | 430 | 400 | 1,1,2,2-tetrachloroethane | U | 400 |
| t-butanol (TBA) | U | 8000 | 1,2,3-trichloropropane | U | 400 |
| 2-butanone (MEK) | U | 2000 | n-propylbenzene | U | 400 |
| 2,2-dichloropropane | U | 400 | bromobenzene | U | 400 |
| cis-1,2-dichloroethene | U | 400 | 1,3,5-trimethylbenzene | U | 400 |
| chloroform | U | 400 | 2-chlorotoluene | U | 400 |
| bromochloromethane | U | 400 | 4-chlorotoluene | U | 400 |
| tetrahydrofuran (THF) | U | 2000 | tert-butylbenzene | U | 400 |
| 1,1,1-trichloroethane | 23000 | 400 | 1,2,4-trimethylbenzene | U | 400 |
| 1,1-dichloropropene | U | 400 | sec-butylbenzene | U | 400 |
| t-amyl-methyl ether (TAME) | U | 400 | 1,3-dichlorobenzene | U | 400 |
| carbon tetrachloride | U | 400 | 4-isopropyltoluene | U | 400 |
| 1,2-dichloroethane | 1600 | 400 | 1,4-dichlorobenzene | U | 400 |
| benzene | U | 400 | 1,2-dichlorobenzene | U | 400 |
| trichloroethene | 57000 | 400 | n-butylbenzene | U | 400 |
| 1,2-dichloropropane | U | 400 | 1,2-dibromo-3-chloropropane | U | 400 |
| bromodichloromethane | U | 400 | 1,2,4-trichlorobenzene | U | 400 |
| dibromomethane | U | 400 | hexachlorobutadiene | U | 400 |
| 4-methyl-2-pentanone (MIBK) | U | 2000 | naphthalene | U | 1000 |
| cis-1,3-dichloropropene | U | 400 | 1,2,3-trichlorobenzene | U | 400 |
| toluene | 2500 | 400 | | | |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 95 | 78-114 | | | |
| toluene-D8 | 108 | 88-110 | | | |
| 4-bromofluorobenzene | 97 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-15
Sample Designation: ESM-03B-S
Date Sampled: 4/7/06
Date Analyzed: 4/13/06
Matrix: Water
Instrument Dilution Factor: 20
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 40 | trans-1,3-dichloropropene | U | 40 |
| chloromethane | U | 40 | 2-hexanone | U | 200 |
| vinyl chloride | U | 40 | 1,1,2-trichloroethane | U | 40 |
| bromomethane | U | 40 | 1,3-dichloropropane | U | 40 |
| chloroethane | U | 40 | tetrachloroethene | 6900 | 40 |
| trichlorofluoromethane | U | 40 | dibromochloromethane | U | 40 |
| diethyl ether | U | 200 | 1,2-dibromoethane | U | 40 |
| acetone | U | 200 | chlorobenzene | U | 40 |
| 1,1-dichloroethene | 1200 | 20 | 1,1,1,2-tetrachloroethane | U | 40 |
| methylene chloride | 2900 | 100 | ethylbenzene | 55 | 40 |
| carbon disulfide | U | 40 | m&p-xylenes | 61 | 40 |
| methyl t-butyl ether (MTBE) | U | 40 | o-xylene | U | 40 |
| trans-1,2-dichloroethene | U | 40 | styrene | U | 40 |
| isopropyl ether (DIPE) | U | 40 | bromoform | U | 40 |
| ethyl t-butyl ether (ETBE) | U | 40 | isopropylbenzene | U | 40 |
| 1,1-dichloroethane | 230 | 40 | 1,1,2,2-tetrachloroethane | U | 40 |
| t-butanol (TBA) | U | 800 | 1,2,3-trichloropropane | U | 40 |
| 2-butanone (MEK) | U | 200 | n-propylbenzene | U | 40 |
| 2,2-dichloropropane | U | 40 | bromobenzene | U | 40 |
| cis-1,2-dichloroethene | 92 | 40 | 1,3,5-trimethylbenzene | U | 40 |
| chloroform | U | 40 | 2-chlorotoluene | U | 40 |
| bromochloromethane | U | 40 | 4-chlorotoluene | U | 40 |
| tetrahydrofuran (THF) | U | 200 | tert-butylbenzene | U | 40 |
| 1,1,1-trichloroethane | 7800 | 40 | 1,2,4-trimethylbenzene | U | 40 |
| 1,1-dichloropropene | U | 40 | sec-butylbenzene | U | 40 |
| t-amyl-methyl ether (TAME) | U | 40 | 1,3-dichlorobenzene | U | 40 |
| carbon tetrachloride | U | 40 | 4-isopropyltoluene | U | 40 |
| 1,2-dichloroethane | 470 | 40 | 1,4-dichlorobenzene | U | 40 |
| benzene | U | 40 | 1,2-dichlorobenzene | 140 | 40 |
| trichloroethene | 19000 D | 100 | n-butylbenzene | U | 40 |
| 1,2-dichloropropane | U | 40 | 1,2-dibromo-3-chloropropane | U | 40 |
| bromodichloromethane | U | 40 | 1,2,4-trichlorobenzene | U | 40 |
| dibromomethane | U | 40 | hexachlorobutadiene | U | 40 |
| 4-methyl-2-pentanone (MIBK) | U | 200 | naphthalene | U | 100 |
| cis-1,3-dichloropropene | U | 40 | 1,2,3-trichlorobenzene | U | 40 |
| toluene | 420 | 40 | | | |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| dibromofluoromethane | 93 | 78-114 |
| toluene-D8 | 104 | 88-110 |
| 4-bromofluorobenzene | 102 | 86-115 |

D = Result obtained by a re-analysis at a dilution.
U = Below quantitation limit

Lab Number: 10154-16
Sample Designation: ESM-01
Date Sampled: 4/7/06
Date Analyzed: 4/13/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | 22 | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | 54 | 2 | tetrachloroethene | 21 | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | 9 | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | 3 | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | 4 | 2 |
| trans-1,2-dichloroethene | 3 | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | 110 | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | 13 | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | 18 | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | 11 | 2 |
| 1,2-dichloroethane | 3 | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | 83 | 2 |
| trichloroethene | 16 | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | 8 | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 92 | 78-114 | | | |
| toluene-D8 | 102 | 88-110 | | | |
| 4-bromofluorobenzene | 111 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-05
 Sample Designation: ESM14
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Extracted: 4/11/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Dilution Factor: 1
 Analyst: AJD

EPH ANALYTICAL RESULTS

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | Concentration ug/L | Quantitation Limit ug/L | Date of Analysis |
|--|-----------------------|----------------------------|------------------|
| Diesel PAH Analytes: | | | |
| naphthalene | U | 0.5 | 4/20/06 |
| 2-methylnaphthalene | U | 0.5 | 4/20/06 |
| phenanthrene | U | 0.5 | 4/20/06 |
| acenaphthene | U | 0.5 | 4/20/06 |
| Other PAH Analytes: | | | |
| acenaphthylene | U | 0.5 | 4/20/06 |
| fluorene | U | 0.5 | 4/20/06 |
| anthracene | U | 0.5 | 4/20/06 |
| fluoranthene | 1.2 | 0.5 | 4/20/06 |
| pyrene | 1.1 | 0.5 | 4/20/06 |
| benzo(a)anthracene | 0.7 | 0.5 | 4/20/06 |
| chrysene | 0.7 | 0.5 | 4/20/06 |
| benzo(b)fluoranthene | 1.2 | 0.5 | 4/20/06 |
| benzo(k)fluoranthene | 0.6 | 0.5 | 4/20/06 |
| benzo(a)pyrene | 0.9 | 0.2 | 4/20/06 |
| indeno(1,2,3-cd)pyrene | 0.7 | 0.5 | 4/20/06 |
| dibenzo(a,h)anthracene | U | 0.5 | 4/20/06 |
| benzo(g,h,i)perylene | 0.8 | 0.5 | 4/20/06 |
| Ranges: | | | |
| Unadjusted C11-C22 Aromatics (1) | U | 200 | 4/20/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | U | 200 | 4/20/06 |
| Extraction Surrogate Recoveries: | | | |
| 1-chloro-octadecane | 10% # | | |
| o-terphenyl | 23% # | | |
| Fractionation Surrogate Recoveries: | | | |
| 2-fluorobiphenyl | 101% | | |
| 2-bromonaphthalene | 99% | | |
| Surrogate Acceptance Range | 40-140% | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

= Surrogate below acceptance criteria. Matrix interference suspected.

RL Resource Laboratories, LLC

Lab Number: 10154-07
 Sample Designation: ESM13
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Extracted: 4/11/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Dilution Factor: 1
 Analyst: AJD

EPH ANALYTICAL RESULTS

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | Concentration ug/L | Quantitation Limit ug/L | Date of Analysis |
|-------------------------------------|-----------------------|----------------------------|------------------|
| Diesel PAH Analytes: | | | |
| naphthalene | 3.3 | 0.6 | 4/20/06 |
| 2-methylnaphthalene | U | 0.6 | 4/20/06 |
| phenanthrene | U | 0.6 | 4/20/06 |
| acenaphthene | U | 0.6 | 4/20/06 |
| Other PAH Analytes: | | | |
| acenaphthylene | U | 0.6 | 4/20/06 |
| fluorene | U | 0.6 | 4/20/06 |
| anthracene | U | 0.6 | 4/20/06 |
| fluoranthene | U | 0.6 | 4/20/06 |
| pyrene | U | 0.6 | 4/20/06 |
| benzo(a)anthracene | U | 0.6 | 4/20/06 |
| chrysene | U | 0.6 | 4/20/06 |
| benzo(b)fluoranthene | U | 0.6 | 4/20/06 |
| benzo(k)fluoranthene | U | 0.6 | 4/20/06 |
| benzo(a)pyrene | U | 0.2 | 4/20/06 |
| Indeno(1,2,3-cd)pyrene | U | 0.6 | 4/20/06 |
| dibenzo(a,h)anthracene | U | 0.6 | 4/20/06 |
| benzo(g,h,i)perylene | U | 0.6 | 4/20/06 |
| Ranges: | | | |
| Unadjusted C11-C22 Aromatics (1) | U | 200 | 4/20/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | U | 200 | 4/20/06 |
| Extraction Surrogate Recoveries: | | | |
| 1-chloro-octadecane | 9% # | | |
| o-terphenyl | 24% # | | |
| Fractionation Surrogate Recoveries: | | | |
| 2-fluorobiphenyl | 106% | | |
| 2-bromonaphthalene | 102% | | |
| Surrogate Acceptance Range | 40-140% | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or Internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

= Surrogate below acceptance criteria. Matrix interference suspected.

RL Resource Laboratories, LLC

V

Lab Number: 10154-08
 Sample Designation: ESM-8B
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Extracted: 4/11/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Dilution Factor: 1
 Analyst: AJD

EPH ANALYTICAL RESULTS

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | Concentration ug/L | Quantitation Limit ug/L | Date of Analysis |
|--|-----------------------|----------------------------|------------------|
| Diesel PAH Analytes: | | | |
| naphthalene | U | 0.5 | 4/20/06 |
| 2-methylnaphthalene | U | 0.5 | 4/20/06 |
| phenanthrene | U | 0.5 | 4/20/06 |
| acenaphthene | U | 0.5 | 4/20/06 |
| Other PAH Analytes: | | | |
| acenaphthylene | U | 0.5 | 4/20/06 |
| fluorene | U | 0.5 | 4/20/06 |
| anthracene | U | 0.5 | 4/20/06 |
| fluoranthene | U | 0.5 | 4/20/06 |
| pyrene | U | 0.5 | 4/20/06 |
| benzo(a)anthracene | U | 0.5 | 4/20/06 |
| chrysene | U | 0.5 | 4/20/06 |
| benzo(b)fluoranthene | U | 0.5 | 4/20/06 |
| benzo(k)fluoranthene | U | 0.5 | 4/20/06 |
| benzo(a)pyrene | U | 0.2 | 4/20/06 |
| indeno(1,2,3-cd)pyrene | U | 0.5 | 4/20/06 |
| dibenzo(a,h)anthracene | U | 0.5 | 4/20/06 |
| benzo(g,h,i)perylene | U | 0.5 | 4/20/06 |
| Ranges: | | | |
| Unadjusted C11-C22 Aromatics (1) | U | 200 | 4/20/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | U | 200 | 4/20/06 |
| Extraction Surrogate Recoveries: | | | |
| 1-chloro-octadecane | 31% # | | |
| o-terphenyl | 54% | | |
| Fractionation Surrogate Recoveries: | | | |
| 2-fluorobiphenyl | 102% | | |
| 2-bromonaphthalene | 102% | | |
| Surrogate Acceptance Range | 40-140% | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

= Surrogate below acceptance criteria. Matrix interference suspected.

RL Resource Laboratories, LLC

Lab Number: 10154-09
 Sample Designation: ESM-8
 Date Sampled: 4/6/06
 Date Received: 4/7/06
 Date Extracted: 4/11/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Dilution Factor: 1
 Analyst: AJD

EPH ANALYTICAL RESULTS

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | Concentration ug/L | Quantitation Limit ug/L | Date of Analysis |
|--|-----------------------|----------------------------|------------------|
| Diesel PAH Analytes: | | | |
| naphthalene | U | 0.8 | 4/20/06 |
| 2-methylnaphthalene | U | 0.8 | 4/20/06 |
| phenanthrene | U | 0.8 | 4/20/06 |
| acenaphthene | U | 0.8 | 4/20/06 |
| Other PAH Analytes: | | | |
| acenaphthylene | U | 0.8 | 4/20/06 |
| fluorene | U | 0.8 | 4/20/06 |
| anthracene | U | 0.8 | 4/20/06 |
| fluoranthene | 1.0 | 0.8 | 4/20/06 |
| pyrene | 1.1 | 0.8 | 4/20/06 |
| benzo(a)anthracene | U | 0.8 | 4/20/06 |
| chrysene | U | 0.8 | 4/20/06 |
| benzo(b)fluoranthene | 1.0 | 0.8 | 4/20/06 |
| benzo(k)fluoranthene | U | 0.8 | 4/20/06 |
| benzo(a)pyrene | 0.5 | 0.3 | 4/20/06 |
| indeno(1,2,3-cd)pyrene | U | 0.8 | 4/20/06 |
| dibenzo(a,h)anthracene | U | 0.8 | 4/20/06 |
| benzo(g,h,i)perylene | U | 0.8 | 4/20/06 |
| Ranges: | | | |
| Unadjusted C11-C22 Aromatics (1) | U | 300 | 4/20/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | U | 300 | 4/20/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | U | 300 | 4/20/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | U | 300 | 4/20/06 |
| Extraction Surrogate Recoveries: | | | |
| 1-chloro-octadecane | 22% # | | |
| o-terphenyl | 43% | | |
| Fractionation Surrogate Recoveries: | | | |
| 2-fluorobiphenyl | 113% | | |
| 2-bromonaphthalene | 109% | | |
| Surrogate Acceptance Range | 40-140% | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

= Surrogate below acceptance criteria. Matrix interference suspected.

RL Resource Laboratories, LLC

Lab Number: 10154-01
 Sample Designation: ESM11
 Date Sampled: 4/6/06
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 65 | 30-150 |
| Decachlorobiphenyl | 87 | 30-150 |

U = Below quantitation limit

V

Lab Number: 10154-03
 Sample Designation: ESM12
 Date Sampled: 4/6/06
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 73 | 30-150 |
| Decachlorobiphenyl | 90 | 30-150 |

U = Below quantitation limit

Lab Number: 10154-04
Sample Designation: ESM02
Date Sampled: 4/6/06
Date Extracted: 4/12/06
Date Analyzed: 4/13/06
Matrix: Water
Dilution Factor: 1
Analyst: AJD

POLYCHLORINATED BIPHENYLS
SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 66 | 30-150 |
| Decachlorobiphenyl | 85 | 30-150 |

U = Below quantitation limit

✓

Lab Number: 10154-05
Sample Designation: ESM14
Date Sampled: 4/6/06
Date Extracted: 4/12/06
Date Analyzed: 4/19/06
Matrix: Water
Dilution Factor: 5
Analyst: AJD

POLYCHLORINATED BIPHENYLS
SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 2 |
| PCB-1242 | 8 | 2 |
| PCB-1221 | U | 2 |
| PCB-1232 | U | 2 |
| PCB-1248 | U | 2 |
| PCB-1254 | U | 2 |
| PCB-1260 | U | 2 |
| PCB-1262 | U | 2 |
| PCB-1268 | U | 2 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 46 | 30-150 |
| Decachlorobiphenyl | 106 | 30-150 |

U = Below quantitation limit

Lab Number: 10154-07
Sample Designation: ESM13
Date Sampled: 4/6/06
Date Extracted: 4/12/06
Date Analyzed: 4/13/06
Matrix: Water
Dilution Factor: 1
Analyst: AJD

POLYCHLORINATED BIPHENYLS
SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 76 | 30-150 |
| Decachlorobiphenyl | 90 | 30-150 |

Note: A non-aroclor like compound was observed in this sample.
U = Below quantitation limit

Lab Number: 10154-08
 Sample Designation: ESM-8B
 Date Sampled: 4/6/06
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 61 | 30-150 |
| Decachlorobiphenyl | 84 | 30-150 |

U = Below quantitation limit

Lab Number: 10154-09
Sample Designation: ESM-8
Date Sampled: 4/6/06
Date Extracted: 4/12/06
Date Analyzed: 4/19/06
Matrix: Water
Dilution Factor: 1
Analyst: AJD

POLYCHLORINATED BIPHENYLS
SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.5 |
| PCB-1242 | 0.8 | 0.5 |
| PCB-1221 | U | 0.5 |
| PCB-1232 | U | 0.5 |
| PCB-1248 | U | 0.5 |
| PCB-1254 | U | 0.5 |
| PCB-1260 | U | 0.5 |
| PCB-1262 | U | 0.5 |
| PCB-1268 | U | 0.5 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 80 | 30-150 |
| Decachlorobiphenyl | 91 | 30-150 |

Note: A non-aroclor like compound was observed in this sample.
U = Below quantitation limit

Lab Number: 10154-10
 Sample Designation: ESM-3
 Date Sampled: 4/7/06
 Date Extracted: 4/12/06
 Date Analyzed: 4/14/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 78 | 30-150 |
| Decachlorobiphenyl | 98 | 30-150 |

U = Below quantitation limit

Lab Number: 10154-13
Sample Designation: ESM-18
Date Sampled: 4/7/06
Date Extracted: 4/12/06
Date Analyzed: 4/14/06
Matrix: Water
Dilution Factor: 1
Analyst: AJD

POLYCHLORINATED BIPHENYLS
SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 71 | 30-150 |
| Decachlorobiphenyl | 96 | 30-150 |

Note: A non-aroclor like compound was observed in this sample.
U = Below quantitation limit

Lab Number: 10154-14
 Sample Designation: ESM-03B-D
 Date Sampled: 4/7/06
 Date Extracted: 4/12/06
 Date Analyzed: 4/14/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS

SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 77 | 30-150 |
| Decachlorobiphenyl | 98 | 30-150 |

U = Below quantitation limit

Lab Number: 10154-15
 Sample Designation: ESM-03B-S
 Date Sampled: 4/7/06
 Date Extracted: 4/12/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 83 | 30-150 |
| Decachlorobiphenyl | 103 | 30-150 |

Note: Non-aroclor like peaks were observed in this sample.
 U = Below quantitation limit

Lab Number: 10154-16
 Sample Designation: ESM-01
 Date Sampled: 4/7/06
 Date Extracted: 4/13/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 71 | 30-150 |
| Decachlorobiphenyl | 87 | 30-150 |

Note: Non-aroclor like peaks were observed in this sample.
 U = Below quantitation limit

Project ID: DND Lewis 2006-056

Lab ID: 10154

Lab Number: 10154-002

Sample ID: ESM11

Matrix: Water

Sampled: 4/7/06 10:30

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|----------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Arsenic | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:31 | SW3005A6010B |
| Barium | 0.20 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:31 | SW3005A6010B |
| Cadmium | < 0.005 | 0.005 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:31 | SW3005A6010B |
| Chromium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:31 | SW3005A6010B |
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:31 | SW3005A6010B |
| Mercury | < 0.0009 | 0.0009 | mg/L | 1 | BJS | 4/12/06 | 4/12/06 | N/A | SW7470A |
| Selenium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:31 | SW3005A6010B |
| Silver | < 0.007 | 0.007 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:31 | SW3005A6010B |

Lab Number: 10154-010

Sample ID: ESM-3

Matrix: Water

Sampled: 4/7/06 9:45

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|----------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Arsenic | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:39 | SW3005A6010B |
| Barium | 0.07 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:39 | SW3005A6010B |
| Cadmium | < 0.005 | 0.005 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:39 | SW3005A6010B |
| Chromium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:39 | SW3005A6010B |
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:39 | SW3005A6010B |
| Mercury | < 0.0009 | 0.0009 | mg/L | 1 | BJS | 4/12/06 | 4/12/06 | N/A | SW7470A |
| Selenium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:39 | SW3005A6010B |
| Silver | < 0.007 | 0.007 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:39 | SW3005A6010B |

Lab Number: 10154-012

Sample ID: ESM-14

Matrix: Water

Sampled: 4/7/06 8:15

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|----------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Arsenic | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:34 | SW3005A6010B |
| Barium | 0.06 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:34 | SW3005A6010B |
| Cadmium | < 0.005 | 0.005 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:34 | SW3005A6010B |
| Chromium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:34 | SW3005A6010B |
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:34 | SW3005A6010B |
| Mercury | < 0.0009 | 0.0009 | mg/L | 1 | BJS | 4/12/06 | 4/12/06 | N/A | SW7470A |
| Selenium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:34 | SW3005A6010B |
| Silver | < 0.007 | 0.007 | mg/L | 1 | BJS | N/A | 4/12/06 | 16:34 | SW3005A6010B |

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative
Lab # 10154

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 5 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines. The following changes were made to the analysis request per the customer: 10154-05 did not request metals analysis, VOC 8260 analysis was added to sample 10154-10.

Method Blank

No exceptions noted.

Surrogate Recoveries

EPH: The following samples had surrogates that did not meet the acceptance criteria: 10154-05, 07, 08, and 09. All other batch QC had acceptable recoveries. It is suspected that this may have been caused by matrix interference. The samples could not be re-extracted due to insufficient sample volume.

Laboratory Control Sample Results

VOC 8260: LCS 10154-53 did not meet acceptance criteria for dichlorodifluoromethane.

Samples associated with this LCS: 10154-07

VOC 8260: LCS 10154-51 did not meet the acceptance limits for 2,2 dichloropropane, and dichlorodifluoromethane. The RPD for styrene did not meet the acceptance criteria, however both recoveries were acceptable. Samples associated with this LCS: 10154-01, 03, 04, 05, 08, 09, 13, 15, and 16.

VOC 8260: LCS 10154-57 did not meet the acceptance limits for 2,2 dichloropropane, and dichlorodifluoromethane. The RPD for styrene did not meet the acceptance criteria due to a low recovery in the LCSD.

These compounds noted with failures are known to be problematic in the method.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Metals: Sample 10154-10 was analyzed as a matrix spike. The results were acceptable and are attached.

Other

EPH: The fractionation check sample (LCS) for the batch of silica gel in use for these samples met the method acceptance criteria.

VPH Target compounds and ranges were determined by GC/MS. Ranges were determined in a similar manner as described in the MADEP APH method of 2/2000.

PCB: The sample peaks most closely resembled Aroclor 1242, however there was also a similarity to Aroclor 1248. Possibly due to weathering, the Aroclor identification is

difficult and not absolute. Chromatograms are provided. Samples that contained non-aroclor peaks have been noted on the report pages. Quantification is quadratic.

VOC: The following compounds were quantified using quadratic fit:
Dichlorodifluoromethane, chloromethane, vinyl chloride, bromomethane, acetone, dibromochloromethane, isopropylbenzene, tert-butyl benzene, and hexachlorobutadiene.

Metals: RCRA metals only were requested by the customer.

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10154

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|-----------|----------|----------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 (x) | Cyanide () | Other () |
| | 8270 () | VPH (x) | 7470/7471 (x) | Other () | Other () |
| | 8082 (x) | EPH (x) | Other () | Other () | Other () |

| | | |
|---|--|-----------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No (x) |
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes () No (x) |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes () No (x) |

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Position: Lab DirectorPrinted Name: Susan C. SylvesterDate: 4-27-06

Lab Number: 10154-54
 Sample Designation: Method Blank
 Date Sampled: N/A
 Date Received: N/A
 Date Analyzed: 4/17/06
 Matrix: Water
 Instrument Dilution Factor: 1
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | U | 100 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | U | 100 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 2 |
| benzene | C5-C8 | U | 1 |
| toluene | C5-C8 | U | 2 |
| ethylbenzene | C9-C12 | U | 2 |
| m&p-xylenes | C9-C12 | U | 2 |
| o-xylene | C9-C12 | U | 2 |
| naphthalene | N/A | U | 5 |
| C5-C8 Aliphatics (1,2) | N/A | U | 100 |
| C9-C12 Aliphatics (1,3) | N/A | U | 100 |
| C9-C10 Aromatics (1) | N/A | U | 100 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 102% | |
| 2,5-dibromotoluene as aliphatic | | 104% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10154-55
 Sample Designation: LCS/LCSD
 Date Sampled: N/A
 Date Received: N/A
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: LMM

VOLATILE ORGANICS MS/MSD LCS/LCSD Report
 MADEP VPH 2004-1.1

| Compound | Spike Added (ug/L) | Sample Concentration (ug/L) | MS/LCS Concentration (ug/L) | % Recovery | QC Lower Limit | QC Upper Limit |
|---------------------------------|--------------------|-----------------------------|-----------------------------|------------|----------------|----------------|
| methyl-t-butyl ether (MTBE) | 150 | | 141 | 94% | 70% | 130% |
| benzene | 50 | | 50 | 99% | 70% | 130% |
| toluene | 150 | | 145 | 97% | 70% | 130% |
| ethylbenzene | 50 | | 48 | 96% | 70% | 130% |
| m&p-xylenes | 200 | | 200 | 100% | 70% | 130% |
| o-xylene | 100 | | 97 | 97% | 70% | 130% |
| naphthalene | 100 | | 92 | 92% | 70% | 130% |
| C5-C8 Aliphatics | 750 | | 675 | 90% | 70% | 130% |
| C9-C12 Aliphatics | 550 | | 475 | 86% | 70% | 130% |
| C9-C10 Aromatics | 100 | | 100 | 100% | 70% | 130% |
| 2,5-dibromotoluene as aromatic | | | | 102% | 70% | 130% |
| 2,5-dibromotoluene as aliphatic | | | | 104% | 70% | 130% |

| Compound | Spike Added (ug/L) | Sample Concentration (ug/L) | MSD/LCSD Concentration (ug/L) | % Recovery | QC Lower Limit | QC Upper Limit | RPD | RPD Limit |
|---------------------------------|--------------------|-----------------------------|-------------------------------|------------|----------------|----------------|-----|-----------|
| methyl-t-butyl ether (MTBE) | 150 | | 141 | 94% | 70% | 130% | 0% | 20 |
| benzene | 50 | | 49 | 98% | 70% | 130% | 1% | 20 |
| toluene | 150 | | 143 | 95% | 70% | 130% | 2% | 20 |
| ethylbenzene | 50 | | 48 | 97% | 70% | 130% | 1% | 20 |
| m&p-xylenes | 200 | | 199 | 100% | 70% | 130% | 1% | 20 |
| o-xylene | 100 | | 96 | 96% | 70% | 130% | 1% | 20 |
| naphthalene | 100 | | 96 | 96% | 70% | 130% | 4% | 20 |
| C5-C8 Aliphatics | 750 | | 664 | 89% | 70% | 130% | 2% | 20 |
| C9-C12 Aliphatics | 550 | | 472 | 86% | 70% | 130% | 1% | 20 |
| C9-C10 Aromatics | 100 | | 102 | 102% | 70% | 130% | 2% | 20 |
| 2,5-dibromotoluene as aromatic | | | | 104% | 70% | 130% | 2% | 20 |
| 2,5-dibromotoluene as aliphatic | | | | 103% | 70% | 130% | 1% | 20 |

Lab Number: 10154-50
 Sample Designation: Method Blank
 Date Sampled: N/A
 Date Analyzed: 4/12/06
 Matrix: Water
 Instrument Dilution Factor: 1
 Analyst: CWC

VOLATILE ORGANICS
 SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 95 | 78-114 | | | |
| toluene-D8 | 102 | 88-110 | | | |
| 4-bromofluorobenzene | 100 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\A\VOA03\041206W3041247.D
Date Analyzed: 4/13/05
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-------------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 13 | 67% | 13 | 67% | 1% |
| chloromethane | 15 | 75% | 15 | 73% | 3% |
| vinyl chloride | 16 | 79% | 16 | 80% | 2% |
| bromomethane | 15 | 74% | 16 | 78% | 5% |
| chloroethane | 15 | 77% | 15 | 77% | 0% |
| trichlorofluoromethane | 17 | 86% | 17 | 87% | 2% |
| diethylether | 18 | 90% | 18 | 88% | 3% |
| acetone | 15 | 76% | 15 | 77% | 1% |
| 1,1-dichloroethane | 18 | 92% | 18 | 90% | 2% |
| methylene chloride | 18 | 89% | 18 | 92% | 4% |
| carbon disulfide | 18 | 90% | 17 | 87% | 3% |
| methyl-t-butyl ether (MTBE) | 34 | 85% | 34 | 84% | 1% |
| trans-1,2-dichloroethene | 19 | 94% | 18 | 89% | 5% |
| isopropyl ether (DIPE) | 18 | 92% | 18 | 91% | 1% |
| ethyl-t-butyl ether (ETBE) | 18 | 91% | 19 | 93% | 2% |
| 1,1-dichloroethane | 17 | 85% | 17 | 83% | 2% |
| t-butanol (TBA) | 73 | 73% | 77 | 77% | 5% |
| 2-butanone (MEK) | 15 | 77% | 16 | 78% | 2% |
| 2,2-dichloropropane | 12 | 62% | 12 | 69% | 5% |
| cis-1,2-dichloroethene | 21 | 103% | 20 | 101% | 2% |
| chloroform | 19 | 95% | 18 | 92% | 4% |
| bromochloromethane | 21 | 105% | 21 | 106% | 0% |
| tetrahydrofuran (THF) | 16 | 79% | 17 | 83% | 4% |
| 1,1,1-trichloroethane | 18 | 91% | 18 | 88% | 3% |
| 1,1-dichloropropane | 20 | 99% | 18 | 91% | 9% |
| i-amyl-methyl ether (TAME) | 19 | 93% | 18 | 90% | 3% |
| carbon tetrachloride | 16 | 81% | 16 | 80% | 1% |
| 1,2-dichloroethane | 17 | 84% | 16 | 82% | 3% |
| benzene | 20 | 101% | 19 | 94% | 7% |
| trichloroethene | 21 | 106% | 21 | 103% | 2% |
| 1,2-dichloropropane | 19 | 97% | 19 | 93% | 4% |
| bromodichloromethane | 16 | 82% | 17 | 83% | 2% |
| dibromomethane | 20 | 102% | 20 | 101% | 1% |
| 4-methyl-2-pentanone (MIBK) | 18 | 90% | 19 | 94% | 4% |
| cis-1,3-dichloropropene | 18 | 90% | 18 | 89% | 1% |
| toluene | 20 | 101% | 21 | 104% | 3% |
| trans-1,3-dichloropropene | 16 | 80% | 16 | 79% | 1% |
| 2-hexanone | 16 | 78% | 16 | 80% | 2% |
| 1,1,2-trichloroethane | 21 | 103% | 21 | 103% | 0% |
| 1,3-dichloropropane | 21 | 106% | 22 | 108% | 2% |
| tetrachloroethene | 26 | 129% | 25 | 125% | 3% |
| dibromochloromethane | 20 | 98% | 21 | 106% | 7% |
| 1,2-dibromoethane (EDB) | 21 | 106% | 22 | 112% | 6% |
| chlorobenzene | 22 | 111% | 23 | 113% | 2% |
| 1,1,1,2-tetrachloroethane | 21 | 106% | 22 | 111% | 4% |
| ethylbenzene | 22 | 109% | 22 | 111% | 2% |
| m&p-xylenes | 46 | 115% | 45 | 114% | 1% |
| o-xylene | 23 | 115% | 24 | 118% | 3% |
| styrene | 22 | 112% | 15 | 75% | 39% |
| bromoform | 21 | 104% | 22 | 110% | 6% |
| isopropylbenzene | 24 | 119% | 25 | 123% | 4% |
| 1,1,2,2-tetrachloroethane | 18 | 89% | 19 | 96% | 7% |
| 1,2,3-trichloropropane | 17 | 87% | 18 | 92% | 5% |
| n-propylbenzene | 20 | 101% | 20 | 101% | 1% |
| bromobenzene | 21 | 105% | 22 | 109% | 4% |
| 1,3,5-trimethylbenzene | 19 | 96% | 19 | 95% | 1% |
| 2-chlorotoluene | 19 | 94% | 19 | 95% | 1% |
| 4-chlorotoluene | 19 | 94% | 19 | 95% | 1% |
| tert-butylbenzene | 18 | 88% | 18 | 88% | 0% |
| 1,2,4-trimethylbenzene | 19 | 97% | 20 | 100% | 3% |
| sec-butylbenzene | 20 | 101% | 20 | 99% | 2% |
| 1,3-dichlorobenzene | 20 | 102% | 21 | 105% | 3% |
| 4-isopropyltoluene | 21 | 104% | 21 | 105% | 1% |
| 1,4-dichlorobenzene | 20 | 101% | 20 | 102% | 1% |
| 1,2-dichlorobenzene | 21 | 104% | 21 | 104% | 0% |
| n-butylbenzene | 20 | 101% | 20 | 102% | 0% |
| 1,2-dibromo-3-chloropropane (| 16 | 80% | 18 | 88% | 9% |
| 1,2,4-trichlorobenzene | 19 | 95% | 19 | 97% | 2% |
| hexachlorobutadiene | 20 | 100% | 19 | 94% | 6% |
| naphthalene | 15 | 77% | 17 | 84% | 9% |
| 1,2,3-trichlorobenzene | 19 | 96% | 20 | 102% | 5% |
| 1,4-dioxane | 34 | 85% | 33 | 84% | 2% |
| SURROGATE STANDARDS | | | | | |
| SS dibromofluoromethane | | 95% | | 96% | |
| SS toluene-D8 | | 104% | | 98% | |
| SS 4-bromofluorobenzene | | 107% | | 113% | |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

+ =The RPD was outside the acceptance limits.

Lab Number: 10154-52
Sample Designation: Method Blank
Date Sampled: N/A
Date Analyzed: 4/12/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: CWC

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 94 | 78-114 | | | |
| toluene-D8 | 102 | 88-110 | | | |
| 4-bromofluorobenzene | 99 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-53
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\VOA03\041206\1V3041223.D
Date Analyzed: 4/12/06
SW 846 Method 5030B/5260B

| Compound | LCS | | LCSD | | RPD |
|-----------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 15 | 77% | 13 | 67% | 14% |
| chloromethane | 17 | 83% | 16 | 78% | 6% |
| vinyl chloride | 18 | 89% | 16 | 82% | 7% |
| bromomethane | 18 | 88% | 16 | 80% | 10% |
| chloroethane | 17 | 87% | 16 | 78% | 12% |
| trichlorofluoromethane | 20 | 98% | 17 | 84% | 15% |
| diethylether | 19 | 95% | 19 | 97% | 2% |
| acetone | 15 | 75% | 14 | 72% | 5% |
| 1,1-dichloroethane | 20 | 99% | 19 | 94% | 5% |
| methylene chloride | 20 | 102% | 19 | 98% | 7% |
| carbon disulfide | 20 | 98% | 19 | 93% | 6% |
| methyl-t-butyl ether (MTBE) | 38 | 95% | 36 | 90% | 6% |
| trans-1,2-dichloroethene | 20 | 98% | 19 | 94% | 4% |
| isopropyl ether (DIPE) | 20 | 102% | 20 | 98% | 2% |
| ethyl-t-butyl ether (ETBE) | 20 | 101% | 19 | 93% | 9% |
| 1,1-dichloroethane | 19 | 93% | 17 | 87% | 6% |
| t-butanol (TBA) | 87 | 87% | 87 | 87% | 0% |
| 2-butanone (MEK) | 18 | 91% | 16 | 82% | 10% |
| 2,2-dichloropropane | 19 | 93% | 17 | 86% | 8% |
| cis-1,2-dichloroethane | 23 | 113% | 21 | 105% | 7% |
| chloroform | 21 | 104% | 20 | 98% | 6% |
| bromochloromethane | 23 | 115% | 22 | 111% | 4% |
| tetrahydrofuran (THF) | 18 | 92% | 17 | 85% | 7% |
| 1,1,1-trichloroethane | 20 | 98% | 18 | 89% | 9% |
| 1,1-dichloropropene | 22 | 108% | 20 | 99% | 9% |
| t-amyl-methyl ether (TAME) | 21 | 103% | 19 | 95% | 8% |
| carbon tetrachloride | 16 | 80% | 16 | 81% | 1% |
| 1,2-dichloroethane | 18 | 90% | 17 | 86% | 5% |
| benzene | 22 | 108% | 21 | 103% | 5% |
| trichloroethene | 22 | 110% | 21 | 103% | 8% |
| 1,2-dichloropropane | 21 | 105% | 21 | 103% | 1% |
| bromodichloromethane | 17 | 85% | 17 | 86% | 1% |
| dibromomethane | 22 | 109% | 21 | 106% | 3% |
| 4-methyl-2-pentanone (MIBK) | 21 | 103% | 19 | 97% | 6% |
| cis-1,3-dichloropropene | 21 | 103% | 19 | 97% | 6% |
| toluene | 22 | 111% | 22 | 108% | 3% |
| trans-1,3-dichloropropene | 18 | 88% | 18 | 90% | 2% |
| 2-hexanone | 19 | 94% | 18 | 92% | 2% |
| 1,1,2-trichloroethane | 22 | 110% | 21 | 107% | 3% |
| 1,3-dichloropropane | 22 | 110% | 22 | 110% | 0% |
| tetrachloroethene | 25 | 124% | 24 | 120% | 4% |
| dibromochloromethane | 19 | 94% | 20 | 101% | 7% |
| 1,2-dibromoethane (EDB) | 22 | 111% | 23 | 115% | 3% |
| chlorobenzene | 23 | 115% | 22 | 112% | 3% |
| 1,1,1,2-tetrachloroethane | 20 | 100% | 21 | 104% | 4% |
| ethylbenzene | 22 | 112% | 22 | 111% | 1% |
| m&p-xylenes | 48 | 119% | 47 | 117% | 2% |
| o-xylene | 24 | 121% | 24 | 119% | 1% |
| styrene | 24 | 118% | 23 | 117% | 1% |
| bromoform | 19 | 96% | 20 | 100% | 4% |
| isopropylbenzene | 25 | 125% | 24 | 119% | 5% |
| 1,1,2,2-tetrachloroethane | 19 | 93% | 19 | 93% | 1% |
| 1,2,3-trichloropropane | 18 | 92% | 18 | 92% | 0% |
| n-propylbenzene | 21 | 107% | 21 | 104% | 3% |
| bromobenzene | 22 | 112% | 22 | 109% | 2% |
| 1,3,5-trimethylbenzene | 20 | 102% | 20 | 98% | 4% |
| 2-chlorotoluene | 19 | 97% | 20 | 98% | 0% |
| 4-chlorotoluene | 20 | 101% | 19 | 94% | 7% |
| tert-butylbenzene | 18 | 91% | 18 | 89% | 2% |
| 1,2,4-trimethylbenzene | 20 | 102% | 20 | 101% | 1% |
| sec-butylbenzene | 20 | 101% | 19 | 97% | 3% |
| 1,3-dichlorobenzene | 21 | 105% | 20 | 101% | 5% |
| 4-isopropyltoluene | 22 | 110% | 22 | 109% | 1% |
| 1,4-dichlorobenzene | 21 | 105% | 21 | 103% | 2% |
| 1,2-dichlorobenzene | 21 | 107% | 21 | 107% | 0% |
| n-butylbenzene | 22 | 108% | 21 | 104% | 4% |
| 1,2-dibromo-3-chloropropane | 17 | 84% | 17 | 85% | 1% |
| 1,2,4-trichlorobenzene | 20 | 100% | 20 | 98% | 2% |
| hexachlorobutadiene | 21 | 103% | 21 | 103% | 0% |
| naphthalene | 17 | 84% | 17 | 86% | 2% |
| 1,2,3-trichlorobenzene | 21 | 103% | 21 | 106% | 3% |
| 1,4-dioxane | 40 | 100% | 39 | 97% | 3% |

SURROGATE STANDARDS

| | | |
|---------------------------|------|------|
| SS dibromodifluoromethane | 98% | 96% |
| SS toluene-DB | 106% | 108% |
| SS 4-bromodifluorobenzene | 107% | 111% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

Lab Number: 10154-56
Sample Designation: Method Blank
Date Sampled: N/A
Date Analyzed: 4/14/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 97 | 78-114 | | | |
| toluene-D8 | 108 | 88-110 | | | |
| 4-bromofluorobenzene | 102 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10154-57
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: C:\DATA\WQA03\041306\1V3041347.D
Date Analyzed: 4/14/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-----------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 15 | 74% | 12 | 62% | 17% |
| chloromethane | 17 | 87% | 16 | 80% | 8% |
| vinyl chloride | 23 | 115% | 20 | 98% | 16% |
| bromomethane | 17 | 85% | 16 | 82% | 4% |
| chloroethane | 17 | 87% | 17 | 83% | 5% |
| trichlorofluoromethane | 19 | 93% | 17 | 84% | 10% |
| diethylether | 18 | 91% | 18 | 92% | 2% |
| acetone | 17 | 85% | 16 | 80% | 5% |
| 1,1-dichloroethane | 19 | 94% | 16 | 81% | 15% |
| methylene chloride | 19 | 94% | 18 | 90% | 4% |
| carbon disulfide | 18 | 90% | 16 | 80% | 12% |
| methyl-t-butyl ether (MTBE) | 36 | 88% | 34 | 86% | 4% |
| trans-1,2-dichloroethane | 19 | 93% | 17 | 86% | 8% |
| isopropyl ether (DIPE) | 19 | 95% | 18 | 89% | 6% |
| ethyl-t-butyl ether (ETBE) | 19 | 95% | 18 | 92% | 3% |
| 1,1-dichloroethane | 18 | 89% | 16 | 82% | 9% |
| t-butanol (TBA) | 84 | 84% | 80 | 80% | 5% |
| 2-butanone (MEK) | 18 | 89% | 17 | 84% | 6% |
| 2,2-dichloropropane | 11 | 57% | 11 | 53% | 8% |
| cis-1,2-dichloroethane | 21 | 105% | 19 | 94% | 10% |
| chloroform | 19 | 97% | 18 | 90% | 8% |
| bromochloromethane | 22 | 109% | 20 | 102% | 7% |
| tetrahydrofuran (THF) | 19 | 94% | 19 | 93% | 1% |
| 1,1,1-trichloroethane | 19 | 95% | 17 | 87% | 9% |
| 1,1-dichloropropene | 20 | 101% | 18 | 91% | 11% |
| 1-amy1-methyl ether (TAME) | 19 | 95% | 18 | 92% | 3% |
| carbon tetrachloride | 17 | 87% | 16 | 80% | 8% |
| 1,2-dichloroethane | 18 | 88% | 16 | 82% | 7% |
| benzene | 21 | 104% | 18 | 92% | 13% |
| trichloroethene | 21 | 106% | 19 | 93% | 13% |
| 1,2-dichloropropane | 20 | 102% | 19 | 94% | 8% |
| bromodichloromethane | 17 | 87% | 16 | 81% | 8% |
| dibromomethane | 21 | 104% | 20 | 98% | 6% |
| 4-methyl-2-pentanone (MIBK) | 20 | 100% | 20 | 99% | 0% |
| cis-1,3-dichloropropene | 19 | 93% | 17 | 83% | 12% |
| toluene | 22 | 109% | 19 | 94% | 15% |
| trans-1,3-dichloropropene | 16 | 80% | 15 | 75% | 7% |
| 2-hexanone | 17 | 87% | 15 | 74% | 16% |
| 1,1,2-trichloroethane | 21 | 104% | 20 | 98% | 6% |
| 1,3-dichloropropane | 21 | 103% | 20 | 99% | 4% |
| tetrachloroethene | 23 | 114% | 22 | 108% | 5% |
| dibromochloromethane | 19 | 97% | 19 | 96% | 1% |
| 1,2-dibromoethane (EDB) | 20 | 102% | 20 | 102% | 0% |
| chlorobenzene | 21 | 107% | 20 | 101% | 6% |
| 1,1,1,2-tetrachloroethane | 21 | 103% | 20 | 99% | 4% |
| ethylbenzene | 21 | 106% | 20 | 100% | 5% |
| m&p-xylenes | 44 | 110% | 42 | 105% | 5% |
| o-xylene | 23 | 113% | 21 | 105% | 7% |
| styrene | 21 | 106% | 13 | 64% | 49% |
| bromoform | 20 | 102% | 20 | 102% | 0% |
| isopropylbenzene | 22 | 110% | 21 | 108% | 4% |
| 1,1,2,2-tetrachloroethane | 18 | 91% | 18 | 91% | 0% |
| 1,2,3-trichloropropane | 18 | 90% | 17 | 86% | 5% |
| n-propylbenzene | 20 | 102% | 18 | 92% | 9% |
| bromobenzene | 21 | 103% | 20 | 98% | 5% |
| 1,3,5-trimethylbenzene | 19 | 97% | 18 | 90% | 8% |
| 2-chlorotoluene | 19 | 93% | 18 | 88% | 6% |
| 4-chlorotoluene | 20 | 98% | 17 | 86% | 13% |
| tert-butylbenzene | 20 | 100% | 19 | 95% | 5% |
| 1,2,4-trimethylbenzene | 19 | 98% | 18 | 92% | 5% |
| sec-butylbenzene | 19 | 97% | 18 | 91% | 6% |
| 1,3-dichlorobenzene | 20 | 100% | 19 | 94% | 7% |
| 4-isopropyltoluene | 21 | 108% | 19 | 94% | 12% |
| 1,4-dichlorobenzene | 20 | 100% | 18 | 91% | 10% |
| 1,2-dichlorobenzene | 21 | 105% | 19 | 96% | 9% |
| n-butylbenzene | 20 | 99% | 18 | 92% | 7% |
| 1,2-dibromo-3-chloropropane | 18 | 92% | 17 | 87% | 5% |
| 1,2,4-trichlorobenzene | 19 | 95% | 17 | 86% | 10% |
| hexachlorobutadiene | 19 | 98% | 18 | 92% | 5% |
| naphthalene | 16 | 81% | 15 | 75% | 7% |
| 1,2,3-trichlorobenzene | 20 | 100% | 18 | 92% | 9% |
| 1,4-dioxane | 39 | 98% | 42 | 104% | 6% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 98% | 98% |
| SS toluene-D8 | 109% | 108% |
| SS 4-bromofluorobenzene | 102% | 103% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

Lab Number: 10154-50
 Sample Designation: Preparation Blank 180 W
 Date Sampled: N/A
 Date Received: N/A
 Date Extracted: 4/11/06
 Matrix: Water
 Containers: N/A
 Sample Preservation: N/A
 Temperature: N/A
 Dilution Factor: 1
 Analyst: A.JD

EPH ANALYTICAL RESULTS

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | Concentration ug/L | Quantitation Limit ug/L | Date of Analysis |
|-------------------------------------|-----------------------|----------------------------|------------------|
| Diesel PAH Analytes: | | | |
| naphthalene | U | 0.5 | 4/19/06 |
| 2-methylnaphthalene | U | 0.5 | 4/19/06 |
| phenanthrene | U | 0.5 | 4/19/06 |
| acenaphthene | U | 0.5 | 4/19/06 |
| Other PAH Analytes: | | | |
| acenaphthylene | U | 0.5 | 4/19/06 |
| fluorene | U | 0.5 | 4/19/06 |
| anthracene | U | 0.5 | 4/19/06 |
| fluoranthene | U | 0.5 | 4/19/06 |
| pyrene | U | 0.5 | 4/19/06 |
| benzo(a)anthracene | U | 0.5 | 4/19/06 |
| chrysene | U | 0.5 | 4/19/06 |
| benzo(b)fluoranthene | U | 0.5 | 4/19/06 |
| benzo(k)fluoranthene | U | 0.5 | 4/19/06 |
| benzo(a)pyrene | U | 0.2 | 4/19/06 |
| indeno(1,2,3-cd)pyrene | U | 0.5 | 4/19/06 |
| dibenzo(a,h)anthracene | U | 0.5 | 4/19/06 |
| benzo(g,h,i)perylene | U | 0.5 | 4/19/06 |
| Ranges: | | | |
| Unadjusted C11-C22 Aromatics (1) | U | 200 | 4/19/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | U | 200 | 4/19/06 |
| Extraction Surrogate Recoveries: | | | |
| 1-chloro-octadecane | 48% | | |
| o-terphenyl | 59% | | |
| Fractionation Surrogate Recoveries: | | | |
| 2-fluorobiphenyl | 112% | | |
| 2-bromonaphthalene | 109% | | |
| Surrogate Acceptance Range | 40-140% | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

Lab Number: 10154-51
 Sample Designation: Lab Control Sample/Duplicate 180 W
 Date Sampled: N/A
 Date Received: N/A
 Date Extracted: 4/11/06
 Matrix: Water
 Containers: N/A
 Sample Preservation: N/A
 Temperature: N/A
 Dilution Factor: 1
 Analyst: AJD

EPH ANALYTICAL RESULTS

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | Amount Added ug/L | Amount Found in LCS ug/L | LCS Recovery (%) | Amount Found in LCSD ug/L | LCSD Recovery (%) | RPD (%) | Acceptance Criteria (%) | Date of Analysis |
|--|-------------------------|-----------------------------------|------------------------|------------------------------------|-------------------------|------------|-------------------------------|------------------|
| Diesel PAH Analytes: | | | | | | | | |
| naphthalene | 60 | 32.45 | 54% | 33.33 | 56% | 3% | <25% | 4/19/06 |
| 2-methylnaphthalene | 60 | 32.92 | 55% | 33.92 | 57% | 3% | <25% | 4/19/06 |
| phenanthrene | 60 | 41.76 | 70% | 39.97 | 67% | -4% | <25% | 4/19/06 |
| acenaphthene | 60 | 40.13 | 67% | 39.03 | 65% | -3% | <25% | 4/19/06 |
| Other PAH Analytes: | | | | | | | | |
| acenaphthylene | 60 | 37.78 | 63% | 36.65 | 61% | -3% | <25% | 4/19/06 |
| fluorene | 60 | 38.77 | 65% | 36.16 | 60% | -7% | <25% | 4/19/06 |
| anthracene | 60 | 44.19 | 74% | 41.52 | 69% | -6% | <25% | 4/19/06 |
| fluoranthene | 60 | 46.38 | 77% | 45.51 | 76% | -2% | <25% | 4/19/06 |
| pyrene | 60 | 42.92 | 72% | 42.60 | 71% | -1% | <25% | 4/19/06 |
| benzo(a)anthracene | 60 | 45.81 | 76% | 44.91 | 75% | -2% | <25% | 4/19/06 |
| chrysene | 60 | 47.91 | 80% | 47.69 | 79% | 0% | <25% | 4/19/06 |
| benzo(b)fluoranthene | 60 | 40.12 | 67% | 36.72 | 61% | -9% | <25% | 4/19/06 |
| benzo(k)fluoranthene | 60 | 53.94 | 90% | 58.13 | 94% | 4% | <25% | 4/19/06 |
| benzo(a)pyrene | 60 | 49.79 | 83% | 48.55 | 81% | -3% | <25% | 4/19/06 |
| indeno(1,2,3-cd)pyrene | 60 | 48.74 | 81% | 47.54 | 79% | -2% | <25% | 4/19/06 |
| dibenzo(a,h)anthracene | 60 | 48.90 | 81% | 47.41 | 79% | -3% | <25% | 4/19/06 |
| benzo(g,h,i)perylene | 60 | 47.95 | 80% | 47.50 | 79% | -1% | <25% | 4/19/06 |
| Ranges: | | | | | | | | |
| Unadjusted C11-C22 Aromatics (1) | 1020 | 808.27 | 79% | 694.32 | 68% | -15% | <25% | 4/19/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | 360 | 249.47 | 69% | 264.95 | 74% | 6% | <25% | 4/20/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | 480 | 424.48 | 88% | 407.22 | 85% | -4% | <25% | 4/20/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | | | | | | | | 4/19/06 |
| Extraction Surrogate Recoveries: | | | | | | | | |
| 1-chloro-octadecane | | 100% | | 58% | | | | |
| o-terphenyl | | 66% | | 48% | | | | |
| Fractionation Surrogate Recoveries: | | | | | | | | |
| 2-fluorobiphenyl | | 129% | | 109% | | | | |
| 2-bromonaphthalene | | 120% | | 106% | | | | |
| Acceptance Range | | 40-140% | 40-140% | 40-140% | 40-140% | | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

Lab Number: 10154-52
 Sample Designation: LCS 180 W
 Date Sampled: N/A
 Date Received: N/A
 Date Extracted: 4/11/06
 Matrix: Water
 Containers: N/A
 Sample Preservation: N/A
 Temperature: N/A
 Dilution Factor: 1
 Analyst: AJD

NAPHTHALENE BREAKTHROUGH CALCULATION

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | LCS Aliphatic Breakthrough (%) | Acceptance Criteria | Date of Analysis |
|---------------------|--------------------------------------|------------------------|------------------|
| naphthalene | 0.5% | <5.0% | 4/19/06 |
| 2-methylnaphthalene | 0.4% | <5.0% | 4/19/06 |

Lab Number: 10154-53
 Sample Designation: LCSD 180 W
 Date Sampled: N/A
 Date Received: N/A
 Date Extracted: 4/11/06
 Matrix: Water
 Containers: N/A
 Sample Preservation: N/A
 Temperature: N/A
 Dilution Factor: 1
 Analyst: AJD

NAPHTHALENE BREAKTHROUGH CALCULATION

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | LCSD | | |
|---------------------|------------------------|------------|------------------|
| | Aliphatic Breakthrough | Acceptance | Date of Analysis |
| | (%) | Criteria | |
| naphthalene | 2.4% | <5.0% | 4/19/06 |
| 2-methylnaphthalene | 0.4% | <5.0% | 4/19/06 |

Lab Number: 10154-50
 Sample Designation: Preparation Blank 209 W
 Date Sampled: N/A
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 74 | 30-150 |
| Decachlorobiphenyl | 86 | 30-150 |

U = Below quantitation limit

Lab Number: 10154-51
 Sample Designation: Laboratory Control Sample/Duplicate 209 W
 Date Sampled: N/A
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: A.J.D

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Amount Added ug/L | Amount Found In LCS ug/L | LCS Recovery (%) | Amount Found In LCSD ug/L | LCSD Recovery (%) | RPD (%) | Date of Analysis |
|----------------------|-------------------------|-----------------------------------|------------------------|------------------------------------|-------------------------|------------|------------------|
| PCB-1016 | 2 | 1.5 | 75% | 1.5 | 75% | 0.0% | 4/13/06 |
| PCB-1242 | | | | | | | 4/13/06 |
| PCB-1221 | | | | | | | 4/13/06 |
| PCB-1232 | | | | | | | 4/13/06 |
| PCB-1248 | | | | | | | 4/13/06 |
| PCB-1254 | | | | | | | 4/13/06 |
| PCB-1260 | 2 | 1.4 | 70% | 1.5 | 75% | 6.9% | 4/13/06 |
| PCB-1262 | | | | | | | 4/13/06 |
| PCB-1268 | | | | | | | 4/13/06 |
| Acceptance Criteria: | | | 40-140% | | 40-140% | 20% | |

| SURROGATE STANDARDS | Recovery (%) | Acceptance Limits (%) | Recovery (%) | Acceptance Limits (%) |
|----------------------|-----------------|--------------------------|-----------------|--------------------------|
| Tetrachloro-m-xylene | 83 | 30-150 | 78 | 30-150 |
| Decachlorobiphenyl | 88 | 30-150 | 92 | 30-150 |

U = Below quantitation limit

Lab Number 10154 METALS QC REPORT
Batch QC Results

Prep Blank

| Analyte | Sample ID # | Result (mg/L) | Reporting Limit (mg/L) |
|----------------|--------------------|--------------------------|-----------------------------------|
| Silver | ICB041206 | < 0.007 | 0.007 |
| Arsenic | ICB041206 | < 0.01 | 0.01 |
| Barium | ICB041206 | < 0.05 | 0.05 |
| Cadmium | ICB041206 | < 0.005 | 0.005 |
| Chromium | ICB041206 | < 0.05 | 0.05 |
| Lead | ICB041206 | < 0.01 | 0.01 |
| Selenium | ICB041206 | < 0.05 | 0.05 |
| Mercury | PB041206 | < 0.0009 | 0.0009 |

Laboratory Control Sample

| Analyte | Sample ID # | Result (mg/L) | True Value (mg/L) | %Recovery Control Limits (85-115%) |
|----------------|--------------------|--------------------------|------------------------------|---|
| Silver | ICV041206 | 1.00 | 1.00 | 102 |
| Arsenic | ICV041206 | 0.99 | 1.00 | 98 |
| Barium | ICV041206 | 0.95 | 1.00 | 100 |
| Cadmium | ICV041206 | 0.95 | 1.00 | 101 |
| Chromium | ICV041206 | 0.95 | 1.00 | 100 |
| Lead | ICV041206 | 0.95 | 1.00 | 97 |
| Selenium | ICV041206 | 0.95 | 1.00 | 99 |
| Mercury | LCS041206 | 0.95 | 1.00 | 92 |

Sample Spike Data (MS)

| Analyte | Sample ID # | Sample Result (mg/L) | Spike Sample Result (mg/L) | Spike Amount (mg/L) | %Recovery Control Limits (75-125%) |
|----------------|--------------------|-------------------------------------|---|------------------------------------|---|
| Silver | 10154-10MS | 0 | 0.4221 | 0.5 | 84 |
| Arsenic | 10154-10MS | 0.0008 | 1.021 | 1.0 | 102 |
| Barium | 10154-10MS | 0.0738 | 0.9737 | 1.0 | 90 |
| Cadmium | 10154-10MS | 0 | 0.9933 | 1.0 | 99 |
| Chromium | 10154-10MS | 0.0014 | 0.9804 | 1.0 | 98 |
| Lead | 10154-10MS | 0.0012 | 0.9086 | 1.0 | 91 |
| Selenium | 10154-10MS | 0.0002 | 1.022 | 1.0 | 102 |
| Mercury | 10154-10MS | 0.0002 | 0.0105 | 0.0 | 103 |

Sample Spike Data (MSD)

| Analyte | Sample ID # | Sample Result (mg/L) | Spike Sample Result (mg/L) | Spike Amount (mg/L) | %Recovery Control Limits (75-125%) |
|----------------|--------------------|-------------------------------------|---|------------------------------------|---|
| Silver | 10154-10MSD | 0 | 0.4238 | 0.5 | 85 |
| Arsenic | 10154-10MSD | 0.0008 | 1.034 | 1.0 | 103 |
| Barium | 10154-10MSD | 0.0738 | 0.9916 | 1.0 | 92 |
| Cadmium | 10154-10MSD | 0 | 1.003 | 1.0 | 100 |
| Chromium | 10154-10MSD | 0.0014 | 0.9915 | 1.0 | 99 |
| Lead | 10154-10MSD | 0.0012 | 0.9131 | 1.0 | 91 |
| Selenium | 10154-10MSD | 0.0002 | 1.044 | 1.0 | 104 |

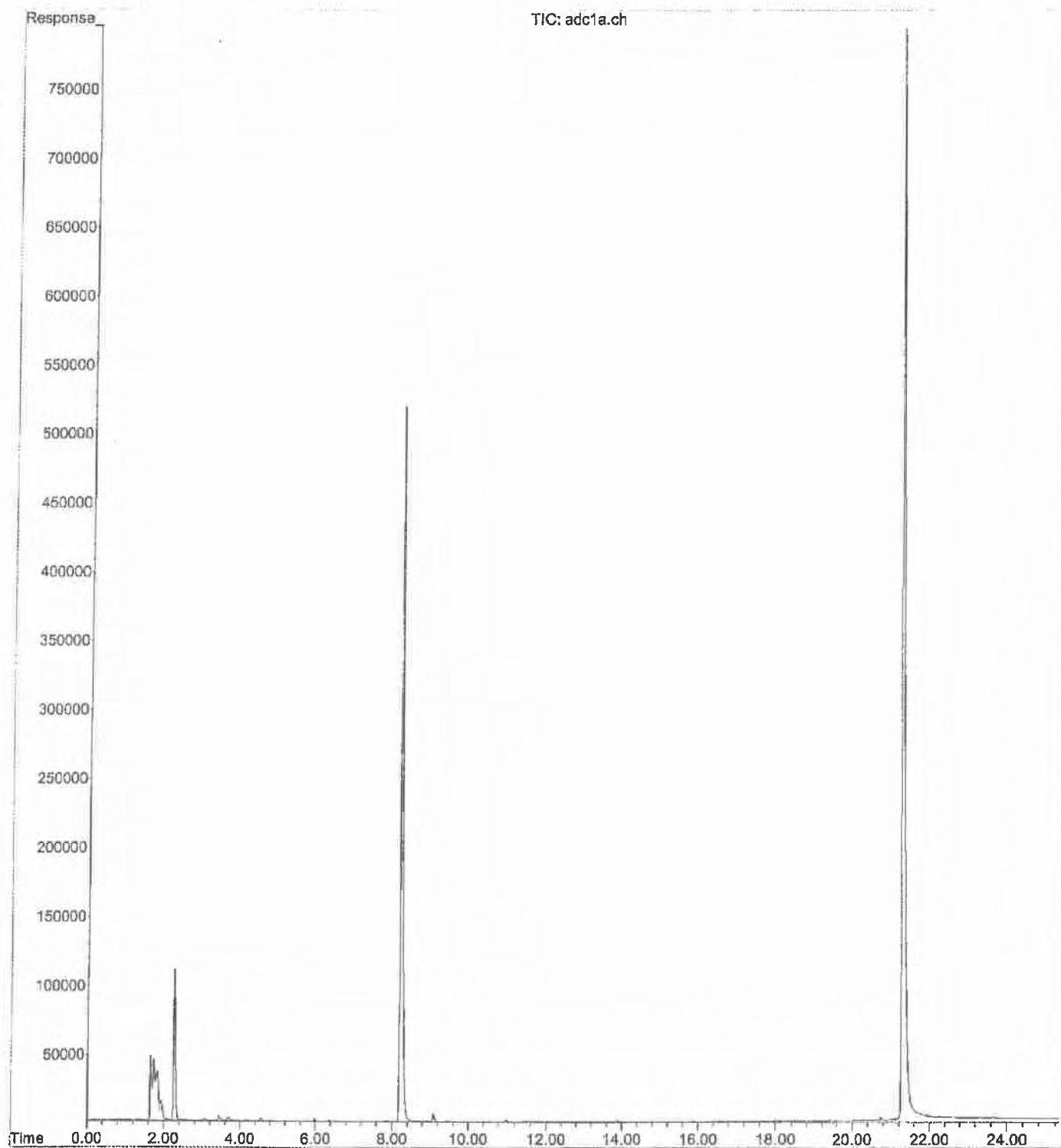
RL Resource Laboratories, LLC

| | | | | | |
|---------|-------------|--------|------|-----|----|
| Mercury | 10154-10MSD | 0.0002 | 0.01 | 0.0 | 98 |
|---------|-------------|--------|------|-----|----|

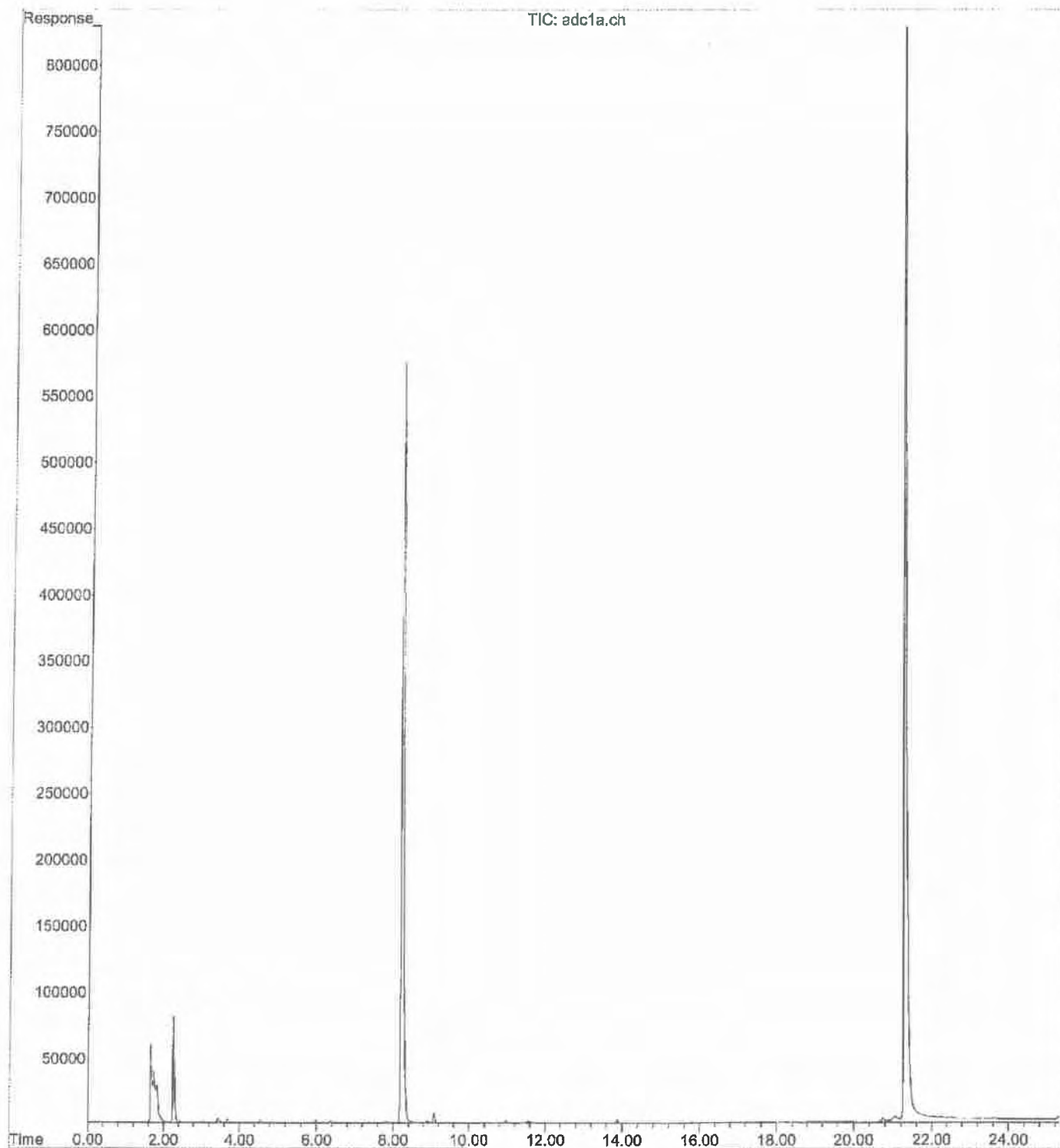
Relative Percent Difference

| Analyte | Sample ID # | MS Recovery % | MSD Recovery % | RPD (+/- 20%) |
|----------------|--------------------|--------------------------|---------------------------|--------------------------|
| Silver | 10154-10 | 84 | 85 | 0 |
| Arsenic | 10154-10 | 102 | 103 | 1 |
| Barium | 10154-10 | 90 | 92 | 2 |
| Cadmium | 10154-10 | 99 | 100 | 1 |
| Chromium | 10154-10 | 98 | 99 | 1 |
| Lead | 10154-10 | 91 | 91 | 0 |
| Selenium | 10154-10 | 102 | 104 | 2 |
| Mercury | 10154-10 | 103 | 98 | 5 |

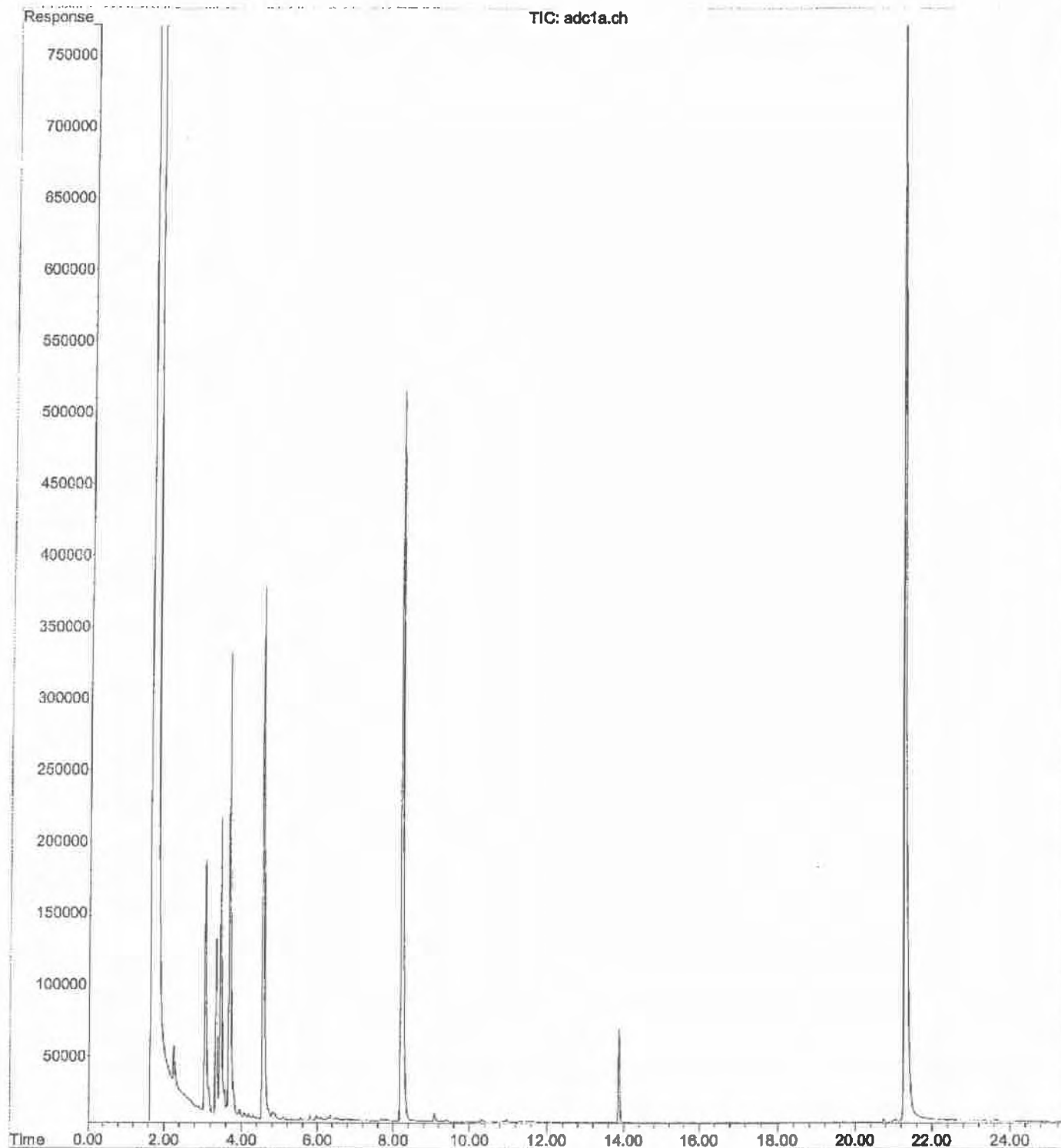
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Instrument : SVOA03
Sample Name: 10154-01 pcb x1
Misc Info :
Vial Number: 13



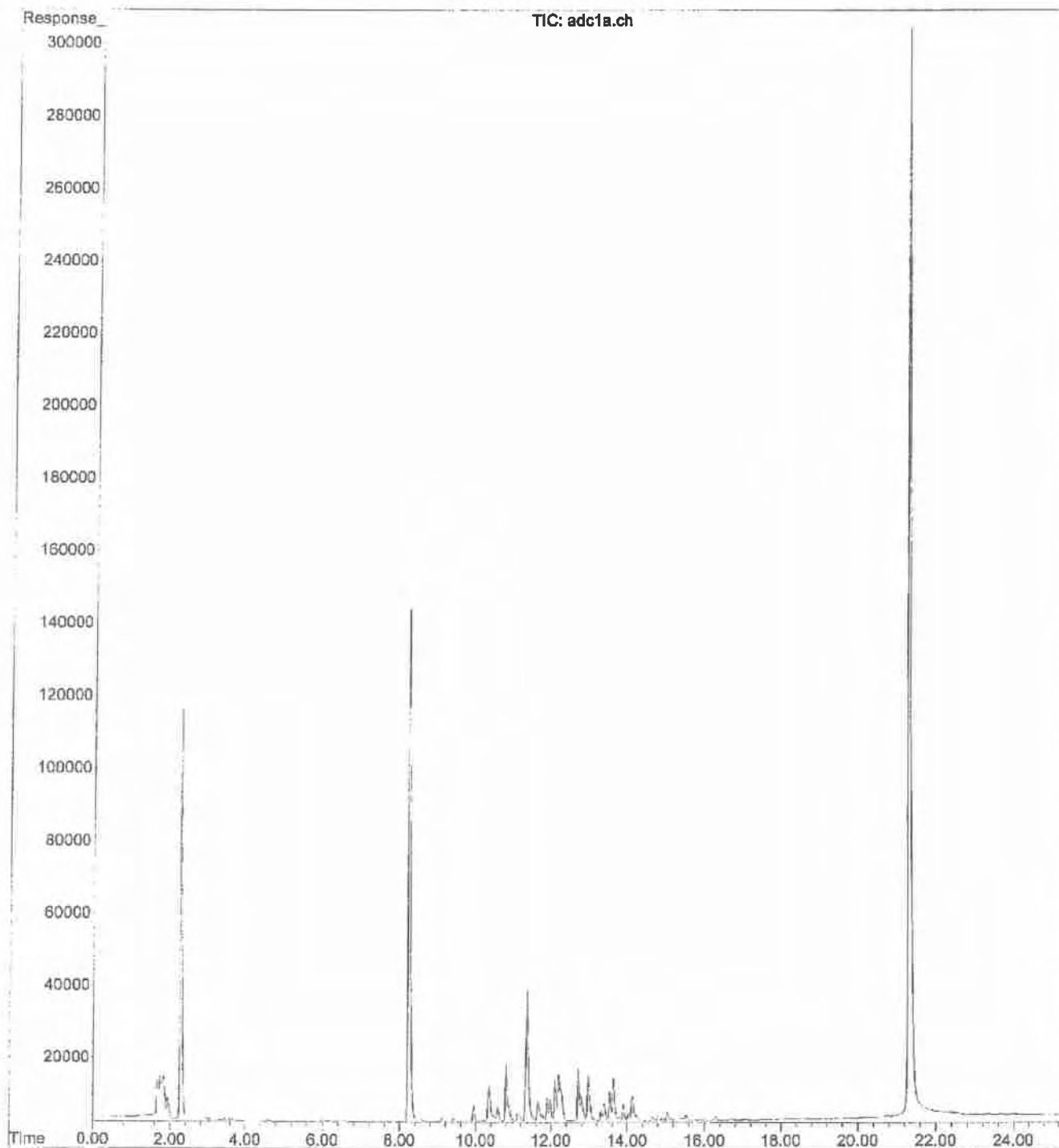
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Instrument : SVOA03
Sample Name: 10154-03 pcb x1
Misc Info :
Vial Number: 14



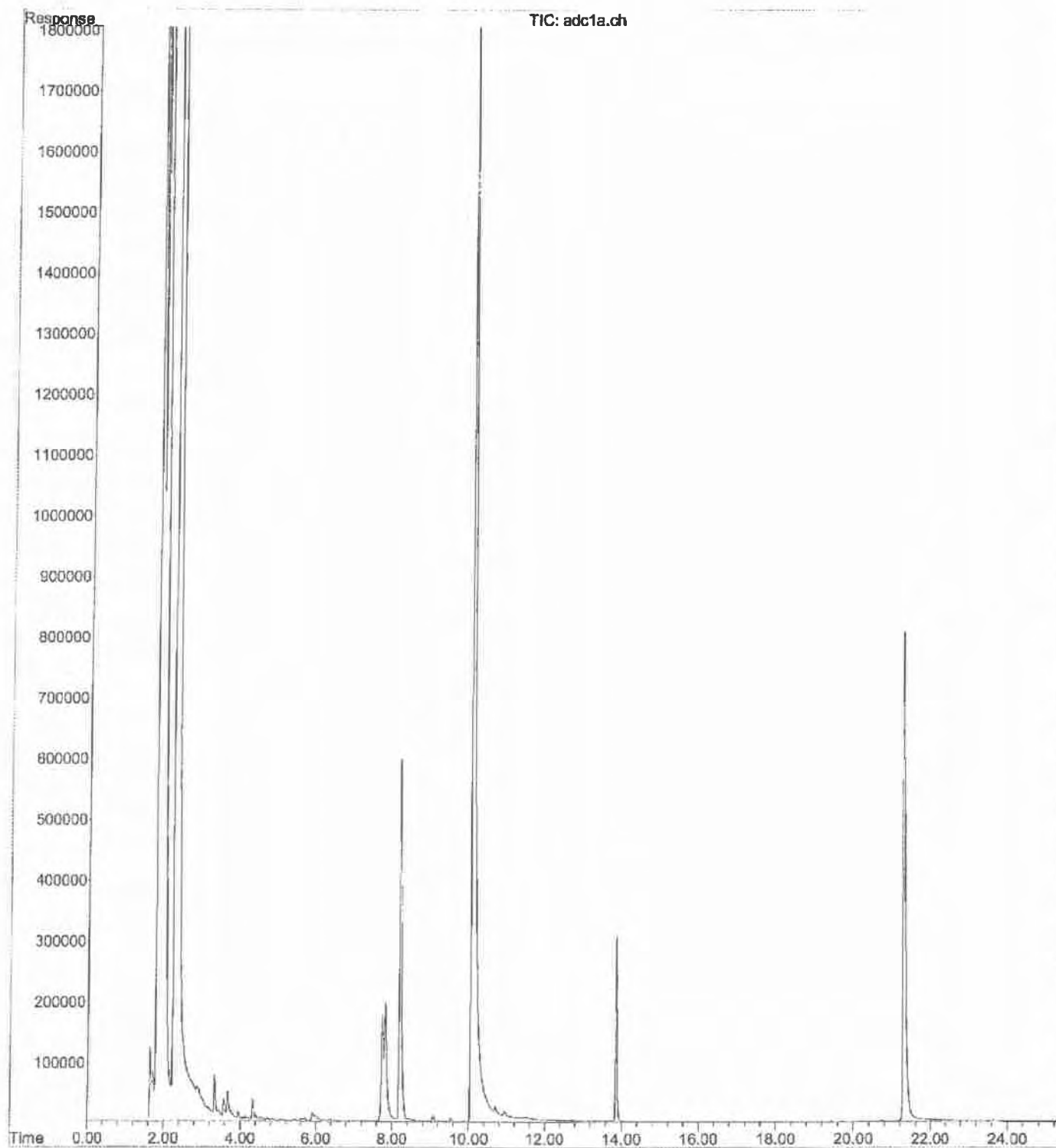
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Instrument : SVOA03
Sample Name: 10154-04 pcb x1
Misc Info :
Vial Number: 15



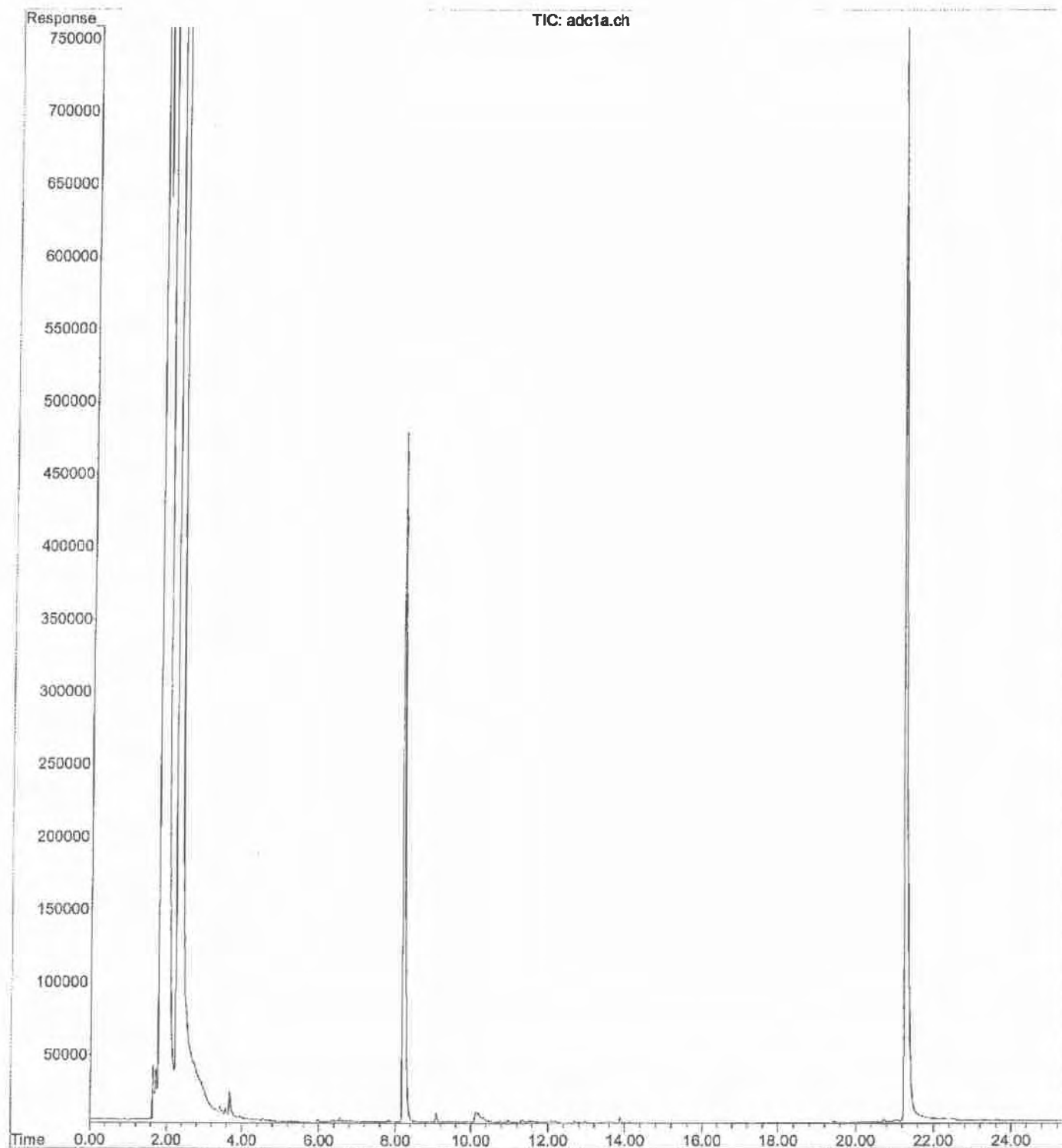
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Instrument : SVOA03
Sample Name: 10154-05rr pcbx5
Misc Info :
Vial Number: 10



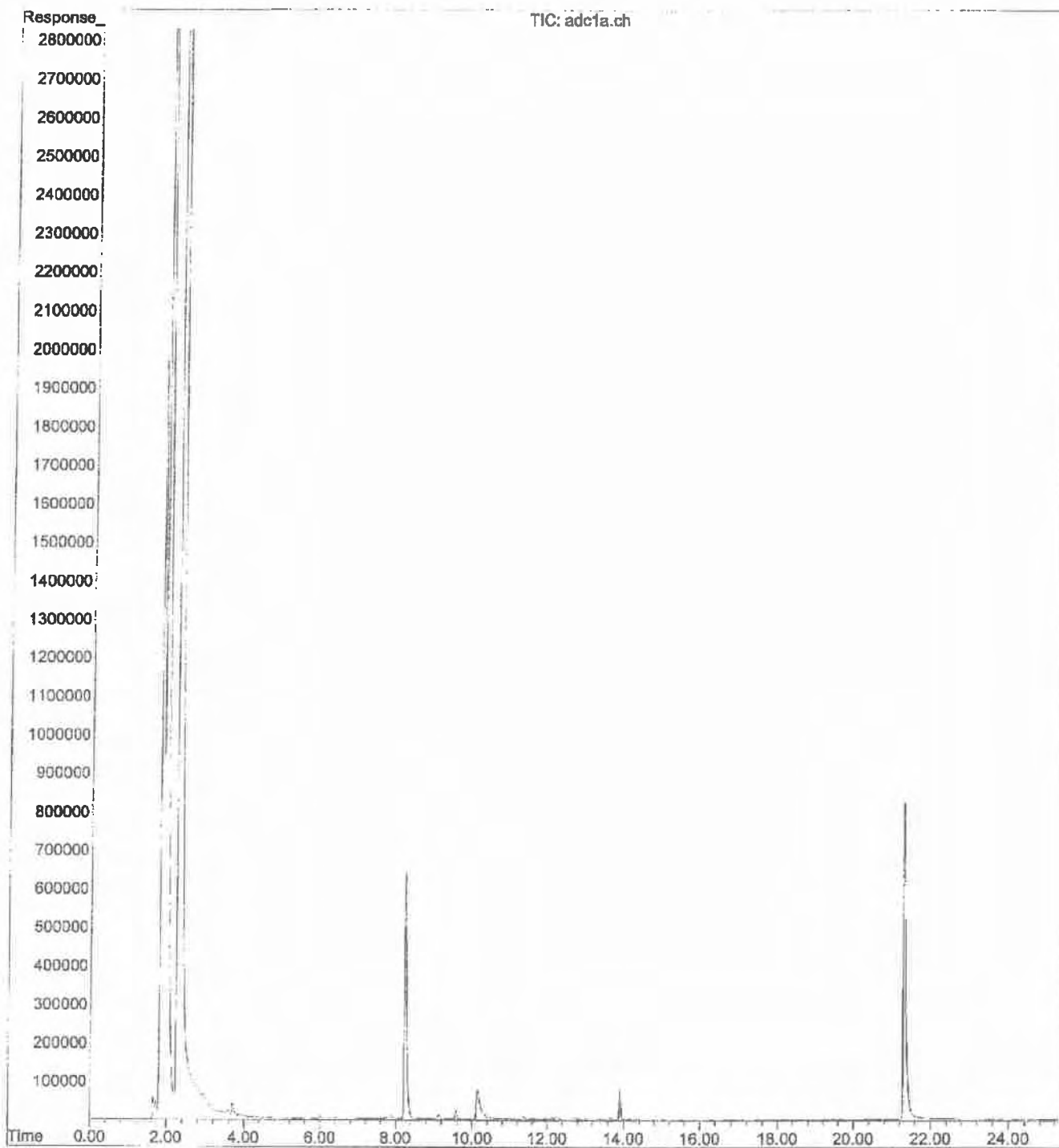
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Instrument : SVOA03
Sample Name: 10154-07 pcb x1
Misc Info :
Vial Number: 17



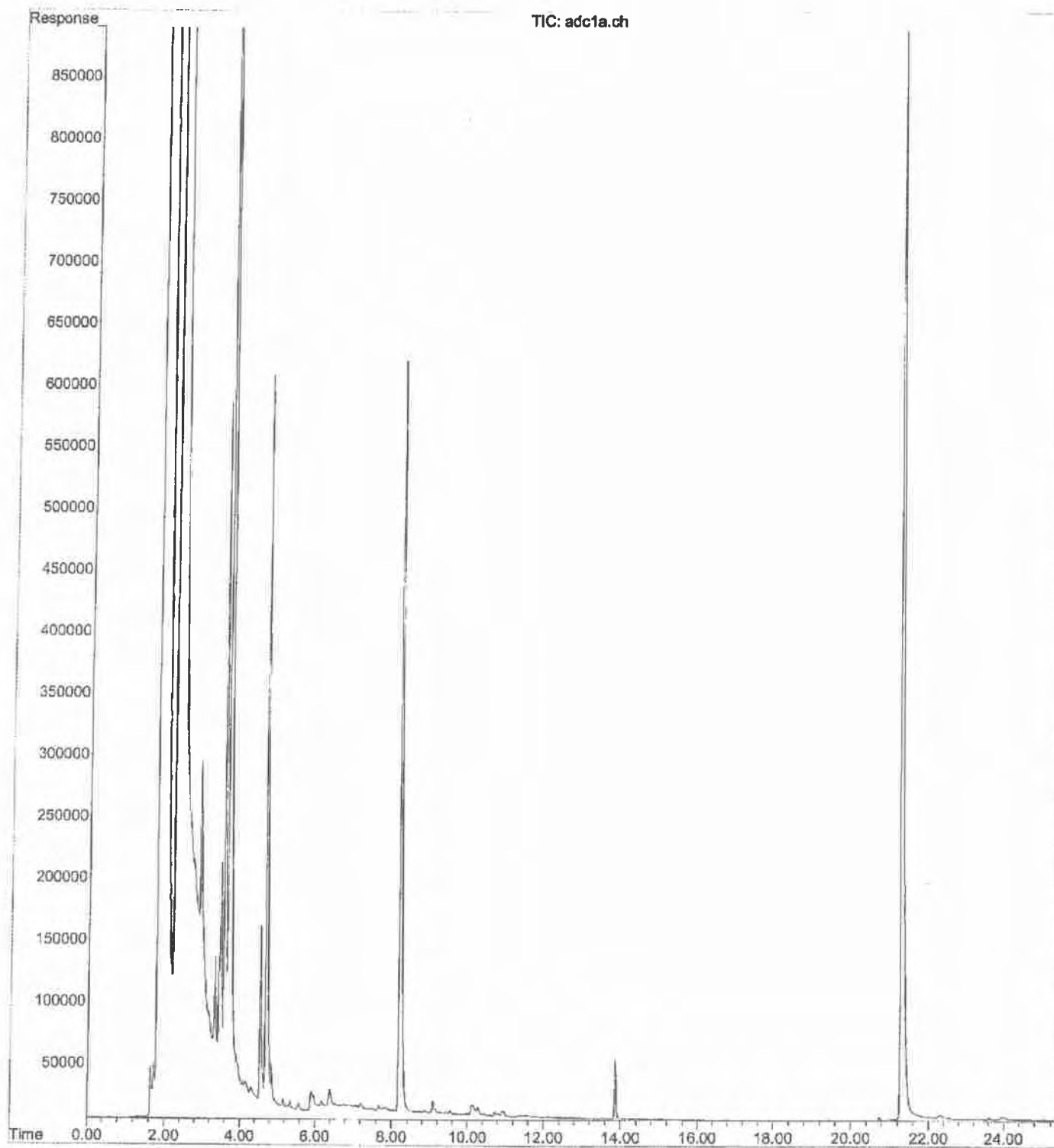
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Misc Info :
Vial Number: 18



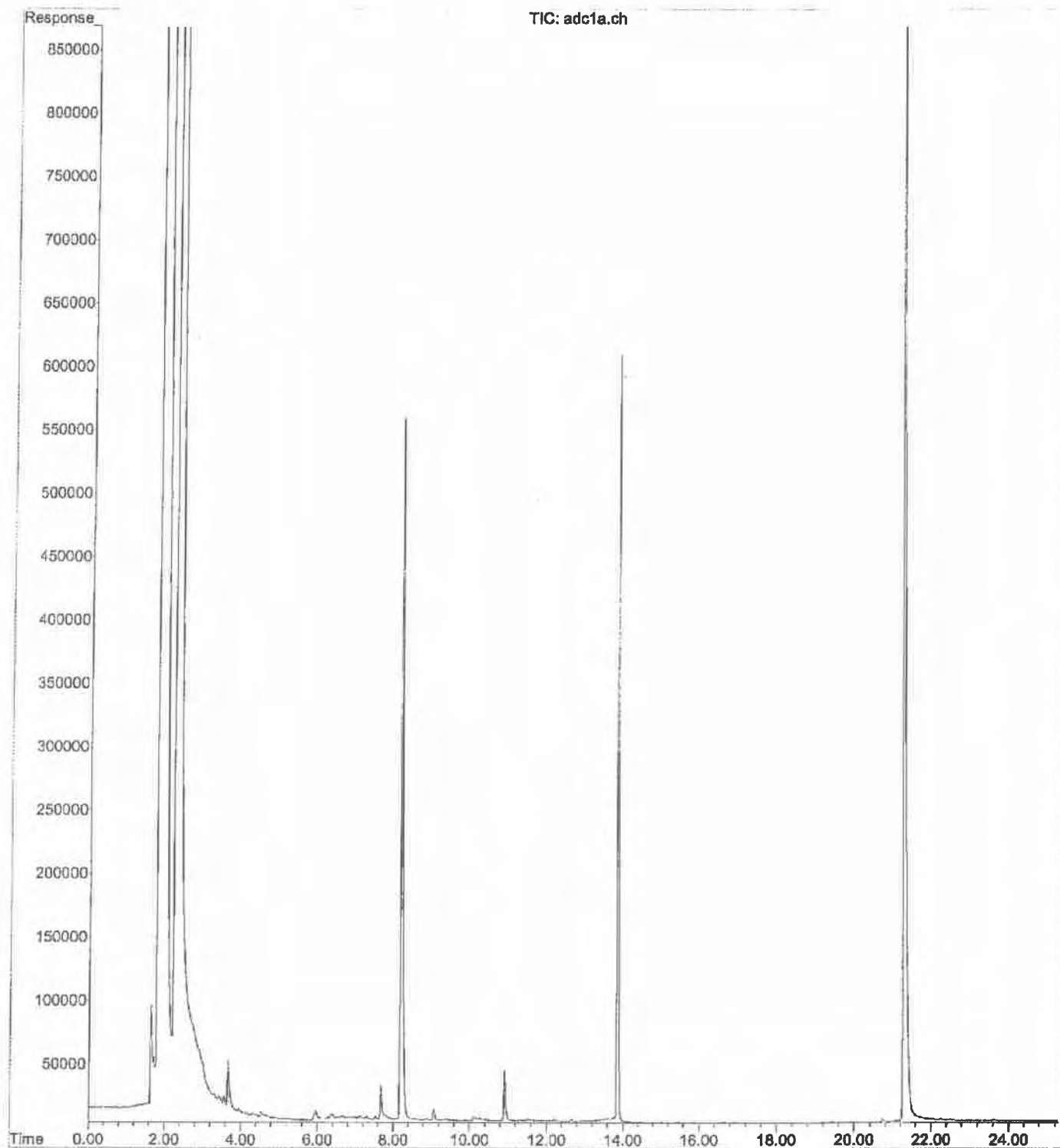
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Instrument : SVOA03
Sample Name: 10154-09rr pcbx1
Misc Info :
Vial Number: 7



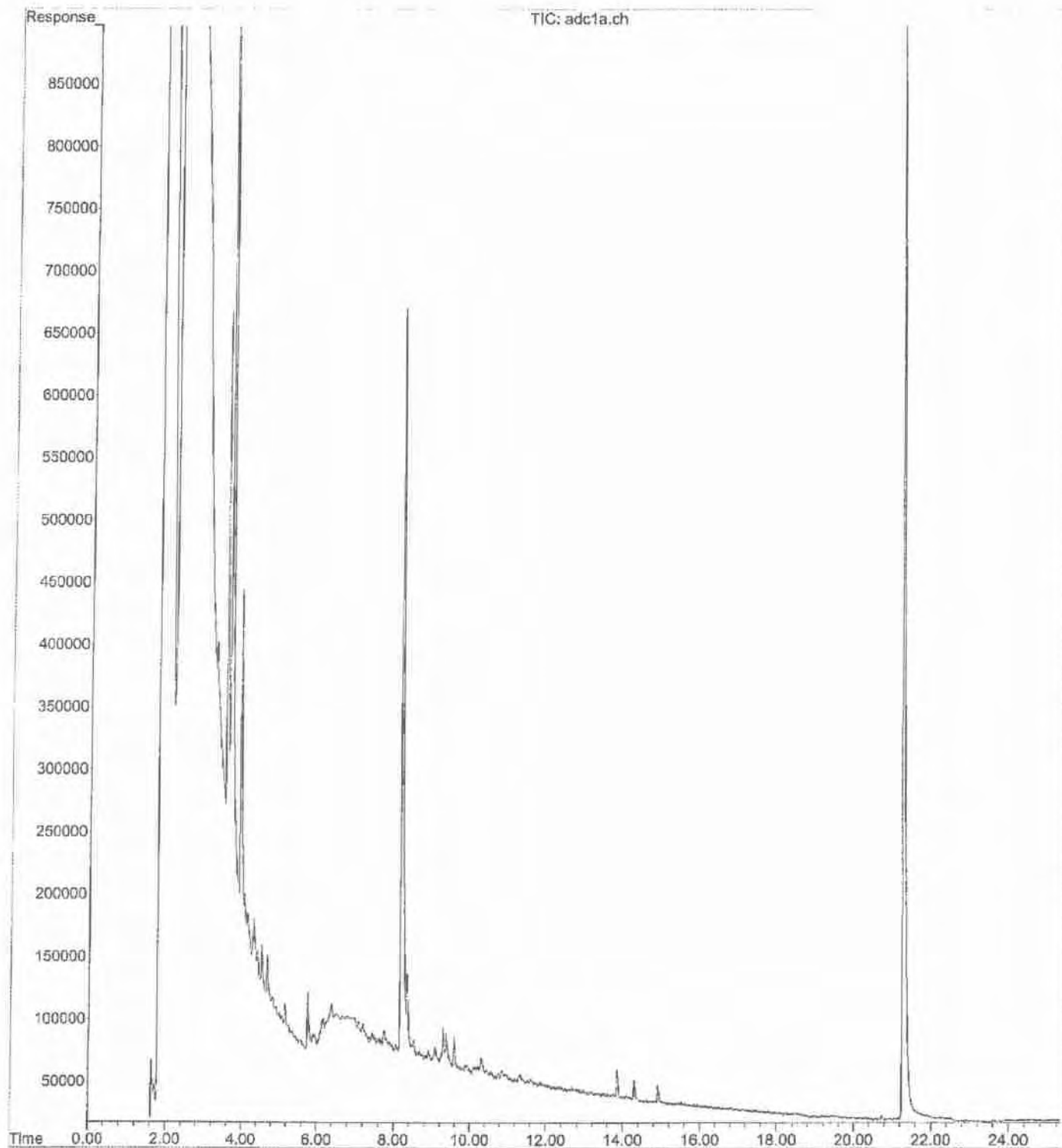
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Instrument : SVOA03
Sample Name: 10154-10 pcb x1
Misc Info :
Vial Number: 22



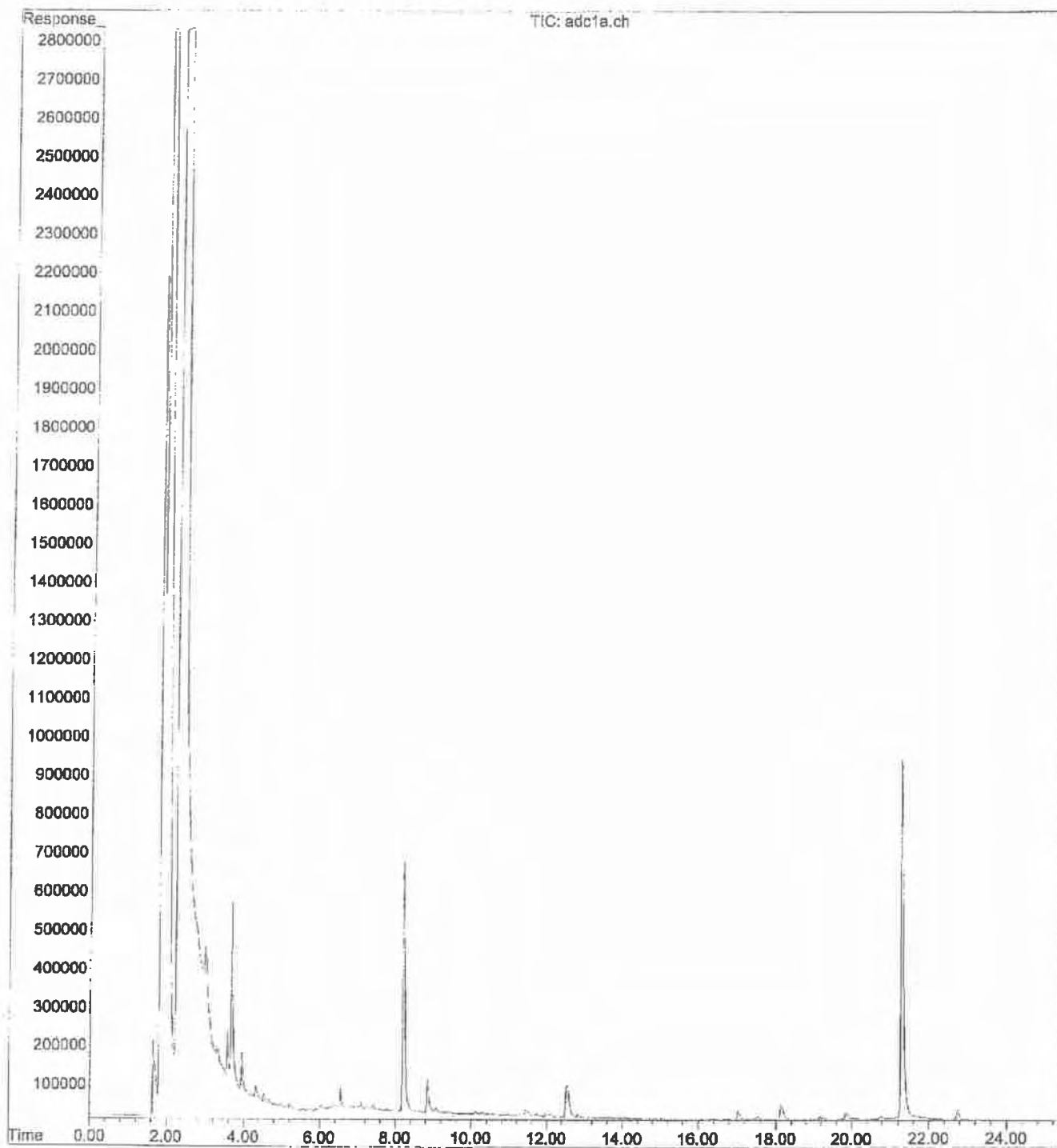
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Sample Name: 10154-13 pcb x1
Misc Info :
Vial Number: 23



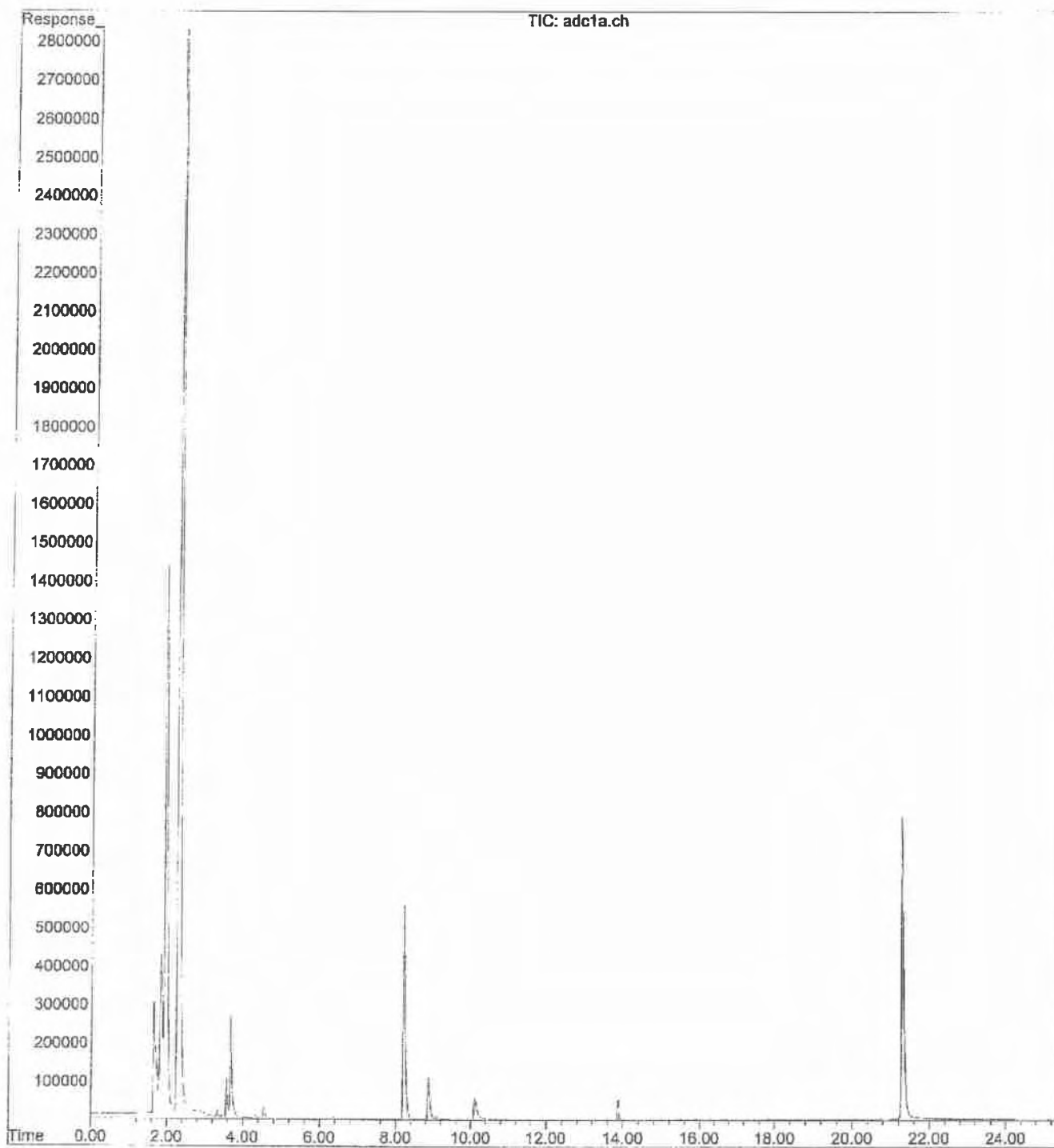
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Instrument : SVOA03
Sample Name: 10154-14 pcb x1
Misc Info :
Vial Number: 24



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Operator : AJD
Acquired : 18 Apr 2006 11:20 am using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10154-15 pcb x1
Misc Info :
Vial Number: 7



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Operator : AJD
Acquired : 18 Apr 2006 02:12 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10154-16 pcb x1
Misc Info :
Vial Number: 13



Resource Laboratories, LLC

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CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

10154

PAGE 1 OF 2

| | | | |
|---|----------|---|--|
| Company Name: ES&H | | Phone #: 603-285-9700 | |
| Company Address: 154 W. Main St. Boston, MA | | FAX #: 603-285-9807 | |
| Project Manager: J. Callahan | | Site Location (City, State): Boston, MA | |
| Invoice To: ES&H | | Project ID / Name: DND LEAS 2006-050 | |
| Protocol: RCRA SDWA NPDES MCP NHDES OTHER | | Preservation Method | |
| Lab Sample ID | Field ID | # CONTAINERS | MATRIX |
| (Lab Use Only) | | | WATER SOLID OTHER |
| | | | HCl HNO ₃ H ₂ SO ₄ NaOH MeOH OTHER (Specify) |
| | | | DATE TIME SAMPLER |
| 10154 | ESM11 | 4 | X |
| 02 | ESM11 | 1 | X |
| 03 | ESM12 | 4 | X |
| 04 | ESM12 | 4 | X |
| 05 | ESM14 | 8 | X |
| 06 | ESM14 | 3 | X |
| 07 | ESM13 | 8 | X |
| 08 | ESM13 | 8 | X |
| 09 | ESM13 | 8 | X |
| 10 | ESM13 | 5 | X |
| 11 | ESM13 | 1 | X |
| SPECIAL INSTRUCTIONS | | | |
| Metals samples field filtered | | | |
| M3/M3D from ESM-3 | | | |
| * Add steel in ESM-3 per J. Callahan | | | |
| REPORTING INSTRUCTIONS | | | |
| <input type="checkbox"/> FAX <input type="checkbox"/> OTHER (specify) <input type="checkbox"/> PDF <input type="checkbox"/> Excel Spreadsheet | | | |
| CUSTODY RECORD | | RECEIVED ON ICE | |
| Relinquished by: J. Callahan | | YES <input checked="" type="checkbox"/> NO <input type="checkbox"/> | |
| Relinquished by: J. Callahan | | TEMPERATURE 5 °C | |
| Date: 7/1/06 Time: 13:55 | | Date: 7/1/06 Time: 13:55 | |
| Received by: J. Callahan | | Received by: J. Callahan | |
| Date: 7/1/06 Time: 13:55 | | Date: 7/1/06 Time: 13:55 | |
| Received by Laboratory: J. Callahan | | Received by Laboratory: J. Callahan | |
| Date: 7/1/06 Time: 13:55 | | Date: 7/1/06 Time: 13:55 | |

- ☐ VOC 8260-NH List ☒ MADEP VPH ☐ MEGRO
- ☒ VOC 8260 ☐ VOC 80156RO ☐ VOC 624
- ☐ VOC 8260 BTEX, MIBE, Naphthalene only
- ☐ VOC 524.2 ☐ VOC 524.2 NH List
- ☐ TPH Fingerprint ☐ MEDRO ☐ DRO 8015
- ☐ 8270PAH ☐ 8270ABN ☐ 625
- ☒ 8082 PCB ☐ 8081 Pesticides ☐ 608
- ☐ O&G 1664 ☐ O&G SM5520F
- ☐ pH ☐ BOD ☐ Conductivity
- ☐ TSS ☐ TDS ☐ TS
- ☒ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals
- ☐ Total Metals-list ☐ Dissolved Metals-list
- ☐ Ammonia ☐ COD
- ☐ T-Phosphatid ☐ Phenol
- ☐ Cyanide ☐ Sulfide
- ☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride
- ☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitibility/FP
- ☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC
- ☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)
- ☐ Standard Drinking Water Test ☐ Bacteria P/A

Grab (G) or Composite (C)

RL

Resource Laboratories, LLC
124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST

ANALYSIS REQUEST

10154

PAGE 2 OF 2

Company Name:

ES+M

Phone #: 508-285-9700

Company Address:

134 W. Main St. North MA 02855

FAX #: 508-285-9757

Site Location (City, State):

Project Manager:

J. Callahan

Project ID / Name:

Boston MA

Invoice To:

ES+M

Protocol:

ROA SDWA NPDES
MGP NHDES OTHER

2006-056

Lab Sample ID

Field ID

CONTAINERS

Matrix

Preservation Method

Sampling

SAMPLER

☐ VOC 8260-NH List ☐ MADEP VPH ☐ MEGRO

☒ VOC 8260 ☐ VOC 8015 GPO ☒ 8260B

☐ VOC 8260 BTEX, MIB, Naphthalene only

☐ VOC 524.2 ☐ VOC 524.2 NH List

☐ TPH Fingerprint ☐ MEDRO ☐ DRO 6015 ☒ FPH

☐ 8270PAH ☐ 8270ABN ☐ 625

☒ 8082 PCB ☐ 8081 Pesticides ☐ 608

☐ O&G 1664 ☐ O&G SM552DF

☐ pH ☐ BOD ☐ Conductivity

☐ TSS ☐ IUS ☐ TS

☒ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals

☐ Total Metals-list ☐ Dissolved Metals-list

☐ Ammonia ☐ COD

☐ T-Phosphate ☐ Phenol

☐ Cyanide ☐ Sulfide

☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride

☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitability/FP

☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC

☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)

☐ Standard Drinking Water Test ☐ Bacteria P/A

Grab (G) or Composite (C)

TAT REQUESTED

Priority (24 hr)

Expected (48 hr)

10 Business Days

Other

Quote #

PO #

SPECIAL INSTRUCTIONS

Metals Samples Filtered in field

REPORTING INSTRUCTIONS

☐ FAX ☐ OTHER (specify)

☐ PDF ☐ Excel Spreadsheet

RECEIVED ON ICE

YES ☒ NO ☐

TEMPERATURE

5 °C

Lab Use Only

CUSTODY RECORD

Relinquished by Sampler:

Relinquished by:

Date 4/7/06 Time 13:55

Received by:

Received by:

Date 4/7/06 Time 13:55

Relinquished by:

Date Time

Received by Laboratory:

Way Bill#:

Date Time

Resource Laboratories, LLC

124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

ANALYSIS REQUEST

Company Name:

ES417

Phone #: 508-285-7700

Company Address:

164 W. Main St. Boston MA

FAX #: 508-285-9957

Site Location (City, State):

Project Manager:

J. Callahan

Project ID / Name:

D202 14-15 2006-0510

Invoice To:

ES417

Protocol:

RCRA SDWA NPDES
MGP NIDES OTHER

| Lab Sample ID (Lab Use Only) | Field ID | # CONTAINERS | Matrix | | | Preservation Method | | | | | Sampling | |
|---------------------------------|----------|--------------|--------|-------|-------|---------------------|------------------|--------------------------------|------|------|---------------|------|
| | | | WATER | SOLID | OTHER | HCl | HNO ₃ | H ₂ SO ₄ | NaOH | MeOH | DATE | TIME |
| ES417 | 4 | 4 | X | | | 3 | 1 | | | | 4/16/06 11:30 | TC |
| ES417 | 1 | 1 | X | | | | | | | | 4/17/06 10:30 | AF |
| ES417 | 4 | 4 | X | | | 3 | | | | | 4/16/06 2:00 | AF |
| ES417 | 4 | 4 | X | | | 3 | | | | | 4/16/06 2:45 | AF |
| ES417 | 3 | 3 | X | | | 3 | | | | | 4/16/06 14:45 | AF |
| ES417 | 3 | 3 | X | | | 3 | | | | | 4/16/06 14:45 | AF |
| ES417 | 3 | 3 | X | | | 3 | | | | | 4/16/06 14:45 | AF |
| ES417 | 3 | 3 | X | | | 3 | | | | | 4/16/06 13:50 | AF |
| ES417 | 3 | 3 | X | | | 3 | | | | | 4/16/06 9:45 | AF |
| ES417 | 1 | 1 | X | | | | | | | | 4/16/06 9:30 | AF |

SPECIAL INSTRUCTIONS

REPORTING INSTRUCTIONS
□ FAX □ OTHER (specify)
□ PDF □ Excel Spreadsheet

RECEIVED ON ICE □ YES □ NO
TEMPERATURE _____ °C

CUSTODY RECORD

Retrieved by Sampler:
Relinquished by:

Date: 4/16/06 Time: 13:35

Received by:

Date: 4/17/06 Time: 12:00

Received by Laboratory:
Way Bill:

Date: 4/17/06 Time: 12:00

- VOC 8260-NH List □ MADEP VPH □ MEGRO
□ VOC 8260 □ VOC 8015GRO □ VOC 624
□ VOC 8260 BTEX, MIBK, Naphthalene only
□ VOC 524.2 □ VOC 524.2 NH List
□ TPH Fingerprint □ MEDRO □ T.DRO 8015 □ EPH
□ B270PAH □ B270A2N □ B25
□ B282 PCB □ B281 Pesticides □ B18
□ O&G 1664 □ O&G SM5520F
□ pH □ BOD □ Conductivity
□ TSS □ TDS □ TS
□ RCRA Metals □ Priority Pollutant Metals □ TAL Metals
□ Total Metals list □ Dissolved Metals-list
□ Ammonia □ COD
□ T-Phosphate □ Phenol
□ Cyanide □ Sulfide
□ NRrate □ NRite □ Orbo P □ Sulfate □ Bromide □ Chloride
□ Corrosivity □ Relative CN □ Reactive S □ Ignitability/FP
□ TCLP Metals □ TCLP VOC □ TCLP SVOC
□ TCLP Pesticide □ TCLP Herbicides (subcontract)
□ Standard Drinking Water Test □ Bacteria P/A

Grab (G) or Composite (C)

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10170
Date Received: 4/11/06

Project: 2006-056 Lewis Chemical

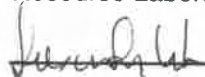
Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

4-27-06

Date

Total number of pages

28

Resource Laboratories, LLC Certifications

New Hampshire NH902
Maine NH903

Connecticut PH-0146
Massachusetts M-NH902

Lab Number: 10170-01
Sample Designation: PZ-03-S
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 20
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 40 | trans-1,3-dichloropropene | U | 40 |
| chloromethane | U | 40 | 2-hexanone | U | 200 |
| vinyl chloride | 3900 | 40 | 1,1,2-trichloroethane | U | 40 |
| bromomethane | U | 40 | 1,3-dichloropropane | U | 40 |
| chloroethane | U | 40 | tetrachloroethene | U | 40 |
| trichlorofluoromethane | U | 40 | dibromochloromethane | U | 40 |
| diethyl ether | U | 200 | 1,2-dibromoethane | U | 40 |
| acetone | U | 200 | chlorobenzene | U | 40 |
| 1,1-dichloroethene | 63 | 20 | 1,1,1,2-tetrachloroethane | U | 40 |
| methylene chloride | U | 100 | ethylbenzene | U | 40 |
| carbon disulfide | U | 40 | m&p-xylenes | 49 | 40 |
| methyl t-butyl ether (MTBE) | U | 40 | o-xylene | U | 40 |
| trans-1,2-dichloroethene | 120 | 40 | styrene | U | 40 |
| isopropyl ether (DIPE) | U | 40 | bromoform | U | 40 |
| ethyl t-butyl ether (ETBE) | U | 40 | Isopropylbenzene | U | 40 |
| 1,1-dichloroethane | 2900 | 40 | 1,1,2,2-tetrachloroethane | U | 40 |
| t-butanol (TBA) | U | 1000 | 1,2,3-trichloropropane | U | 40 |
| 2-butanone (MEK) | U | 200 | n-propylbenzene | U | 40 |
| 2,2-dichloropropane | U | 40 | bromobenzene | U | 40 |
| cis-1,2-dichloroethene | 11000 | 40 | 1,3,5-trimethylbenzene | U | 40 |
| chloroform | U | 40 | 2-chlorotoluene | U | 40 |
| bromochloromethane | U | 40 | 4-chlorotoluene | U | 40 |
| tetrahydrofuran (THF) | U | 200 | tert-butylbenzene | U | 40 |
| 1,1,1-trichloroethane | 1700 | 40 | 1,2,4-trimethylbenzene | U | 40 |
| 1,1-dichloropropene | U | 40 | sec-butylbenzene | U | 40 |
| t-amyl-methyl ether (TAME) | U | 40 | 1,3-dichlorobenzene | U | 40 |
| carbon tetrachloride | U | 40 | 4-isopropyltoluene | U | 40 |
| 1,2-dichloroethane | 410 | 40 | 1,4-dichlorobenzene | U | 40 |
| benzene | U | 40 | 1,2-dichlorobenzene | 260 | 40 |
| trichloroethene | U | 40 | n-butylbenzene | U | 40 |
| 1,2-dichloropropane | U | 40 | 1,2-dibromo-3-chloropropane | U | 40 |
| bromodichloromethane | U | 40 | 1,2,4-trichlorobenzene | U | 40 |
| dibromomethane | U | 40 | hexachlorobutadiene | U | 40 |
| 4-methyl-2-pentanone (MIBK) | U | 200 | naphthalene | U | 100 |
| cis-1,3-dichloropropene | U | 40 | 1,2,3-trichlorobenzene | U | 40 |
| toluene | 260 | 40 | 1,4-dioxane | U | 1000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dlbromofluoromethane | 96 | 78-114 | | | |
| toluene-D8 | 106 | 88-110 | | | |
| 4-bromofluorobenzene | 102 | 86-115 | | | |

U = Below quantitation limit

RL Resource Laboratories, LLC

Lab Number: 10170-02
Sample Designation: PZ-01-D
Date Sampled: 4/10/06
Date Analyzed: 4/18/06
Matrix: Water
Instrument Dilution Factor: 5
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 10 | trans-1,3-dichloropropene | U | 10 |
| chloromethane | U | 10 | 2-hexanone | U | 50 |
| vinyl chloride | 1700 D | 20 | 1,1,2-trichloroethane | U | 10 |
| bromomethane | U | 10 | 1,3-dichloropropane | U | 10 |
| chloroethane | U | 10 | tetrachloroethene | 28 | 10 |
| trichlorofluoromethane | U | 10 | dibromochloromethane | U | 10 |
| diethyl ether | U | 50 | 1,2-dibromoethane | U | 10 |
| acetone | U | 50 | chlorobenzene | U | 10 |
| 1,1-dichloroethene | 11 | 5 | 1,1,1,2-tetrachloroethane | U | 10 |
| methylene chloride | U | 30 | ethylbenzene | 36 | 10 |
| carbon disulfide | U | 10 | m&p-xylenes | 80 | 10 |
| methyl t-butyl ether (MTBE) | U | 10 | o-xylene | 33 | 10 |
| trans-1,2-dichloroethene | 20 | 10 | styrene | U | 10 |
| isopropyl ether (DIPE) | U | 10 | bromoform | U | 10 |
| ethyl t-butyl ether (ETBE) | U | 10 | isopropylbenzene | U | 10 |
| 1,1-dichloroethane | 890 | 10 | 1,1,2,2-tetrachloroethane | U | 10 |
| t-butanol (TBA) | U | 300 | 1,2,3-trichloropropane | U | 10 |
| 2-butanone (MEK) | U | 50 | n-propylbenzene | U | 10 |
| 2,2-dichloropropane | U | 10 | bromobenzene | U | 10 |
| cis-1,2-dichloroethene | 2400 | 10 | 1,3,5-trimethylbenzene | U | 10 |
| chloroform | U | 10 | 2-chlorotoluene | U | 10 |
| bromochloromethane | U | 10 | 4-chlorotoluene | U | 10 |
| tetrahydrofuran (THF) | U | 50 | tert-butylbenzene | U | 10 |
| 1,1,1-trichloroethane | 990 | 10 | 1,2,4-trimethylbenzene | 10 | 10 |
| 1,1-dichloropropene | U | 10 | sec-butylbenzene | U | 10 |
| t-amyl-methyl ether (TAME) | U | 10 | 1,3-dichlorobenzene | U | 10 |
| carbon tetrachloride | U | 10 | 4-isopropyltoluene | U | 10 |
| 1,2-dichloroethane | 57 | 10 | 1,4-dichlorobenzene | U | 10 |
| benzene | 11 | 10 | 1,2-dichlorobenzene | 45 | 10 |
| trichloroethene | 33 | 10 | n-butylbenzene | U | 10 |
| 1,2-dichloropropane | U | 10 | 1,2-dibromo-3-chloropropane | U | 10 |
| bromodichloromethane | U | 10 | 1,2,4-trichlorobenzene | U | 10 |
| dibromomethane | U | 10 | hexachlorobutadiene | U | 10 |
| 4-methyl-2-pentanone (MIBK) | U | 50 | naphthalene | U | 30 |
| cis-1,3-dichloropropene | U | 10 | 1,2,3-trichlorobenzene | U | 10 |
| toluene | 670 | 10 | 1,4-dioxane | U | 300 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 92 | 78-114 | | | |
| toluene-D8 | 103 | 88-110 | | | |
| 4-bromofluorobenzene | 105 | 86-115 | | | |

D = Result obtained from a re-analysis at a dilution.
U = Below quantitation limit

Lab Number: 10170-03
Sample Designation: PZ-01-S
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 50
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 100 | trans-1,3-dichloropropene | U | 100 |
| chloromethane | U | 100 | 2-hexanone | U | 500 |
| vinyl chloride | 6400 | 100 | 1,1,2-trichloroethane | U | 100 |
| bromomethane | U | 100 | 1,3-dichloropropane | U | 100 |
| chloroethane | U | 100 | tetrachloroethene | U | 100 |
| trichlorofluoromethane | U | 100 | dibromochloromethane | U | 100 |
| diethyl ether | U | 500 | 1,2-dibromoethane | U | 100 |
| acetone | U | 500 | chlorobenzene | U | 100 |
| 1,1-dichloroethene | 92 | 50 | 1,1,1,2-tetrachloroethane | U | 100 |
| methylene chloride | U | 300 | ethylbenzene | 330 | 100 |
| carbon disulfide | U | 100 | m&p-xylenes | 940 | 100 |
| methyl t-butyl ether (MTBE) | U | 100 | o-xylene | 290 | 100 |
| trans-1,2-dichloroethene | 200 | 100 | styrene | U | 100 |
| isopropyl ether (DIPE) | U | 100 | bromoform | U | 100 |
| ethyl t-butyl ether (ETBE) | U | 100 | isopropylbenzene | U | 100 |
| 1,1-dichloroethane | 4000 | 100 | 1,1,1,2-tetrachloroethane | U | 100 |
| t-butanol (TBA) | U | 3000 | 1,2,3-trichloropropane | U | 100 |
| 2-butanone (MEK) | U | 500 | n-propylbenzene | U | 100 |
| 2,2-dichloropropane | U | 100 | bromobenzene | U | 100 |
| cis-1,2-dichloroethene | 24000 | 100 | 1,3,5-trimethylbenzene | U | 100 |
| chloroform | U | 100 | 2-chlorotoluene | U | 100 |
| bromochloromethane | U | 100 | 4-chlorotoluene | U | 100 |
| tetrahydrofuran (THF) | U | 500 | tert-butylbenzene | U | 100 |
| 1,1,1-trichloroethane | 4700 | 100 | 1,2,4-trimethylbenzene | U | 100 |
| 1,1-dichloropropene | U | 100 | sec-butylbenzene | U | 100 |
| t-amyl-methyl ether (TAME) | U | 100 | 1,3-dichlorobenzene | U | 100 |
| carbon tetrachloride | U | 100 | 4-isopropyltoluene | U | 100 |
| 1,2-dichloroethane | 100 | 100 | 1,4-dichlorobenzene | U | 100 |
| benzene | U | 100 | 1,2-dichlorobenzene | 200 | 100 |
| trichloroethene | 200 | 100 | n-butylbenzene | U | 100 |
| 1,2-dichloropropane | U | 100 | 1,2-dibromo-3-chloropropane | U | 100 |
| bromodichloromethane | U | 100 | 1,2,4-trichlorobenzene | U | 100 |
| dibromomethane | U | 100 | hexachlorobutadiene | U | 100 |
| 4-methyl-2-pentanone (MIBK) | U | 500 | naphthalene | U | 300 |
| cis-1,3-dichloropropene | U | 100 | 1,2,3-trichlorobenzene | U | 100 |
| toluene | 24000 | 100 | 1,4-dioxane | U | 3000 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 97 | 78-114 | | | |
| toluene-D8 | 106 | 88-110 | | | |
| 4-bromofluorobenzene | 103 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10170-04
Sample Designation: PZ-02-S
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 50
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 100 | trans-1,3-dichloropropene | U | 100 |
| chloromethane | U | 100 | 2-hexanone | U | 500 |
| vinyl chloride | 740 | 100 | 1,1,2-trichloroethane | 140 | 100 |
| bromomethane | U | 100 | 1,3-dichloropropane | U | 100 |
| chloroethane | U | 100 | tetrachloroethene | 240 | 100 |
| trichlorofluoromethane | 200 | 100 | dibromochloromethane | U | 100 |
| diethyl ether | U | 500 | 1,2-dibromoethane | U | 100 |
| acetone | U | 500 | chlorobenzene | U | 100 |
| 1,1-dichloroethene | 740 | 50 | 1,1,1,2-tetrachloroethane | U | 100 |
| methylene chloride | 5800 | 300 | ethylbenzene | 350 | 100 |
| carbon disulfide | U | 100 | m&p-xylenes | 890 | 100 |
| methyl t-butyl ether (MTBE) | U | 100 | o-xylene | 310 | 100 |
| trans-1,2-dichloroethene | 240 | 100 | styrene | U | 100 |
| isopropyl ether (DIPE) | U | 100 | bromoform | U | 100 |
| ethyl t-butyl ether (ETBE) | U | 100 | isopropylbenzene | U | 100 |
| 1,1-dichloroethane | 3200 | 100 | 1,1,2,2-tetrachloroethane | U | 100 |
| t-butanol (TBA) | U | 3000 | 1,2,3-trichloropropane | U | 100 |
| 2-butanone (MEK) | U | 500 | n-propylbenzene | U | 100 |
| 2,2-dichloropropane | U | 100 | bromobenzene | U | 100 |
| cis-1,2-dichloroethene | 48000 D | 400 | 1,3,5-trimethylbenzene | U | 100 |
| chloroform | U | 100 | 2-chlorotoluene | U | 100 |
| bromochloromethane | U | 100 | 4-chlorotoluene | U | 100 |
| tetrahydrofuran (THF) | U | 500 | tert-butylbenzene | U | 100 |
| 1,1,1-trichloroethane | 37000 | 100 | 1,2,4-trimethylbenzene | 180 | 100 |
| 1,1-dichloropropene | U | 100 | sec-butylbenzene | U | 100 |
| t-amyl-methyl ether (TAME) | U | 100 | 1,3-dichlorobenzene | U | 100 |
| carbon tetrachloride | U | 100 | 4-isopropyltoluene | U | 100 |
| 1,2-dichloroethane | 3300 | 100 | 1,4-dichlorobenzene | U | 100 |
| benzene | 220 | 100 | 1,2-dichlorobenzene | 120 | 100 |
| trichloroethene | 220 | 100 | n-butylbenzene | U | 100 |
| 1,2-dichloropropane | U | 100 | 1,2-dibromo-3-chloropropane | U | 100 |
| bromodichloromethane | U | 100 | 1,2,4-trichlorobenzene | U | 100 |
| dibromomethane | U | 100 | hexachlorobutadiene | U | 100 |
| 4-methyl-2-pentanone (MIBK) | 560 | 500 | naphthalene | U | 300 |
| cis-1,3-dichloropropene | U | 100 | 1,2,3-trichlorobenzene | U | 100 |
| toluene | 10000 | 100 | 1,4-dioxane | U | 3000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 98 | 78-114 | | | |
| toluene-D8 | 105 | 88-110 | | | |
| 4-bromofluorobenzene | 104 | 86-115 | | | |

D = Result obtained from a re-analysis at a dilution.
U = Below quantitation limit

Lab Number: 10170-05
Sample Designation: PZ-02-D
Date Sampled: 4/10/06
Date Analyzed: 4/18/06
Matrix: Water
Instrument Dilution Factor: 10
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 20 | trans-1,3-dichloropropene | U | 20 |
| chloromethane | U | 20 | 2-hexanone | U | 100 |
| vinyl chloride | 130 | 20 | 1,1,2-trichloroethane | U | 20 |
| bromomethane | U | 20 | 1,3-dichloropropane | U | 20 |
| chloroethane | U | 20 | tetrachloroethene | 53 | 20 |
| trichlorofluoromethane | U | 20 | dibromochloromethane | U | 20 |
| diethyl ether | U | 100 | 1,2-dibromoethane | U | 20 |
| acetone | U | 100 | chlorobenzene | U | 20 |
| 1,1-dichloroethene | 50 | 10 | 1,1,1,2-tetrachloroethane | U | 20 |
| methylene chloride | 150 | 50 | ethylbenzene | 34 | 20 |
| carbon disulfide | U | 20 | m&p-xylenes | 56 | 20 |
| methyl t-butyl ether (MTBE) | U | 20 | o-xylene | 22 | 20 |
| trans-1,2-dichloroethene | U | 20 | styrene | U | 20 |
| isopropyl ether (DIPE) | U | 20 | bromoform | U | 20 |
| ethyl t-butyl ether (ETBE) | U | 20 | isopropylbenzene | U | 20 |
| 1,1-dichloroethane | 260 | 20 | 1,1,2,2-tetrachloroethane | U | 20 |
| t-butanol (TBA) | U | 500 | 1,2,3-trichloropropane | U | 20 |
| 2-butanone (MEK) | U | 100 | n-propylbenzene | U | 20 |
| 2,2-dichloropropane | U | 20 | bromobenzene | U | 20 |
| cis-1,2-dichloroethene | 3300 | 20 | 1,3,5-trimethylbenzene | U | 20 |
| chloroform | U | 20 | 2-chlorotoluene | U | 20 |
| bromochloromethane | U | 20 | 4-chlorotoluene | U | 20 |
| tetrahydrofuran (THF) | U | 100 | tert-butylbenzene | U | 20 |
| 1,1,1-trichloroethane | 1500 | 20 | 1,2,4-trimethylbenzene | 27 | 20 |
| 1,1-dichloropropene | U | 20 | sec-butylbenzene | U | 20 |
| t-amyl-methyl ether (TAME) | U | 20 | 1,3-dichlorobenzene | U | 20 |
| carbon tetrachloride | U | 20 | 4-isopropyltoluene | U | 20 |
| 1,2-dichloroethane | 150 | 20 | 1,4-dichlorobenzene | U | 20 |
| benzene | 23 | 20 | 1,2-dichlorobenzene | U | 20 |
| trichloroethene | 62 | 20 | n-butylbenzene | U | 20 |
| 1,2-dichloropropane | U | 20 | 1,2-dibromo-3-chloropropane | U | 20 |
| bromodichloromethane | U | 20 | 1,2,4-trichlorobenzene | U | 20 |
| dibromomethane | U | 20 | hexachlorobutadiene | U | 20 |
| 4-methyl-2-pentanone (MIBK) | U | 100 | naphthalene | U | 50 |
| cis-1,3-dichloropropene | U | 20 | 1,2,3-trichlorobenzene | U | 20 |
| toluene | 500 | 20 | 1,4-dioxane | U | 500 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 96 | 78-114 | | | |
| toluene-D8 | 103 | 88-110 | | | |
| 4-bromofluorobenzene | 103 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10170-06
Sample Designation: PZ-03-D
Date Sampled: 4/10/06
Date Analyzed: 4/18/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | 92 | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | 280 | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | 5 | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | 4 | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | 93 | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | 580 | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | 180 | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | 32 | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | 63 | 2 |
| trichloroethene | 73 | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | 3 | 2 | 1,4-dioxane | U | 50 |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 96 | 78-114 | | | |
| toluene-D8 | 105 | 88-110 | | | |
| 4-bromofluorobenzene | 106 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10170-01
 Sample Designation: PZ-03-S
 Date Sampled: 4/10/06
 Date Extracted: 4/13/06
 Date Analyzed: 4/19/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.5 |
| PCB-1242 | 1.6 | 0.5 |
| PCB-1221 | U | 0.5 |
| PCB-1232 | U | 0.5 |
| PCB-1248 | U | 0.5 |
| PCB-1254 | U | 0.5 |
| PCB-1260 | U | 0.5 |
| PCB-1262 | U | 0.5 |
| PCB-1268 | U | 0.5 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 93 | 30-150 |
| Decachlorobiphenyl | 109 | 30-150 |

U = Below quantitation limit

Lab Number: 10170-04
 Sample Designation: PZ-02-S
 Date Sampled: 4/10/06
 Date Extracted: 4/13/06
 Date Analyzed: 4/20/06
 Matrix: Water
 Dilution Factor: 50
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 20 |
| PCB-1242 | 95 | 20 |
| PCB-1221 | U | 20 |
| PCB-1232 | U | 20 |
| PCB-1248 | U | 20 |
| PCB-1254 | U | 20 |
| PCB-1260 | U | 20 |
| PCB-1262 | U | 20 |
| PCB-1268 | U | 20 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | DOR | 30-150 |
| Decachlorobiphenyl | DOR | 30-150 |

DOR = Diluted out of range.
 U = Below quantitation limit

Lab Number: 10170-05
 Sample Designation: PZ-02-D
 Date Sampled: 4/10/06
 Date Extracted: 4/13/06
 Date Analyzed: 4/20/06
 Matrix: Water
 Dilution Factor: 5
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 2 |
| PCB-1242 | 9 | 2 |
| PCB-1221 | U | 2 |
| PCB-1232 | U | 2 |
| PCB-1248 | U | 2 |
| PCB-1254 | U | 2 |
| PCB-1260 | U | 2 |
| PCB-1262 | U | 2 |
| PCB-1268 | U | 2 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 90 | 30-150 |
| Decachlorobiphenyl | 139 | 30-150 |

U = Below quantitation limit

Lab Number: 10170-06
 Sample Designation: PZ-03-D
 Date Sampled: 4/10/06
 Date Extracted: 4/13/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.6 |
| PCB-1242 | U | 0.6 |
| PCB-1221 | U | 0.6 |
| PCB-1232 | U | 0.6 |
| PCB-1248 | U | 0.6 |
| PCB-1254 | U | 0.6 |
| PCB-1260 | U | 0.6 |
| PCB-1262 | U | 0.6 |
| PCB-1268 | U | 0.6 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 69 | 30-150 |
| Decachlorobiphenyl | 95 | 30-150 |

U = Below quantitation limit

Project ID: Lewis Chemical 2006-056

Lab ID: 10170

Lab Number: 10170-001

Sample ID: PZ-03-S

Matrix: Water

Sampled: 4/10/06 14:20

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 16:09 | SW3005A6010B |

Lab Number: 10170-002

Sample ID: PZ-01-D

Matrix: Water

Sampled: 4/10/06 14:45

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 16:13 | SW3005A6010B |

Lab Number: 10170-003

Sample ID: PZ-01-S

Matrix: Water

Sampled: 4/10/06 13:00

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 16:18 | SW3005A6010B |

Lab Number: 10170-004

Sample ID: PZ-02-S

Matrix: Water

Sampled: 4/10/06 15:15

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 16:23 | SW3005A6010B |

Lab Number: 10170-005

Sample ID: PZ-02-D

Matrix: Water

Sampled: 4/10/06 15:30

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/17/06 | 12:17 | SW3005A6010B |

Lab Number: 10170-006

Sample ID: PZ-03-D

Matrix: Water

Sampled: 4/10/06 15:45

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 16:27 | SW3005A6010B |

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative
Lab # 10170

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 5 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

VOC: The continuing calibration verification for naphthalene did not meet acceptance criteria (69%). Samples 10170-01, 03, and -04 were analyzed in this window. All calibration check compounds (CCC) met acceptance criteria. No further action required.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

VOC 8260: LCS 10170-51 did not meet acceptance criteria for dichlorodifluoromethane and acetone.

VOC 8260: LCS 10170-53 did not meet the acceptance limits for 2,2 dichloropropane, bromomethane, styrene and dichlorodifluoromethane. The RPD for styrene did not meet the acceptance criteria.

These compounds noted with failures are known to be problematic in the method.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

PCB: The sample peaks most closely resembled Aroclor 1242, however there was also a similarity to Aroclor 1248. Possibly due to weathering, the Aroclor identification is difficult and not absolute. Chromatograms are provided. Samples that contained non-aroclor peaks have been noted on the report pages. Quantification is quadratic.

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10170

(continued)

PCB: A sulfur clean up was performed on samples 10170-01 and -05.

VOC: The following compounds were quantified using quadratic fit:
Dichlorodifluoromethane, chloromethane, vinyl chloride, bromomethane, acetone,
dibromochloromethane, isopropylbenzene, tert-butyl benzene, and hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages.

Metals: Lead only was requested by the customer.

No other exceptions noted.

| MADEP MCP Analytical Method Report Certification Form | | | | | |
|--|--|-----------|---------------|--------------------------|-----------|
| Laboratory Name: Resource Laboratories, LLC | | | | Lab # 10170 | |
| Project Location Hyde Park | | Project # | | MADEP RTN (if available) | |
| This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers) | | | | | |
| Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other: | | | | | |
| MCP SW-846 Methods Used | 8260 (x) | 8081 () | 6010 (x) | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 (x) | EPH () | Other () | Other () | Other () |
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | | | Yes (x) No () | |
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | | | Yes (x) No () | |
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | | | Yes (x) No () | |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | | | Yes () No () NA | |
| E | Were all QC performance standards and recommendations for the specified methods achieved? | | | Yes () No (x) | |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | | | Yes () No (x) | |
| I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete. | | | | | |

Signature: 

Position: Lab Director

Printed Name: Susan C. Sylvester

Date: 4-27-06

Lab Number: 10170-50
 Sample Designation: Preparation Blank 209 W
 Date Sampled: N/A
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 74 | 30-150 |
| Decachlorobiphenyl | 86 | 30-150 |

U = Below quantitation limit

Lab Number: 10170-51
 Sample Designation: Laboratory Control Sample/Duplicate 209 W
 Date Sampled: N/A
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS

SW 846 Method 3510C/8082A.

| | Amount Added ug/L | Amount Found in LCS ug/L | LCS Recovery (%) | Amount Found in LCSD ug/L | LCSD Recovery (%) | RPD (%) | Date of Analysis |
|----------|-------------------------|-----------------------------------|------------------------|------------------------------------|-------------------------|------------|------------------|
| PCB-1016 | 2 | 1.5 | 75% | 1.5 | 75% | 0.0% | 4/13/06 |
| PCB-1242 | | | | | | | 4/13/06 |
| PCB-1221 | | | | | | | 4/13/06 |
| PCB-1232 | | | | | | | 4/13/06 |
| PCB-1248 | | | | | | | 4/13/06 |
| PCB-1254 | | | | | | | 4/13/06 |
| PCB-1260 | 2 | 1.4 | 70% | 1.5 | 75% | 6.9% | 4/13/06 |
| PCB-1262 | | | | | | | 4/13/06 |
| PCB-1268 | | | | | | | 4/13/06 |

Acceptance Criteria: 40-140% 40-140% 20%

| SURROGATE STANDARDS | Recovery (%) | Acceptance Limits (%) | Recovery (%) | Acceptance Limits (%) |
|----------------------|-----------------|--------------------------|-----------------|--------------------------|
| Tetrachloro-m-xylene | 83 | 30-150 | 78 | 30-150 |
| Decachlorobiphenyl | 88 | 30-150 | 92 | 30-150 |

U = Below quantitation limit

Lab Number 10170
Batch QC Results

METALS QC REPORT

Prep Blank

| Analyte | Sample ID # | Result (mg/L) | Reporting Limit (mg/L) |
|---------|-------------|------------------|---------------------------|
| Lead | ICB041306 | < 0.01 | 0.01 |

Laboratory Control Sample

| Analyte | Sample ID # | Result (mg/L) | True Value (mg/L) | %Recovery Control Limits (85-115%) |
|---------|-------------|------------------|----------------------|--|
| Lead | ICV041306 | 0.97 | 1.00 | 97 |

RL Resource Laboratories, LLC

Lab Number: 10170-50
Sample Designation: Method blank 041806A
Date Sampled: N/A
Date Analyzed: 4/18/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |

| SURROGATE STANDARDS | Recovery (%) | Acceptance Limits (%) |
|----------------------|--------------|-----------------------|
| dibromofluoromethane | 94 | 78-114 |
| toluene-D8 | 107 | 88-110 |
| 4-bromofluorobenzene | 98 | 86-115 |

U = Below quantitation limit

Lab Number: 10170-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: X:\DATA\VOA03\2005\APR05\041806\V3041822.D
Date Analyzed: 4/18/05
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|--------------------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 14 | 70% | 14 | 72% | 3% |
| chloromethane | 16 | 80% | 18 | 88% | 9% |
| vinyl chloride | 17 | 88% | 18 | 92% | 7% |
| bromomethane | 16 | 82% | 18 | 88% | 7% |
| chloroethane | 17 | 86% | 18 | 91% | 5% |
| trichlorofluoromethane | 18 | 90% | 19 | 97% | 7% |
| diethylether | 18 | 92% | 18 | 91% | 1% |
| acetone | 13 | 67% | 14 | 72% | 7% |
| 1,1-dichloroethene | 17 | 86% | 19 | 95% | 10% |
| methylene chloride | 19 | 97% | 19 | 97% | 0% |
| carbon disulfide | 18 | 90% | 19 | 94% | 4% |
| methyl- <i>t</i> -butyl ether (MTBE) | 35 | 89% | 36 | 90% | 1% |
| trans-1,2-dichloroethene | 18 | 91% | 20 | 98% | 7% |
| isopropyl ether (DIPE) | 18 | 92% | 19 | 97% | 5% |
| ethyl- <i>t</i> -butyl ether (ETBE) | 18 | 91% | 19 | 96% | 6% |
| 1,1-dichloroethane | 17 | 86% | 18 | 92% | 7% |
| <i>n</i> -butanol (TBA) | 74 | 74% | 78 | 78% | 5% |
| 2-butanone (MEK) | 17 | 83% | 16 | 81% | 3% |
| 2,2-dichloropropane | 17 | 83% | 18 | 89% | 7% |
| cis-1,2-dichloroethene | 20 | 101% | 21 | 107% | 7% |
| chloroform | 19 | 95% | 20 | 102% | 7% |
| bromochloromethane | 21 | 104% | 21 | 107% | 3% |
| tetrahydrofuran (THF) | 17 | 85% | 18 | 88% | 4% |
| 1,1,1-trichloroethane | 17 | 87% | 19 | 98% | 9% |
| 1,1-dichloropropene | 19 | 95% | 20 | 102% | 7% |
| <i>t</i> -amyl-methyl ether (TAME) | 19 | 94% | 19 | 96% | 3% |
| carbon tetrachloride | 17 | 83% | 18 | 90% | 8% |
| 1,2-dichloroethane | 18 | 88% | 18 | 90% | 2% |
| benzene | 20 | 100% | 21 | 106% | 5% |
| trichloroethene | 20 | 101% | 21 | 107% | 6% |
| 1,2-dichloropropane | 20 | 99% | 21 | 104% | 6% |
| bromodichloromethane | 17 | 84% | 18 | 89% | 5% |
| dibromomethane | 20 | 101% | 21 | 105% | 4% |
| 4-methyl-2-pentanone (MIBK) | 19 | 95% | 19 | 97% | 2% |
| cis-1,3-dichloropropene | 20 | 98% | 20 | 102% | 4% |
| toluene | 21 | 105% | 22 | 110% | 5% |
| trans-1,3-dichloropropene | 17 | 85% | 18 | 90% | 6% |
| 2-hexanone | 17 | 84% | 18 | 89% | 6% |
| 1,1,2-trichloroethane | 20 | 102% | 21 | 106% | 4% |
| 1,3-dichloropropane | 21 | 105% | 22 | 108% | 3% |
| tetrachloroethene | 22 | 111% | 24 | 120% | 8% |
| dibromochloromethane | 20 | 99% | 20 | 100% | 2% |
| 1,2-dibromoethane (EDB) | 21 | 106% | 21 | 106% | 1% |
| chlorobenzene | 21 | 107% | 22 | 112% | 4% |
| 1,1,1,2-tetrachloroethane | 20 | 102% | 21 | 107% | 4% |
| ethylbenzene | 21 | 106% | 22 | 108% | 2% |
| <i>m</i> & <i>p</i> -xylenes | 44 | 109% | 45 | 114% | 4% |
| <i>o</i> -xylene | 22 | 111% | 23 | 116% | 4% |
| styrene | 21 | 106% | 22 | 110% | 4% |
| bromoform | 20 | 102% | 21 | 103% | 1% |
| isopropylbenzene | 22 | 112% | 24 | 119% | 5% |
| 1,1,2,2-tetrachloroethane | 18 | 92% | 18 | 91% | 2% |
| 1,2,3-trichloropropane | 18 | 88% | 17 | 87% | 1% |
| <i>n</i> -propylbenzene | 21 | 103% | 21 | 107% | 4% |
| bromobenzene | 21 | 103% | 22 | 108% | 5% |
| 1,3,5-trimethylbenzene | 19 | 97% | 20 | 100% | 2% |
| 2-chlorotoluene | 19 | 95% | 20 | 99% | 4% |
| 4-chlorotoluene | 19 | 97% | 20 | 100% | 4% |
| tert-butylbenzene | 17 | 87% | 18 | 89% | 3% |
| 1,2,4-trimethylbenzene | 19 | 94% | 20 | 100% | 6% |
| sec-butylbenzene | 19 | 97% | 20 | 102% | 6% |
| 1,3-dichlorobenzene | 20 | 100% | 20 | 102% | 2% |
| 4-isopropyltoluene | 21 | 104% | 22 | 110% | 6% |
| 1,4-dichlorobenzene | 19 | 96% | 20 | 100% | 4% |
| 1,2-dichlorobenzene | 20 | 102% | 21 | 104% | 2% |
| <i>n</i> -butylbenzene | 20 | 100% | 21 | 107% | 7% |
| 1,2-dibromo-3-chloropropane | 17 | 83% | 17 | 87% | 5% |
| 1,2,4-trichlorobenzene | 18 | 90% | 19 | 97% | 8% |
| hexachlorobutadiene | 19 | 97% | 21 | 104% | 7% |
| naphthalene | 14 | 72% | 16 | 78% | 8% |
| 1,2,3-trichlorobenzene | 18 | 91% | 19 | 96% | 5% |
| 1,4-dioxane | 37 | 92% | 33 | 82% | 12% |
| SURROGATE STANDARDS | | | | | |
| SS dibromofluoromethane | | 92% | | 99% | |
| SS toluene-D8 | | 105% | | 108% | |
| SS 4-bromofluorobenzene | | 107% | | 101% | |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

Lab Number: 10170-52
Sample Designation: Method Blank 041906
Date Sampled: N/A
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 93 | 78-114 | | | |
| toluene-D8 | 100 | 88-110 | | | |
| 4-bromofluorobenzene | 99 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10170-53
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: X:\DATA\VOA03\2006\APR06\041806\V3041844.D
Data Analyzed: 4/19/06
SW 846 Method 5030B/8260B

| Compound | LCS | | LCSD | | RPD |
|-------------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 14 | 70% | 13 | 66% | 5% |
| chloromethane | 17 | 84% | 16 | 78% | 6% |
| vinyl chloride | 17 | 87% | 17 | 85% | 3% |
| bromomethane | 11 | 54% | 13 | 63% | 15% |
| chloroethane | 10 | 88% | 17 | 84% | 4% |
| trichlorofluoromethane | 19 | 93% | 18 | 88% | 5% |
| diethylether | 19 | 93% | 19 | 94% | 1% |
| acetone | 18 | 88% | 18 | 88% | 1% |
| 1,1-dichloroethane | 17 | 87% | 18 | 88% | 1% |
| methylene chloride | 20 | 99% | 19 | 94% | 5% |
| carbon disulfide | 18 | 88% | 17 | 84% | 5% |
| methyl-t-butyl ether (MTBE) | 38 | 95% | 37 | 93% | 3% |
| trans-1,2-dichloroethane | 18 | 91% | 18 | 90% | 2% |
| isopropyl ether (DIPE) | 19 | 95% | 18 | 90% | 5% |
| ethyl-t-butyl ether (ETBE) | 18 | 97% | 18 | 94% | 4% |
| 1,1-dichloromethane | 18 | 91% | 17 | 86% | 5% |
| t-butanol (TBA) | 90 | 90% | 88 | 88% | 3% |
| 2-butanone (MEK) | 19 | 93% | 18 | 89% | 4% |
| 2,2-dichloropropane | 13 | 67% | 13 | 66% | 2% |
| cis-1,2-dichloroethane | 21 | 104% | 20 | 99% | 5% |
| chloroform | 19 | 95% | 19 | 96% | 1% |
| bromochloromethane | 21 | 106% | 21 | 104% | 2% |
| tetrahydrofuran (THF) | 19 | 96% | 19 | 95% | 1% |
| 1,1,1-trichloroethane | 20 | 98% | 18 | 91% | 7% |
| 1,1-dichloropropene | 19 | 97% | 19 | 96% | 1% |
| t-amyl-methyl ether (TAME) | 20 | 100% | 19 | 94% | 6% |
| carbon tetrachloride | 17 | 87% | 17 | 83% | 4% |
| 1,2-dichloroethane | 18 | 88% | 18 | 89% | 2% |
| benzene | 20 | 98% | 20 | 99% | 1% |
| trichloroethane | 21 | 103% | 20 | 101% | 2% |
| 1,2-dichloropropane | 20 | 102% | 20 | 100% | 2% |
| bromodichloromethane | 18 | 88% | 17 | 87% | 1% |
| dibromomethane | 21 | 103% | 21 | 104% | 1% |
| 4-methyl-2-pentanone (MIBK) | 21 | 104% | 20 | 100% | 4% |
| cis-1,3-dichloropropene | 18 | 92% | 18 | 89% | 3% |
| toluene | 20 | 102% | 21 | 103% | 1% |
| trans-1,3-dichloropropene | 17 | 85% | 16 | 81% | 4% |
| 2-hexanone | 16 | 82% | 17 | 85% | 4% |
| 1,1,2-trichloroethane | 21 | 106% | 20 | 101% | 5% |
| 1,3-dichloropropane | 20 | 102% | 21 | 104% | 3% |
| tetrachloroethane | 22 | 108% | 22 | 111% | 3% |
| dibromochloromethane | 19 | 96% | 20 | 98% | 2% |
| 1,2-dibromoethane (EDB) | 20 | 99% | 21 | 104% | 4% |
| chlorobenzene | 20 | 98% | 20 | 102% | 4% |
| 1,1,1,2-tetrachloroethane | 20 | 99% | 20 | 100% | 2% |
| ethylbenzene | 20 | 98% | 20 | 100% | 3% |
| m,p-xylenes | 41 | 102% | 43 | 106% | 4% |
| o-xylene | 20 | 102% | 21 | 106% | 4% |
| styrene | 8 | 42% # | 14 | 72% | 53% |
| bromoform | 19 | 97% | 20 | 101% | 3% |
| isopropylbenzene | 21 | 104% | 22 | 108% | 4% |
| 1,1,2,2-tetrachloroethane | 18 | 90% | 18 | 94% | 4% |
| 1,2,3-trichloropropane | 18 | 88% | 18 | 89% | 1% |
| n-propylbenzene | 19 | 93% | 20 | 100% | 7% |
| bromobenzene | 19 | 96% | 21 | 103% | 7% |
| 1,3,5-trimethylbenzene | 17 | 87% | 18 | 97% | 10% |
| 2-chlorotoluene | 17 | 86% | 19 | 94% | 9% |
| 4-chlorotoluene | 18 | 88% | 19 | 95% | 8% |
| tert-butylbenzene | 17 | 83% | 17 | 87% | 4% |
| 1,2,4-trimethylbenzene | 18 | 90% | 19 | 97% | 7% |
| sec-butylbenzene | 18 | 91% | 18 | 92% | 0% |
| 1,3-dichlorobenzene | 19 | 94% | 20 | 99% | 5% |
| 4-isopropyltoluene | 19 | 94% | 20 | 102% | 8% |
| 1,4-dichlorobenzene | 18 | 92% | 19 | 96% | 5% |
| 1,2-dichlorobenzene | 19 | 96% | 20 | 100% | 4% |
| n-butylbenzene | 19 | 95% | 20 | 100% | 5% |
| 1,2-dibromo-3-chloropropane (| 17 | 85% | 18 | 94% | 10% |
| 1,2,4-trichlorobenzene | 17 | 86% | 19 | 94% | 8% |
| hexachlorobutadiene | 18 | 90% | 19 | 95% | 5% |
| naphthalene | 15 | 74% | 16 | 80% | 8% |
| 1,2,3-trichlorobenzene | 18 | 91% | 20 | 100% | 9% |
| 1,4-dioxane | 35 | 89% | 38 | 96% | 8% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 99% | 99% |
| SS toluene-D8 | 108% | 108% |
| SS 4-bromofluorobenzene | 107% | 101% |

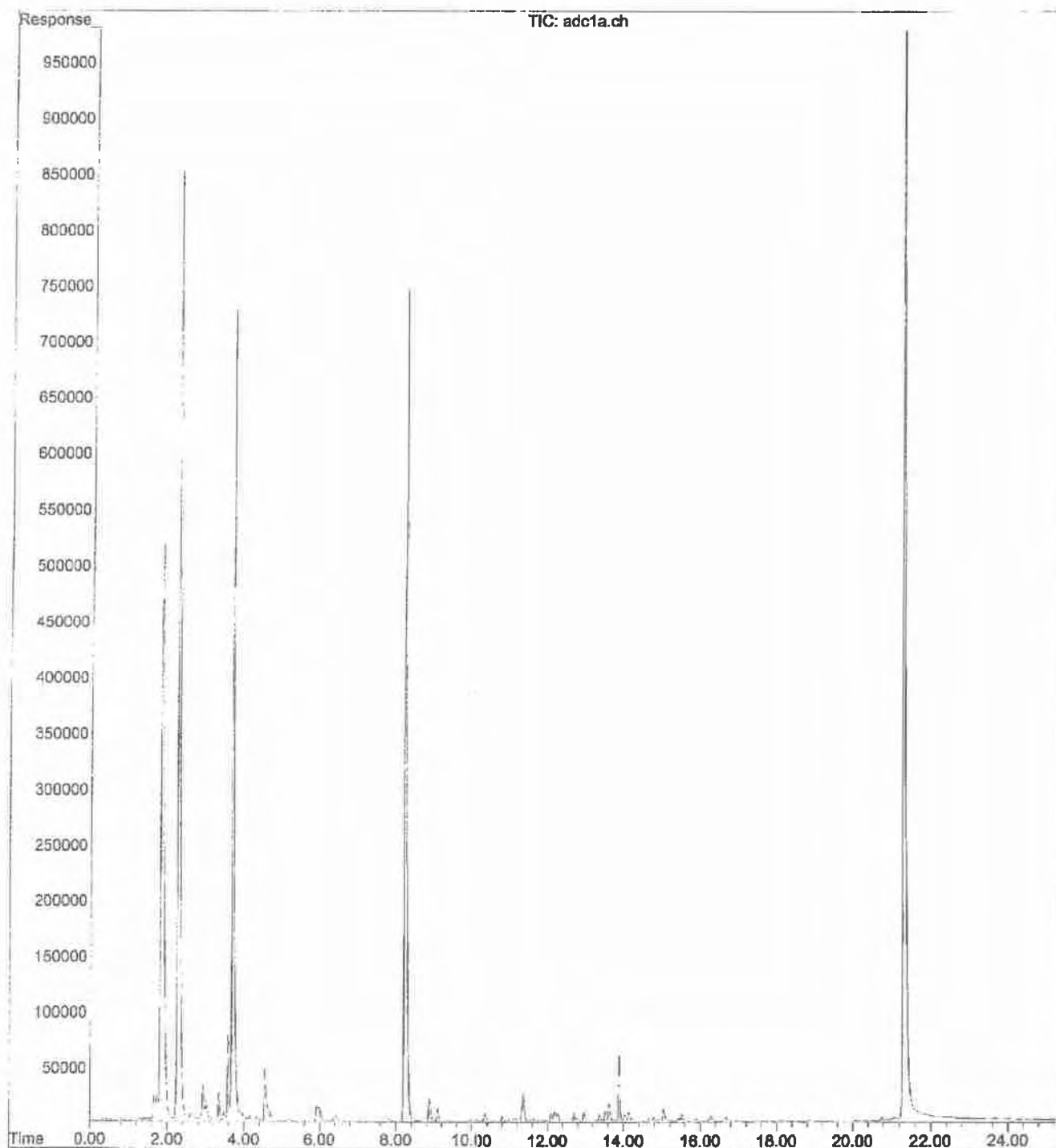
Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

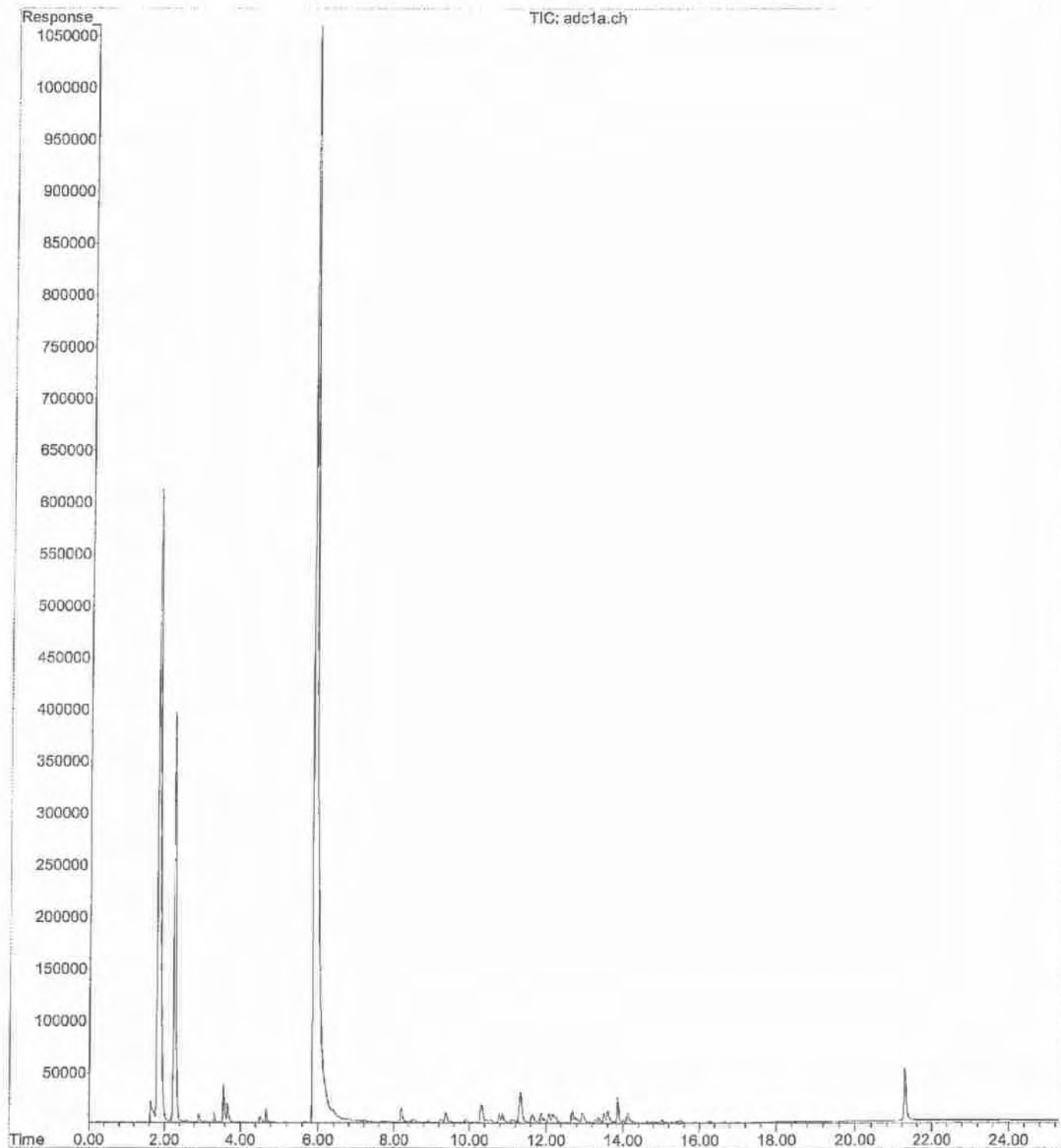
= This analyte showed recovery outside the acceptance limits.

+ = The RPD was above the acceptance limit.

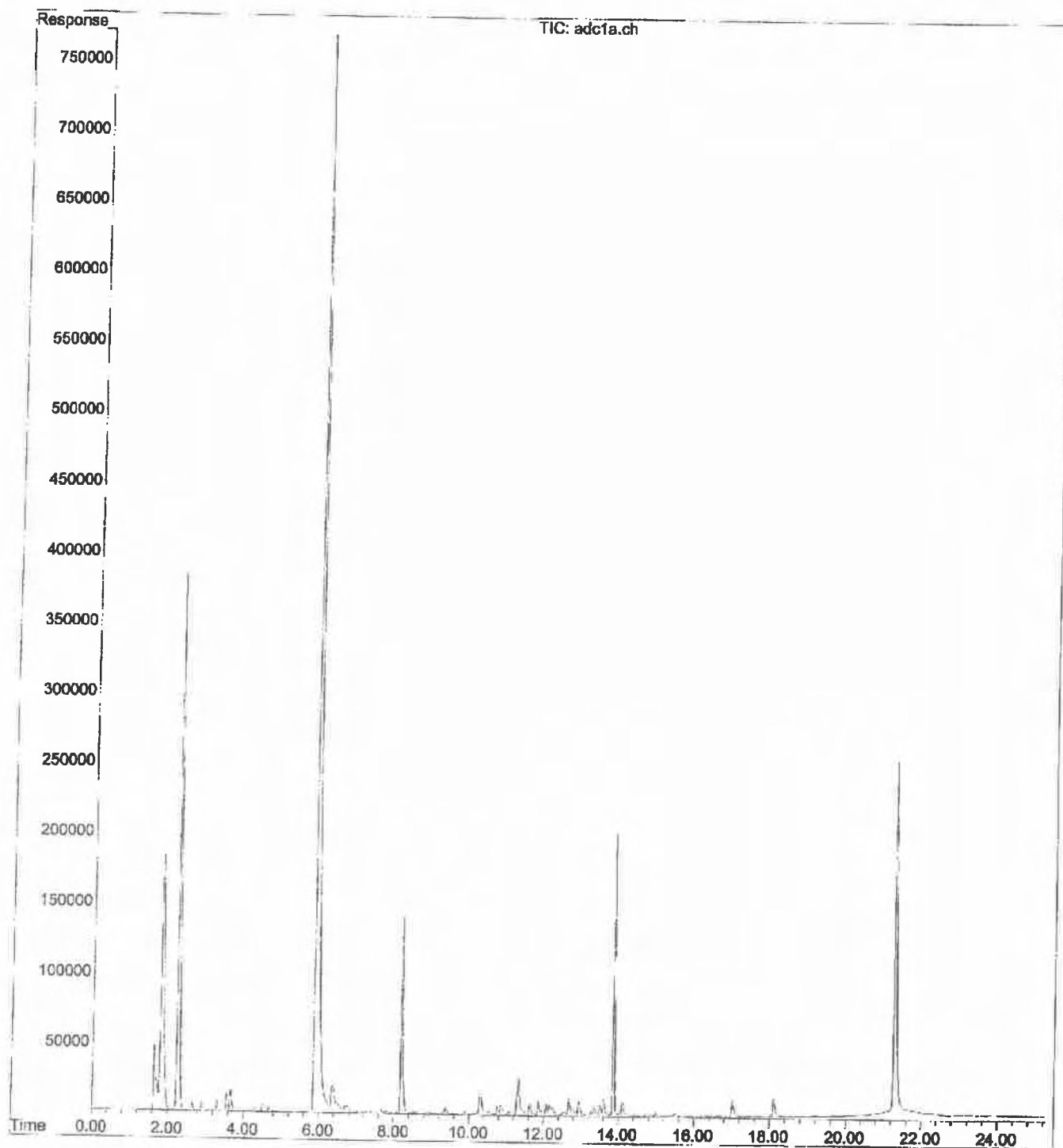
File : X:\DATA\SVOA03\2006\APR06\041906\009F0101.D
Operator : AJD
Acquired : 19 Apr 2006 05:40 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10170-01rr pcbx1
Misc Info :
Vial Number: 9



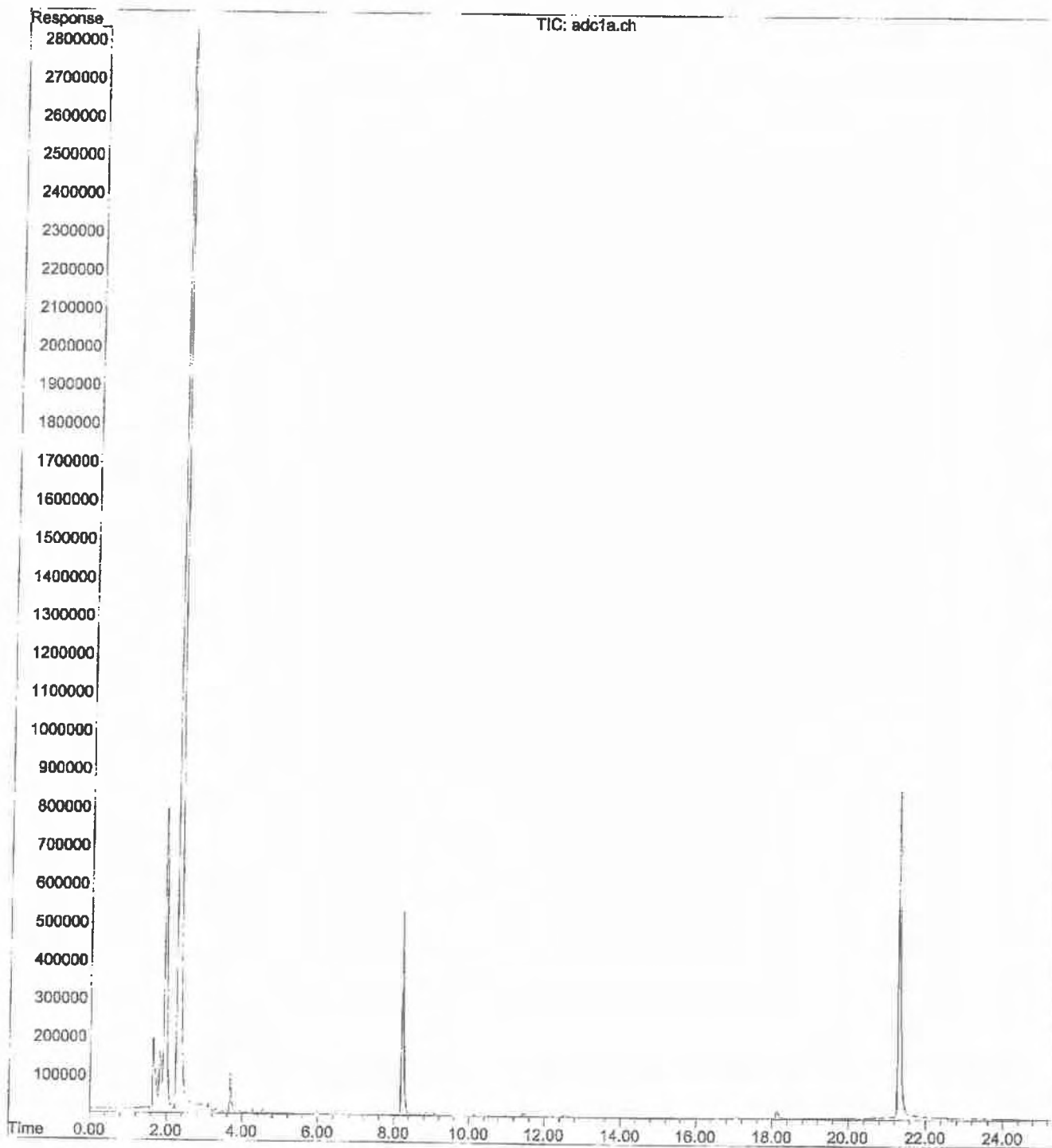
File : X:\DATA\SVOA03\2006\APR06\042006\014F0101.D
Operator : AJD
Acquired : 20 Apr 2006 04:32 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10170-08 pcb x50
Misc Info : 04 *gil*
Vial Number: 14



File : X:\DATA\SVOA03\2006\APR06\042006\013F0101.D
Operator : AJD
Acquired : 20 Apr 2006 04:04 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10170-05 pcb x5
Misc Info :
Vial Number: 13



File : X:\DATA\SVOA03\2006\APR06\041806\009F0101.D
Operator : AJD
Acquired : 18 Apr 2006 12:17 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10170-06 pcb x1
Misc Info :
Vial Number: 9



Resource Laboratories, LLC

124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

PAGE 1 OF 1
10170

Company Name: **ESM**
Phone #: **603 285 9700**
FAX #: **603 285 9957**

Company Address: **141 West Main St North**
Site Location (City, State): **Hyde Park MA**

Project Manager: **Joe Callahan**
Project ID / Name: **2016-056 / Lead Chemical**

Invoice To: **ESM**
Protocol: **RCRA SDWA NPDES NHDES OTHER**
(MCP)

| Lab Sample ID (Lab Use Only) | Field ID | # CONTAINERS | Matrix | | | Preservation Method | | | | | | Sampling | |
|---------------------------------|----------|--------------|--------|-------|-------|---------------------|------------------|--------------------------------|------|------|-----------------|----------|---------|
| | | | WATER | SOLID | OTHER | HCl | HNO ₃ | H ₂ SO ₄ | NaOH | MeOH | OTHER (Specify) | DATE | TIME |
| 10170-01 | PZ-03-S | 4 | X | | | 2 | 1 | | | | | 4/10/06 | 1420 AC |
| 02 | PZ-01-D | 3 | X | | | 2 | 1 | | | | | | 1445 LM |
| 03 | PZ-01-S | 3 | X | | | 2 | 1 | | | | | | 1300 LM |
| 04 | PZ-02-S | 4 | X | | | 2 | 1 | | | | | | 1515 09 |
| 05 | PZ-02-D | 4 | X | | | 2 | 1 | | | | | | 1530 AC |
| 06 | PZ-03-D | 4 | X | | | 2 | 1 | | | | | | 1545 AC |

| | | |
|--|---|---------------------------------------|
| <input type="checkbox"/> VOC 8260-NH List | <input type="checkbox"/> MADEP VPH | <input type="checkbox"/> MEGRO |
| <input checked="" type="checkbox"/> VOC 8260 | <input checked="" type="checkbox"/> VOC 815GRU | <input type="checkbox"/> VOC 624 |
| <input type="checkbox"/> VOC 8260 BTCX, MIBE, Naphthalene only | | |
| <input type="checkbox"/> VOC 524.2 | <input type="checkbox"/> VOC 524.2 NH List | |
| <input type="checkbox"/> TPH Fingerprint | <input type="checkbox"/> MEDRO | <input type="checkbox"/> DRO 8015 |
| <input type="checkbox"/> 8270PAH | <input type="checkbox"/> 8270ABM | <input type="checkbox"/> 625 |
| <input type="checkbox"/> 8062 PCB | <input type="checkbox"/> 8081 Pesticides | <input type="checkbox"/> 606 |
| <input type="checkbox"/> O&G 1664 | <input type="checkbox"/> O&G SM5520F | |
| <input type="checkbox"/> pH | <input type="checkbox"/> BOD | <input type="checkbox"/> Conductivity |
| <input type="checkbox"/> TSS | <input type="checkbox"/> TDS | <input type="checkbox"/> TS |
| <input type="checkbox"/> RCRA Metals | <input type="checkbox"/> Priority Pollutant Metals | <input type="checkbox"/> TAL Metals |
| <input type="checkbox"/> Total Metals-list | <input checked="" type="checkbox"/> Dissolved Metals-list | Lead |
| <input type="checkbox"/> Ammonia | <input type="checkbox"/> COD | |
| <input type="checkbox"/> T-Phosphate | <input type="checkbox"/> Phenol | |
| <input type="checkbox"/> Cyanide | <input type="checkbox"/> Sulfide | |
| <input type="checkbox"/> Nitrate | <input type="checkbox"/> Nitrite | <input type="checkbox"/> Ortho P |
| <input type="checkbox"/> Corrosivity | <input type="checkbox"/> Reactive CN | <input type="checkbox"/> Reactive S- |
| <input type="checkbox"/> TCLP Metals | <input type="checkbox"/> TCLP VOC | <input type="checkbox"/> TCLP SVOC |
| <input type="checkbox"/> TCLP Pesticide | <input type="checkbox"/> TCLP Herbicides (subcontract) | |
| <input type="checkbox"/> Standard Drinking Water Test | <input type="checkbox"/> Bacteria P/A | |

Grab (G) or Composite (C)

TEST REQUESTED

Priority (24 hr) ☐
Expedited (48 hr) ☐
10 Business Days ☒
Other: _____

REPORTING INSTRUCTIONS
☐ FAX ☐ OTHER (specify) _____
☐ PDF ☐ Excel Spreadsheet

E-Mail Address: _____
Quote #: _____
PO #: _____

RECEIVED ON ICE ☒ YES ☐ NO
TEMPERATURE **5** °C

CUSTODY RECORD

Relinquished by Sampler: **Joe Callahan**
Relinquished by: _____
Relinquished by: _____

Received by: **Joe Keo**
Received by: _____
Received by Laboratory: _____

Date: **4/10/06** Time: **11:45**
Date: _____ Time: _____
Date: _____ Time: _____

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 10171
Date Received: 4/11/06

Project: 2006-056 Lewis Chemical

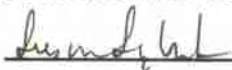
Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

4-27-06

Date

Total number of pages

50

Resource Laboratories, LLC Certifications

New Hampshire NH902
Maine NH903

Connecticut PH-0146
Massachusetts M-NH902

Lab Number: 10171-09
 Sample Designation: ESM-10
 Date Sampled: 4/10/06
 Date Received: 4/11/06
 Date Analyzed: 4/19/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Instrument Dilution Factor: 20
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | 12000 | 2000 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | 23000 | 2000 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 40 |
| benzene | C5-C8 | U | 20 |
| toluene | C5-C8 | 1900 | 40 |
| ethylbenzene | C9-C12 | 8300 | 40 |
| m&p-xylenes | C9-C12 | 12000 | 40 |
| o-xylene | C9-C12 | 1900 | 40 |
| naphthalene | N/A | U | 100 |
| C5-C8 Aliphatics (1,2) | N/A | 10000 | 2000 |
| C9-C12 Aliphatics (1,3) | N/A | U | 2000 |
| C9-C10 Aromatics (1) | N/A | U | 2000 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 101% | |
| 2,5-dibromotoluene as aliphatic | | 105% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10171-01
Sample Designation: ESM-04
Date Sampled: 4/10/06
Date Analyzed: 4/18/06
Matrix: Water
Instrument Dilution Factor: 10
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 20 | trans-1,3-dichloropropene | U | 20 |
| chloromethane | U | 20 | 2-hexanone | U | 100 |
| vinyl chloride | 25 | 20 | 1,1,2-trichloroethane | U | 20 |
| bromomethane | U | 20 | 1,3-dichloropropane | U | 20 |
| chloroethane | U | 20 | tetrachloroethene | U | 20 |
| trichlorofluoromethane | U | 20 | dibromochloromethane | U | 20 |
| diethyl ether | U | 100 | 1,2-dibromoethane | U | 20 |
| acetone | U | 100 | chlorobenzene | U | 20 |
| 1,1-dichloroethene | U | 10 | 1,1,1,2-tetrachloroethane | U | 20 |
| methylene chloride | U | 50 | ethylbenzene | U | 20 |
| carbon disulfide | U | 20 | m&p-xylenes | U | 20 |
| methyl t-butyl ether (MTBE) | U | 20 | o-xylene | U | 20 |
| trans-1,2-dichloroethene | 53 | 20 | styrene | U | 20 |
| isopropyl ether (DIPE) | U | 20 | bromoform | U | 20 |
| ethyl t-butyl ether (ETBE) | U | 20 | isopropylbenzene | U | 20 |
| 1,1-dichloroethane | 500 | 20 | 1,1,2,2-tetrachloroethane | U | 20 |
| t-butanol (TBA) | U | 500 | 1,2,3-trichloropropane | U | 20 |
| 2-butanone (MEK) | U | 100 | n-propylbenzene | U | 20 |
| 2,2-dichloropropane | U | 20 | bromobenzene | U | 20 |
| cis-1,2-dichloroethene | 2700 | 20 | 1,3,5-trimethylbenzene | U | 20 |
| chloroform | U | 20 | 2-chlorotoluene | U | 20 |
| bromochloromethane | U | 20 | 4-chlorotoluene | U | 20 |
| tetrahydrofuran (THF) | U | 100 | tert-butylbenzene | U | 20 |
| 1,1,1-trichloroethane | 900 | 20 | 1,2,4-trimethylbenzene | U | 20 |
| 1,1-dichloropropene | U | 20 | sec-butylbenzene | U | 20 |
| t-amyl-methyl ether (TAME) | U | 20 | 1,3-dichlorobenzene | U | 20 |
| carbon tetrachloride | U | 20 | 4-isopropyltoluene | U | 20 |
| 1,2-dichloroethane | U | 20 | 1,4-dichlorobenzene | U | 20 |
| benzene | U | 20 | 1,2-dichlorobenzene | U | 20 |
| trichloroethene | U | 20 | n-butylbenzene | U | 20 |
| 1,2-dichloropropane | U | 20 | 1,2-dibromo-3-chloropropane | U | 20 |
| bromodichloromethane | U | 20 | 1,2,4-trichlorobenzene | U | 20 |
| dibromomethane | U | 20 | hexachlorobutadiene | U | 20 |
| 4-methyl-2-pentanone (MIBK) | U | 100 | naphthalene | U | 50 |
| cis-1,3-dichloropropene | U | 20 | 1,2,3-trichlorobenzene | U | 20 |
| toluene | U | 20 | 1,4-dioxane | U | 500 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 96 | 78-114 | | | |
| toluene-D8 | 102 | 88-110 | | | |
| 4-bromofluorobenzene | 99 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-02
Sample Designation: ESM-09
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 50
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 100 | trans-1,3-dichloropropene | U | 100 |
| chloromethane | U | 100 | 2-hexanone | U | 500 |
| vinyl chloride | 1700 | 100 | 1,1,2-trichloroethane | U | 100 |
| bromomethane | U | 100 | 1,3-dichloropropane | U | 100 |
| chloroethane | U | 100 | tetrachloroethene | U | 100 |
| trichlorofluoromethane | U | 100 | dibromochloromethane | U | 100 |
| diethyl ether | U | 500 | 1,2-dibromoethane | U | 100 |
| acetone | U | 500 | chlorobenzene | U | 100 |
| 1,1-dichloroethene | 470 | 50 | 1,1,1,2-tetrachloroethane | U | 100 |
| methylene chloride | U | 300 | ethylbenzene | 140 | 100 |
| carbon disulfide | U | 100 | m&p-xylenes | 270 | 100 |
| methyl t-butyl ether (MTBE) | U | 100 | o-xylene | U | 100 |
| trans-1,2-dichloroethene | 240 | 100 | styrene | U | 100 |
| isopropyl ether (DIPE) | U | 100 | bromoform | U | 100 |
| ethyl t-butyl ether (ETBE) | U | 100 | isopropylbenzene | U | 100 |
| 1,1-dichloroethane | 3500 | 100 | 1,1,2,2-tetrachloroethane | U | 100 |
| t-butanol (TBA) | U | 3000 | 1,2,3-trichloropropane | U | 100 |
| 2-butanone (MEK) | U | 500 | n-propylbenzene | U | 100 |
| 2,2-dichloropropane | U | 100 | bromobenzene | U | 100 |
| cis-1,2-dichloroethene | 26000 | 100 | 1,3,5-trimethylbenzene | U | 100 |
| chloroform | U | 100 | 2-chlorotoluene | U | 100 |
| bromochloromethane | U | 100 | 4-chlorotoluene | U | 100 |
| tetrahydrofuran (THF) | U | 500 | tert-butylbenzene | U | 100 |
| 1,1,1-trichloroethane | 35000 | 100 | 1,2,4-trimethylbenzene | U | 100 |
| 1,1-dichloropropene | U | 100 | sec-butylbenzene | U | 100 |
| t-amyl-methyl ether (TAME) | U | 100 | 1,3-dichlorobenzene | U | 100 |
| carbon tetrachloride | U | 100 | 4-isopropyltoluene | U | 100 |
| 1,2-dichloroethane | 290 | 100 | 1,4-dichlorobenzene | U | 100 |
| benzene | U | 100 | 1,2-dichlorobenzene | U | 100 |
| trichloroethene | U | 100 | n-butylbenzene | U | 100 |
| 1,2-dichloropropane | U | 100 | 1,2-dibromo-3-chloropropane | U | 100 |
| bromodichloromethane | U | 100 | 1,2,4-trichlorobenzene | U | 100 |
| dibromomethane | U | 100 | hexachlorobutadiene | U | 100 |
| 4-methyl-2-pentanone (MIBK) | U | 500 | naphthalene | U | 300 |
| cis-1,3-dichloropropene | U | 100 | 1,2,3-trichlorobenzene | U | 100 |
| toluene | 5900 | 100 | 1,4-dioxane | U | 3000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 94 | 78-114 | | | |
| toluene-D8 | 103 | 88-110 | | | |
| 4-bromofluorobenzene | 100 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-03
Sample Designation: DUP-2
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 50
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 100 | trans-1,3-dichloropropene | U | 100 |
| chloromethane | U | 100 | 2-hexanone | U | 500 |
| vinyl chloride | 1600 | 100 | 1,1,2-trichloroethane | U | 100 |
| bromomethane | U | 100 | 1,3-dichloropropane | U | 100 |
| chloroethane | U | 100 | tetrachloroethene | U | 100 |
| trichlorofluoromethane | U | 100 | dibromochloromethane | U | 100 |
| diethyl ether | U | 500 | 1,2-dibromoethane | U | 100 |
| acetone | U | 500 | chlorobenzene | U | 100 |
| 1,1-dichloroethene | 480 | 50 | 1,1,1,2-tetrachloroethane | U | 100 |
| methylene chloride | U | 300 | ethylbenzene | 140 | 100 |
| carbon disulfide | U | 100 | m&p-xylenes | 270 | 100 |
| methyl t-butyl ether (MTBE) | U | 100 | o-xylene | U | 100 |
| trans-1,2-dichloroethene | 230 | 100 | styrene | U | 100 |
| isopropyl ether (DIPE) | U | 100 | bromoform | U | 100 |
| ethyl t-butyl ether (ETBE) | U | 100 | isopropylbenzene | U | 100 |
| 1,1-dichloroethane | 3400 | 100 | 1,1,2,2-tetrachloroethane | U | 100 |
| t-butanol (TBA) | U | 3000 | 1,2,3-trichloropropane | U | 100 |
| 2-butanone (MEK) | U | 500 | n-propylbenzene | U | 100 |
| 2,2-dichloropropane | U | 100 | bromobenzene | U | 100 |
| cis-1,2-dichloroethene | 25000 | 100 | 1,3,5-trimethylbenzene | U | 100 |
| chloroform | U | 100 | 2-chlorotoluene | U | 100 |
| bromochloromethane | U | 100 | 4-chlorotoluene | U | 100 |
| tetrahydrofuran (THF) | U | 500 | tert-butylbenzene | U | 100 |
| 1,1,1-trichloroethane | 34000 | 100 | 1,2,4-trimethylbenzene | U | 100 |
| 1,1-dichloropropene | U | 100 | sec-butylbenzene | U | 100 |
| t-amyl-methyl ether (TAME) | U | 100 | 1,3-dichlorobenzene | U | 100 |
| carbon tetrachloride | U | 100 | 4-isopropyltoluene | U | 100 |
| 1,2-dichloroethane | 290 | 100 | 1,4-dichlorobenzene | U | 100 |
| benzene | U | 100 | 1,2-dichlorobenzene | U | 100 |
| trichloroethene | U | 100 | n-butylbenzene | U | 100 |
| 1,2-dichloropropane | U | 100 | 1,2-dibromo-3-chloropropane | U | 100 |
| bromodichloromethane | U | 100 | 1,2,4-trichlorobenzene | U | 100 |
| dibromomethane | U | 100 | hexachlorobutadiene | U | 100 |
| 4-methyl-2-pentanone (MIBK) | U | 500 | naphthalene | U | 300 |
| cis-1,3-dichloropropene | U | 100 | 1,2,3-trichlorobenzene | U | 100 |
| toluene | 5800 | 100 | 1,4-dioxane | U | 3000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 95 | 78-114 | | | |
| toluene-D8 | 105 | 88-110 | | | |
| 4-bromofluorobenzene | 101 | 86-115 | | | |

U = Below quantitation limit

RL Resource Laboratories, LLC

Lab Number: 10171-04
Sample Designation: ESM-05
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 200
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 400 | trans-1,3-dichloropropene | U | 400 |
| chloromethane | U | 400 | 2-hexanone | U | 2000 |
| vinyl chloride | 490 | 400 | 1,1,2-trichloroethane | U | 400 |
| bromomethane | U | 400 | 1,3-dichloropropane | U | 400 |
| chloroethane | U | 400 | tetrachloroethene | 21000 | 400 |
| trichlorofluoromethane | U | 400 | dibromochloromethane | U | 400 |
| diethyl ether | U | 2000 | 1,2-dibromoethane | U | 400 |
| acetone | U | 2000 | chlorobenzene | 2300 | 400 |
| 1,1-dichloroethene | 2800 | 200 | 1,1,1,2-tetrachloroethane | U | 400 |
| methylene chloride | 1300 | 1000 | ethylbenzene | U | 400 |
| carbon disulfide | U | 400 | m&p-xylenes | 610 | 400 |
| methyl t-butyl ether (MTBE) | U | 400 | o-xylene | U | 400 |
| trans-1,2-dichloroethene | 460 | 400 | styrene | U | 400 |
| isopropyl ether (DIPE) | U | 400 | bromoform | U | 400 |
| ethyl t-butyl ether (ETBE) | U | 400 | isopropylbenzene | U | 400 |
| 1,1-dichloroethane | 4900 | 400 | 1,1,2,2-tetrachloroethane | U | 400 |
| t-butanol (TBA) | U | 10000 | 1,2,3-trichloropropane | U | 400 |
| 2-butanone (MEK) | U | 2000 | n-propylbenzene | U | 400 |
| 2,2-dichloropropane | U | 400 | bromobenzene | U | 400 |
| cis-1,2-dichloroethene | 110000 | 400 | 1,3,5-trimethylbenzene | U | 400 |
| chloroform | U | 400 | 2-chlorotoluene | U | 400 |
| bromochloromethane | U | 400 | 4-chlorotoluene | U | 400 |
| tetrahydrofuran (THF) | U | 2000 | tert-butylbenzene | U | 400 |
| 1,1,1-trichloroethane | 280000 D | 4000 | 1,2,4-trimethylbenzene | U | 400 |
| 1,1-dichloropropene | U | 400 | sec-butylbenzene | U | 400 |
| t-amyl-methyl ether (TAME) | U | 400 | 1,3-dichlorobenzene | U | 400 |
| carbon tetrachloride | U | 400 | 4-isopropyltoluene | U | 400 |
| 1,2-dichloroethane | 1300 | 400 | 1,4-dichlorobenzene | U | 400 |
| benzene | U | 400 | 1,2-dichlorobenzene | U | 400 |
| trichloroethene | 360000 D | 4000 | n-butylbenzene | U | 400 |
| 1,2-dichloropropane | U | 400 | 1,2-dibromo-3-chloropropane | U | 400 |
| bromodichloromethane | U | 400 | 1,2,4-trichlorobenzene | U | 400 |
| dibromomethane | U | 400 | hexachlorobutadiene | U | 400 |
| 4-methyl-2-pentanone (MIBK) | U | 2000 | naphthalene | U | 1000 |
| cis-1,3-dichloropropene | U | 400 | 1,2,3-trichlorobenzene | U | 400 |
| toluene | 38000 | 400 | 1,4-dioxane | U | 10000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 101 | 78-114 | | | |
| toluene-D8 | 107 | 88-110 | | | |
| 4-bromofluorobenzene | 100 | 86-115 | | | |

D = Result provided from a re-analysis at a dilution.
U = Below quantitation limit

Lab Number: 10171-05
Sample Designation: ESM-05B
Date Sampled: 4/10/06
Date Analyzed: 4/18/06
Matrix: Water
Instrument Dilution Factor: 10
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 20 | trans-1,3-dichloropropene | U | 20 |
| chloromethane | U | 20 | 2-hexanone | U | 100 |
| vinyl chloride | U | 20 | 1,1,2-trichloroethane | U | 20 |
| bromomethane | U | 20 | 1,3-dichloropropane | U | 20 |
| chloroethane | U | 20 | tetrachloroethene | 2700 | 20 |
| trichlorofluoromethane | U | 20 | dibromochloromethane | U | 20 |
| diethyl ether | U | 100 | 1,2-dibromoethane | U | 20 |
| acetone | U | 100 | chlorobenzene | U | 20 |
| 1,1-dichloroethene | 320 | 10 | 1,1,1,2-tetrachloroethane | U | 20 |
| methylene chloride | 1200 | 50 | ethylbenzene | 32 | 20 |
| carbon disulfide | U | 20 | m&p-xylenes | U | 20 |
| methyl t-butyl ether (MTBE) | U | 20 | o-xylene | 30 | 20 |
| trans-1,2-dichloroethene | U | 20 | styrene | U | 20 |
| isopropyl ether (DIPE) | U | 20 | bromoform | U | 20 |
| ethyl t-butyl ether (ETBE) | U | 20 | isopropylbenzene | U | 20 |
| 1,1-dichloroethane | 180 | 20 | 1,1,2,2-tetrachloroethane | U | 20 |
| t-butanol (TBA) | U | 500 | 1,2,3-trichloropropane | U | 20 |
| 2-butanone (MEK) | U | 100 | n-propylbenzene | U | 20 |
| 2,2-dichloropropane | U | 20 | bromobenzene | U | 20 |
| cis-1,2-dichloroethene | 62 | 20 | 1,3,5-trimethylbenzene | U | 20 |
| chloroform | U | 20 | 2-chlorotoluene | U | 20 |
| bromochloromethane | U | 20 | 4-chlorotoluene | U | 20 |
| tetrahydrofuran (THF) | U | 100 | tert-butylbenzene | U | 20 |
| 1,1,1-trichloroethane | 6300 | 20 | 1,2,4-trimethylbenzene | U | 20 |
| 1,1-dichloropropene | U | 20 | sec-butylbenzene | U | 20 |
| t-amyl-methyl ether (TAME) | U | 20 | 1,3-dichlorobenzene | U | 20 |
| carbon tetrachloride | U | 20 | 4-isopropyltoluene | U | 20 |
| 1,2-dichloroethane | 120 | 20 | 1,4-dichlorobenzene | U | 20 |
| benzene | U | 20 | 1,2-dichlorobenzene | 93 | 20 |
| trichloroethene | 7100 | 20 | n-butylbenzene | U | 20 |
| 1,2-dichloropropane | U | 20 | 1,2-dibromo-3-chloropropane | U | 20 |
| bromodichloromethane | U | 20 | 1,2,4-trichlorobenzene | U | 20 |
| dibromomethane | U | 20 | hexachlorobutadiene | U | 20 |
| 4-methyl-2-pentanone (MIBK) | U | 100 | naphthalene | U | 50 |
| cis-1,3-dichloropropene | U | 20 | 1,2,3-trichlorobenzene | U | 20 |
| toluene | 430 | 20 | 1,4-dioxane | U | 500 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 94 | 78-114 | | | |
| toluene-D8 | 102 | 88-110 | | | |
| 4-bromofluorobenzene | 102 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-06
Sample Designation: ESM-06
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 20
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 40 | trans-1,3-dichloropropene | U | 40 |
| chloromethane | U | 40 | 2-hexanone | U | 200 |
| vinyl chloride | U | 40 | 1,1,2-trichloroethane | U | 40 |
| bromomethane | U | 40 | 1,3-dichloropropane | U | 40 |
| chloroethane | U | 40 | tetrachloroethene | 1600 | 40 |
| trichlorofluoromethane | U | 40 | dibromochloromethane | U | 40 |
| diethyl ether | U | 200 | 1,2-dibromoethane | U | 40 |
| acetone | U | 200 | chlorobenzene | U | 40 |
| 1,1-dichloroethene | 140 | 20 | 1,1,1,2-tetrachloroethane | U | 40 |
| methylene chloride | U | 100 | ethylbenzene | U | 40 |
| carbon disulfide | U | 40 | m&p-xylenes | U | 40 |
| methyl t-butyl ether (MTBE) | U | 40 | o-xylene | U | 40 |
| trans-1,2-dichloroethene | 43 | 40 | styrene | U | 40 |
| isopropyl ether (DIPE) | U | 40 | bromoform | U | 40 |
| ethyl t-butyl ether (ETBE) | U | 40 | isopropylbenzene | U | 40 |
| 1,1-dichloroethane | 280 | 40 | 1,1,2,2-tetrachloroethane | U | 40 |
| t-butanol (TBA) | U | 1000 | 1,2,3-trichloropropane | U | 40 |
| 2-butanone (MEK) | U | 200 | n-propylbenzene | U | 40 |
| 2,2-dichloropropane | U | 40 | bromobenzene | U | 40 |
| cis-1,2-dichloroethene | 8800 | 40 | 1,3,5-trimethylbenzene | U | 40 |
| chloroform | U | 40 | 2-chlorotoluene | U | 40 |
| bromochloromethane | U | 40 | 4-chlorotoluene | U | 40 |
| tetrahydrofuran (THF) | U | 200 | tert-butylbenzene | U | 40 |
| 1,1,1-trichloroethane | 11000 | 40 | 1,2,4-trimethylbenzene | U | 40 |
| 1,1-dichloropropene | U | 40 | sec-butylbenzene | U | 40 |
| t-amyl-methyl ether (TAME) | U | 40 | 1,3-dichlorobenzene | U | 40 |
| carbon tetrachloride | U | 40 | 4-isopropyltoluene | U | 40 |
| 1,2-dichloroethane | U | 40 | 1,4-dichlorobenzene | U | 40 |
| benzene | U | 40 | 1,2-dichlorobenzene | U | 40 |
| trichloroethene | 11000 | 40 | n-butylbenzene | U | 40 |
| 1,2-dichloropropane | U | 40 | 1,2-dibromo-3-chloropropane | U | 40 |
| bromodichloromethane | U | 40 | 1,2,4-trichlorobenzene | U | 40 |
| dibromomethane | U | 40 | hexachlorobutadiene | U | 40 |
| 4-methyl-2-pentanone (MIBK) | U | 200 | naphthalene | U | 100 |
| cis-1,3-dichloropropene | U | 40 | 1,2,3-trichlorobenzene | U | 40 |
| toluene | 110 | 40 | 1,4-dioxane | U | 1000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 95 | 78-114 | | | |
| toluene-D8 | 102 | 88-110 | | | |
| 4-bromofluorobenzene | 101 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-07
Sample Designation: ESM-15
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 50
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 100 | trans-1,3-dichloropropene | U | 100 |
| chloromethane | U | 100 | 2-hexanone | U | 500 |
| vinyl chloride | 920 | 100 | 1,1,2-trichloroethane | 130 | 100 |
| bromomethane | U | 100 | 1,3-dichloropropane | U | 100 |
| chloroethane | U | 100 | tetrachloroethene | 11000 | 100 |
| trichlorofluoromethane | U | 100 | dibromochloromethane | U | 100 |
| diethyl ether | U | 500 | 1,2-dibromoethane | U | 100 |
| acetone | U | 500 | chlorobenzene | U | 100 |
| 1,1-dichloroethene | 920 | 50 | 1,1,1,2-tetrachloroethane | U | 100 |
| methylene chloride | 2100 | 300 | ethylbenzene | 460 | 100 |
| carbon disulfide | U | 100 | m&p-xylenes | 1300 | 100 |
| methyl t-butyl ether (MTBE) | U | 100 | o-xylene | 450 | 100 |
| trans-1,2-dichloroethene | 210 | 100 | styrene | U | 100 |
| isopropyl ether (DIPE) | U | 100 | bromoform | U | 100 |
| ethyl t-butyl ether (ETBE) | U | 100 | isopropylbenzene | U | 100 |
| 1,1-dichloroethane | 2600 | 100 | 1,1,1,2-tetrachloroethane | U | 100 |
| t-butanol (TBA) | U | 3000 | 1,2,3-trichloropropane | U | 100 |
| 2-butanone (MEK) | U | 500 | n-propylbenzene | U | 100 |
| 2,2-dichloropropane | U | 100 | bromobenzene | U | 100 |
| cis-1,2-dichloroethene | 27000 | 100 | 1,3,5-trimethylbenzene | U | 100 |
| chloroform | U | 100 | 2-chlorotoluene | U | 100 |
| bromochloromethane | U | 100 | 4-chlorotoluene | U | 100 |
| tetrahydrofuran (THF) | U | 500 | tert-butylbenzene | U | 100 |
| 1,1,1-trichloroethane | 66000 D | 1000 | 1,2,4-trimethylbenzene | U | 100 |
| 1,1-dichloropropene | U | 100 | sec-butylbenzene | U | 100 |
| t-amyl-methyl ether (TAME) | U | 100 | 1,3-dichlorobenzene | U | 100 |
| carbon tetrachloride | U | 100 | 4-isopropyltoluene | U | 100 |
| 1,2-dichloroethane | 2900 | 100 | 1,4-dichlorobenzene | U | 100 |
| benzene | U | 100 | 1,2-dichlorobenzene | U | 100 |
| trichloroethene | 93000 D | 1000 | n-butylbenzene | U | 100 |
| 1,2-dichloropropane | U | 100 | 1,2-dibromo-3-chloropropane | U | 100 |
| bromodichloromethane | U | 100 | 1,2,4-trichlorobenzene | U | 100 |
| dibromomethane | U | 100 | hexachlorobutadiene | U | 100 |
| 4-methyl-2-pentanone (MIBK) | U | 500 | naphthalene | U | 300 |
| cis-1,3-dichloropropene | U | 100 | 1,2,3-trichlorobenzene | U | 100 |
| toluene | 9900 | 100 | 1,4-dioxane | U | 3000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 98 | 78-114 | | | |
| toluene-D8 | 103 | 88-110 | | | |
| 4-bromofluorobenzene | 107 | 86-115 | | | |

D = Result provided from a re-analysis at a dilution.
U = Below quantitation limit

Lab Number: 10171-08
Sample Designation: ESM-07
Date Sampled: 4/10/06
Date Analyzed: 4/18/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | 16 | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | 37 | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | 3 | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | 58 | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | 110 | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | 63 | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | 90 | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dlbromofluoromethane | 97 | 78-114 | | | |
| toluene-D8 | 106 | 88-110 | | | |
| 4-bromofluorobenzene | 98 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-09
Sample Designation: ESM-10
Date Sampled: 4/10/06
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 20
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 40 | trans-1,3-dichloropropene | U | 40 |
| chloromethane | U | 40 | 2-hexanone | U | 200 |
| vinyl chloride | 120 | 40 | 1,1,2-trichloroethane | U | 40 |
| bromomethane | U | 40 | 1,3-dichloropropane | U | 40 |
| chloroethane | U | 40 | tetrachloroethene | 180 | 40 |
| trichlorofluoromethane | U | 40 | dibromochloromethane | U | 40 |
| diethyl ether | U | 200 | 1,2-dibromoethane | U | 40 |
| acetone | U | 200 | chlorobenzene | U | 40 |
| 1,1-dichloroethene | U | 20 | 1,1,1,2-tetrachloroethane | U | 40 |
| methylene chloride | 410 | 100 | ethylbenzene | 11000 | 40 |
| carbon disulfide | U | 40 | m&p-xylenes | 17000 | 40 |
| methyl t-butyl ether (MTBE) | U | 40 | o-xylene | 2300 | 40 |
| trans-1,2-dichloroethene | U | 40 | styrene | U | 40 |
| isopropyl ether (DIPE) | U | 40 | bromoform | U | 40 |
| ethyl t-butyl ether (ETBE) | U | 40 | isopropylbenzene | U | 40 |
| 1,1-dichloroethane | 91 | 40 | 1,1,2,2-tetrachloroethane | U | 40 |
| t-butanol (TBA) | U | 1000 | 1,2,3-trichloropropane | U | 40 |
| 2-butanone (MEK) | U | 200 | n-propylbenzene | U | 40 |
| 2,2-dichloropropane | U | 40 | bromobenzene | U | 40 |
| cis-1,2-dichloroethene | 610 | 40 | 1,3,5-trimethylbenzene | U | 40 |
| chloroform | U | 40 | 2-chlorotoluene | U | 40 |
| bromochloromethane | U | 40 | 4-chlorotoluene | U | 40 |
| tetrahydrofuran (THF) | U | 200 | tert-butylbenzene | U | 40 |
| 1,1,1-trichloroethane | U | 40 | 1,2,4-trimethylbenzene | U | 40 |
| 1,1-dichloropropene | U | 40 | sec-butylbenzene | U | 40 |
| t-amyl-methyl ether (TAME) | U | 40 | 1,3-dichlorobenzene | U | 40 |
| carbon tetrachloride | U | 40 | 4-isopropyltoluene | 620 | 40 |
| 1,2-dichloroethane | U | 40 | 1,4-dichlorobenzene | U | 40 |
| benzene | U | 40 | 1,2-dichlorobenzene | U | 40 |
| trichloroethene | 120 | 40 | n-butylbenzene | U | 40 |
| 1,2-dichloropropane | U | 40 | 1,2-dibromo-3-chloropropane | U | 40 |
| bromodichloromethane | U | 40 | 1,2,4-trichlorobenzene | U | 40 |
| dibromomethane | U | 40 | hexachlorobutadiene | U | 40 |
| 4-methyl-2-pentanone (MIBK) | U | 200 | naphthalene | U | 100 |
| cis-1,3-dichloropropene | U | 40 | 1,2,3-trichlorobenzene | U | 40 |
| toluene | 2100 | 40 | 1,4-dioxane | U | 1000 |
| SURROGATE STANDARDS | | | | | |
| | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 94 | 78-114 | | | |
| toluene-D8 | 105 | 88-110 | | | |
| 4-bromofluorobenzene | 106 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-10
Sample Designation: Trip Blank
Date Sampled: 4/10/06
Date Analyzed: 4/17/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 92 | 78-114 | | | |
| toluene-D8 | 104 | 88-110 | | | |
| 4-bromofluorobenzene | 102 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-09
 Sample Designation: ESM-10
 Date Sampled: 4/10/06
 Date Received: 4/11/06
 Date Extracted: 4/19/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: pH ≤ 2
 Temperature: Received on Ice at 4±2°C
 Dilution Factor: 1
 Analyst: AJD

EPH ANALYTICAL RESULTS

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | Concentration ug/L | Quantitation Limit ug/L | Date of Analysis |
|-------------------------------------|-----------------------|----------------------------|------------------|
| Diesel PAH Analytes: | | | |
| naphthalene | 3.1 | 0.5 | 4/20/06 |
| 2-methylnaphthalene | U | 0.5 | 4/20/06 |
| phenanthrene | U | 0.5 | 4/20/06 |
| acenaphthene | U | 0.5 | 4/20/06 |
| Other PAH Analytes: | | | |
| acenaphthylene | U | 0.5 | 4/20/06 |
| fluorene | U | 0.5 | 4/20/06 |
| anthracene | U | 0.5 | 4/20/06 |
| fluoranthene | U | 0.5 | 4/20/06 |
| pyrene | U | 0.5 | 4/20/06 |
| benzo(a)anthracene | U | 0.5 | 4/20/06 |
| chrysene | U | 0.5 | 4/20/06 |
| benzo(b)fluoranthene | U | 0.5 | 4/20/06 |
| benzo(k)fluoranthene | U | 0.5 | 4/20/06 |
| benzo(a)pyrene | U | 0.2 | 4/20/06 |
| indeno(1,2,3-cd)pyrene | U | 0.5 | 4/20/06 |
| dibenzo(a,h)anthracene | U | 0.5 | 4/20/06 |
| benzo(g,h,i)perylene | U | 0.5 | 4/20/06 |
| Ranges: | | | |
| Unadjusted C11-C22 Aromatics (1) | U | 200 | 4/20/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | U | 200 | 4/20/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | U | 200 | 4/20/06 |
| Extraction Surrogate Recoveries: | | | |
| 1-chloro-octadecane | 37% # | | |
| o-terphenyl | 49% | | |
| Fractionation Surrogate Recoveries: | | | |
| 2-fluorobiphenyl | 82% | | |
| 2-bromonaphthalene | 79% | | |
| Surrogate Acceptance Range | 40-140% | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

= Surrogate below acceptance criteria. Matrix Interference suspected.

RL Resource Laboratories, LLC

Lab Number: 10171-01
 Sample Designation: ESM-04
 Date Sampled: 4/10/06
 Date Extracted: 4/13/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 44 | 30-150 |
| Decachlorobiphenyl | 64 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-02
 Sample Designation: ESM-09
 Date Sampled: 4/10/06
 Date Extracted: 4/14/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 78 | 30-150 |
| Decachlorobiphenyl | 91 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-04
 Sample Designation: ESM-05
 Date Sampled: 4/10/06
 Date Extracted: 4/14/06
 Date Analyzed: 4/19/06
 Matrix: Water
 Dilution Factor: 20
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 6 |
| PCB-1242 | 26 | 6 |
| PCB-1221 | U | 6 |
| PCB-1232 | U | 6 |
| PCB-1248 | U | 6 |
| PCB-1254 | U | 6 |
| PCB-1260 | U | 6 |
| PCB-1262 | U | 6 |
| PCB-1268 | U | 6 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 36 | 30-150 |
| Decachlorobiphenyl | 132 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-05
 Sample Designation: ESM-05B
 Date Sampled: 4/10/06
 Date Extracted: 4/14/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 68 | 30-150 |
| Decachlorobiphenyl | 96 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-08
 Sample Designation: ESM-08
 Date Sampled: 4/10/06
 Date Extracted: 4/14/06
 Date Analyzed: 4/19/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | 3.5 | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 80 | 30-150 |
| Decachlorobiphenyl | 94 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-07
 Sample Designation: ESM-15
 Date Sampled: 4/10/06
 Date Extracted: 4/14/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 69 | 30-150 |
| Decachlorobiphenyl | 94 | 30-150 |

Note: Non-aroclor like peaks were observed in this sample.
 U = Below quantitation limit

Lab Number: 10171-08
 Sample Designation: ESM-07
 Date Sampled: 4/10/06
 Date Extracted: 4/14/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 83 | 30-150 |
| Decachlorobiphenyl | 96 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-09
 Sample Designation: ESM-10
 Date Sampled: 4/10/06
 Date Extracted: 4/14/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 61 | 30-150 |
| Decachlorobiphenyl | 78 | 30-150 |

Note: A non-aroclor like compound was observed in this sample.
 U = Below quantitation limit

Project ID: DND Lewis 2006-056

Lab ID: 10171

Lab Number: 10171-004

Sample ID: ESM-05

Matrix: Water

Sampled: 4/10/06 9:30

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|----------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Arsenic | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:24 | SW3005A6010B |
| Barium | 0.18 | 0.05 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:24 | SW3005A6010B |
| Cadmium | < 0.005 | 0.005 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:24 | SW3005A6010B |
| Chromium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:24 | SW3005A6010B |
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:24 | SW3005A6010B |
| Mercury | < 0.0009 | 0.0009 | mg/L | 1 | BJS | 4/12/06 | 4/12/06 | N/A | SW7470A |
| Selenium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:24 | SW3005A6010B |
| Silver | < 0.007 | 0.007 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:24 | SW3005A6010B |

Lab Number: 10171-005

Sample ID: ESM-05B

Matrix: Water

Sampled: 4/10/06 10:00

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|------------|----------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| Arsenic | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:40 | SW3005A6010B |
| Barium | 0.22 | 0.05 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:40 | SW3005A6010B |
| Cadmium | < 0.005 | 0.005 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:40 | SW3005A6010B |
| Chromium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:40 | SW3005A6010B |
| Lead | < 0.01 | 0.01 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:40 | SW3005A6010B |
| Mercury | < 0.0009 | 0.0009 | mg/L | 1 | BJS | 4/12/06 | 4/12/06 | N/A | SW7470A |
| Selenium | < 0.05 | 0.05 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:40 | SW3005A6010B |
| Silver | < 0.007 | 0.007 | mg/L | 1 | BJS | N/A | 4/13/06 | 15:40 | SW3005A6010B |

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 10171

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 4 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

VOC: The continuing calibration verification for naphthalene did not meet acceptance criteria (69%). Samples 10171-02,03,04,06,07, and 09 were analyzed in this window. All calibration check compounds (CCC) met acceptance criteria. No further action required.

Method Blank

No exceptions noted.

Surrogate Recoveries

EPH: The following sample had a surrogate that did not meet the acceptance criteria: 10171-09. All other batch QC had acceptable recoveries. It is suspected that this may have been caused by matrix interference. The samples could not be re-extracted due to insufficient sample volume.

Laboratory Control Sample Results

VOC 8260: LCS 10171-51 did not meet acceptance criteria for dichlorodifluoromethane and acetone.

VOC 8260: LCS 10171-53 did not meet the acceptance limits for 2,2 dichloropropane, bromomethane, styrene and dichlorodifluoromethane. The RPD for styrene did not meet the acceptance criteria.

These compounds noted with failures are known to be problematic in the method.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

EPH: The fractionation check sample (LCS) for the batch of silica gel in use for these samples met the method acceptance criteria.

VPH Target compounds and ranges were determined by GC/MS. Ranges were determined in a similar manner as described in the MADEP APH method of 2/2000.

RESOURCE LABORATORIES, LLC.

(continued)

PCB: The sample peaks most closely resembled Aroclor 1242, however there was also a similarity to Aroclor 1248. Possibly due to weathering, the Aroclor identification is difficult and not absolute. Chromatograms are provided. Samples that contained non-aroclor peaks have been noted on the report pages. Quantification is quadratic.

VOC: The following compounds were quantified using quadratic fit:
Dichlorodifluoromethane, chloromethane, vinyl chloride, bromomethane, acetone, dibromochloromethane, isopropylbenzene, tert-butyl benzene, and hexachlorobutadiene.

Dilutions performed during the analysis are noted on the result pages.

Metals: RCRA metals only were requested by the customer.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 10171

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|-----------|----------|----------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 (x) | Cyanide () | Other () |
| | 8270 () | VPH (x) | 7470/7471 (x) | Other () | Other () |
| | 8082 (x) | EPH (x) | Other () | Other () | Other () |

| | | |
|---|--|-----------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No (x) |
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes () No (x) |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes () No (x) |

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Position: Lab DirectorPrinted Name: Susan C. SylvesterDate: 4-27-06

Lab Number: 10171-50
Sample Designation: Method blank 041806A
Date Sampled: N/A
Date Analyzed: 4/18/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethene | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 94 | 78-114 | | | |
| toluene-D8 | 107 | 88-110 | | | |
| 4-bromofluorobenzene | 98 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-51
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: X:\DATA\VOA03\2006\APR06\041806\W3041822.D
Date Analyzed: 4/18/06
SW 846 Method 5030B/8260B

| Compound | LCS | | | LCSD | | RPD |
|-----------------------------|--------------|------------|---|--------------|------------|-----|
| | Amount Found | % Recovery | | Amount Found | % Recovery | |
| dichlorodifluoromethane | 14 | 70% | * | 14 | 72% | 3% |
| chloromethane | 16 | 80% | | 18 | 88% | 9% |
| vinyl chloride | 17 | 86% | | 18 | 92% | 7% |
| bromomethane | 16 | 82% | | 18 | 88% | 7% |
| chloroethane | 17 | 86% | | 18 | 91% | 5% |
| trichlorofluoromethane | 18 | 90% | | 19 | 97% | 7% |
| diethylether | 18 | 92% | | 18 | 91% | 1% |
| acetone | 13 | 67% | * | 14 | 72% | 7% |
| 1,1-dichloroethene | 17 | 86% | | 19 | 95% | 10% |
| methylene chloride | 19 | 97% | | 19 | 97% | 0% |
| carbon disulfide | 18 | 90% | | 19 | 94% | 4% |
| methyl-1-butyl ether (MTBE) | 35 | 89% | | 38 | 90% | 1% |
| trans-1,2-dichloroethene | 18 | 91% | | 20 | 98% | 7% |
| isopropyl ether (DIPE) | 18 | 92% | | 19 | 97% | 5% |
| ethyl-1-butyl ether (ETBE) | 18 | 91% | | 19 | 96% | 8% |
| 1,1-dichloroethane | 17 | 88% | | 18 | 92% | 7% |
| 1-butanol (TBA) | 74 | 74% | | 78 | 78% | 6% |
| 2-butanone (MEK) | 17 | 83% | | 16 | 81% | 3% |
| 2,2-dichloropropane | 17 | 83% | | 18 | 89% | 7% |
| cis-1,2-dichloroethene | 20 | 101% | | 21 | 107% | 7% |
| chloroform | 19 | 95% | | 20 | 102% | 7% |
| bromochloromethane | 21 | 104% | | 21 | 107% | 3% |
| tetrahydrofuran (THF) | 17 | 85% | | 18 | 88% | 4% |
| 1,1,1-trichloroethane | 17 | 87% | | 19 | 96% | 9% |
| 1,1-dichloropropene | 19 | 95% | | 20 | 102% | 7% |
| 1-amyl-methyl ether (TAME) | 19 | 94% | | 19 | 96% | 3% |
| carbon tetrachloride | 17 | 83% | | 18 | 90% | 8% |
| 1,2-dichloroethane | 18 | 88% | | 18 | 90% | 2% |
| benzene | 20 | 100% | | 21 | 106% | 5% |
| trichloroethene | 20 | 101% | | 21 | 107% | 6% |
| 1,2-dichloropropane | 20 | 99% | | 21 | 104% | 5% |
| bromodichloromethane | 17 | 84% | | 18 | 89% | 6% |
| dibromomethane | 20 | 101% | | 21 | 105% | 4% |
| 4-methyl-2-pentanone (MIBK) | 19 | 95% | | 19 | 97% | 2% |
| cis-1,3-dichloropropene | 20 | 98% | | 20 | 102% | 4% |
| toluene | 21 | 105% | | 22 | 110% | 5% |
| trans-1,3-dichloropropene | 17 | 85% | | 18 | 90% | 6% |
| 2-hexanone | 17 | 84% | | 18 | 89% | 6% |
| 1,1,2-trichloroethane | 20 | 102% | | 21 | 106% | 4% |
| 1,3-dichloropropane | 21 | 105% | | 22 | 108% | 3% |
| tetrachloroethene | 22 | 111% | | 24 | 120% | 8% |
| dibromochloromethane | 20 | 99% | | 20 | 100% | 2% |
| 1,2-dibromoethane (EDB) | 21 | 106% | | 21 | 106% | 1% |
| chlorobenzene | 21 | 107% | | 22 | 112% | 4% |
| 1,1,1,2-tetrachloroethane | 20 | 102% | | 21 | 107% | 4% |
| ethylbenzene | 21 | 106% | | 22 | 108% | 2% |
| m,p-xylenes | 44 | 109% | | 45 | 114% | 4% |
| o-xylene | 22 | 111% | | 23 | 116% | 4% |
| styrene | 21 | 106% | | 22 | 110% | 4% |
| bromoform | 20 | 102% | | 21 | 103% | 1% |
| isopropylbenzene | 22 | 112% | | 24 | 119% | 6% |
| 1,1,2,2-tetrachloroethane | 18 | 92% | | 18 | 91% | 2% |
| 1,2,3-trichloropropane | 18 | 88% | | 17 | 87% | 1% |
| n-propylbenzene | 21 | 103% | | 21 | 107% | 4% |
| bromobenzene | 21 | 103% | | 22 | 108% | 5% |
| 1,3,5-trimethylbenzene | 19 | 97% | | 20 | 100% | 2% |
| 2-chlorotoluene | 19 | 95% | | 20 | 99% | 4% |
| 4-chlorotoluene | 19 | 97% | | 20 | 100% | 4% |
| tert-butylbenzene | 17 | 87% | | 18 | 89% | 3% |
| 1,2,4-trimethylbenzene | 19 | 94% | | 20 | 100% | 6% |
| sec-butylbenzene | 19 | 97% | | 20 | 102% | 6% |
| 1,3-dichlorobenzene | 20 | 100% | | 20 | 102% | 2% |
| 4-isopropyltoluene | 21 | 104% | | 22 | 110% | 6% |
| 1,4-dichlorobenzene | 19 | 96% | | 20 | 100% | 4% |
| 1,2-dichlorobenzene | 20 | 102% | | 21 | 104% | 2% |
| n-butylbenzene | 20 | 100% | | 21 | 107% | 7% |
| 1,2-dibromo-3-chloropropane | 17 | 83% | | 17 | 87% | 5% |
| 1,2,4-trichlorobenzene | 18 | 90% | | 19 | 97% | 8% |
| hexachlorobutadiene | 19 | 97% | | 21 | 104% | 7% |
| naphthalene | 14 | 72% | | 16 | 78% | 8% |
| 1,2,3-trichlorobenzene | 18 | 91% | | 19 | 98% | 5% |
| 1,4-dioxane | 37 | 92% | | 33 | 82% | 12% |
| SURROGATE STANDARDS | | | | | | |
| SS dibromofluoromethane | | 92% | | | 99% | |
| SS toluene-D8 | | 105% | | | 108% | |
| SS 4-bromofluorobenzene | | 107% | | | 101% | |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

Lab Number: 10171-52
Sample Designation: Method Blank 041906
Date Sampled: N/A
Date Analyzed: 4/19/06
Matrix: Water
Instrument Dilution Factor: 1
Analyst: LMM

VOLATILE ORGANICS
SW 846 Method 5030B/8260B

| | Concentration | Quantitation Limit | | Concentration | Quantitation Limit |
|-----------------------------|---------------|--------------------|-----------------------------|---------------|--------------------|
| | ug/L | ug/L | | ug/L | ug/L |
| dichlorodifluoromethane | U | 2 | trans-1,3-dichloropropene | U | 2 |
| chloromethane | U | 2 | 2-hexanone | U | 10 |
| vinyl chloride | U | 2 | 1,1,2-trichloroethane | U | 2 |
| bromomethane | U | 2 | 1,3-dichloropropane | U | 2 |
| chloroethane | U | 2 | tetrachloroethene | U | 2 |
| trichlorofluoromethane | U | 2 | dibromochloromethane | U | 2 |
| diethyl ether | U | 10 | 1,2-dibromoethane | U | 2 |
| acetone | U | 10 | chlorobenzene | U | 2 |
| 1,1-dichloroethene | U | 1 | 1,1,1,2-tetrachloroethane | U | 2 |
| methylene chloride | U | 5 | ethylbenzene | U | 2 |
| carbon disulfide | U | 2 | m&p-xylenes | U | 2 |
| methyl t-butyl ether (MTBE) | U | 2 | o-xylene | U | 2 |
| trans-1,2-dichloroethene | U | 2 | styrene | U | 2 |
| Isopropyl ether (DIPE) | U | 2 | bromoform | U | 2 |
| ethyl t-butyl ether (ETBE) | U | 2 | isopropylbenzene | U | 2 |
| 1,1-dichloroethane | U | 2 | 1,1,2,2-tetrachloroethane | U | 2 |
| t-butanol (TBA) | U | 50 | 1,2,3-trichloropropane | U | 2 |
| 2-butanone (MEK) | U | 10 | n-propylbenzene | U | 2 |
| 2,2-dichloropropane | U | 2 | bromobenzene | U | 2 |
| cis-1,2-dichloroethane | U | 2 | 1,3,5-trimethylbenzene | U | 2 |
| chloroform | U | 2 | 2-chlorotoluene | U | 2 |
| bromochloromethane | U | 2 | 4-chlorotoluene | U | 2 |
| tetrahydrofuran (THF) | U | 10 | tert-butylbenzene | U | 2 |
| 1,1,1-trichloroethane | U | 2 | 1,2,4-trimethylbenzene | U | 2 |
| 1,1-dichloropropene | U | 2 | sec-butylbenzene | U | 2 |
| t-amyl-methyl ether (TAME) | U | 2 | 1,3-dichlorobenzene | U | 2 |
| carbon tetrachloride | U | 2 | 4-isopropyltoluene | U | 2 |
| 1,2-dichloroethane | U | 2 | 1,4-dichlorobenzene | U | 2 |
| benzene | U | 2 | 1,2-dichlorobenzene | U | 2 |
| trichloroethene | U | 2 | n-butylbenzene | U | 2 |
| 1,2-dichloropropane | U | 2 | 1,2-dibromo-3-chloropropane | U | 2 |
| bromodichloromethane | U | 2 | 1,2,4-trichlorobenzene | U | 2 |
| dibromomethane | U | 2 | hexachlorobutadiene | U | 2 |
| 4-methyl-2-pentanone (MIBK) | U | 10 | naphthalene | U | 5 |
| cis-1,3-dichloropropene | U | 2 | 1,2,3-trichlorobenzene | U | 2 |
| toluene | U | 2 | 1,4-dioxane | U | 50 |
| | | | | | |
| SURROGATE STANDARDS | Recovery | Acceptance Limits | | | |
| | (%) | (%) | | | |
| dibromofluoromethane | 93 | 78-114 | | | |
| toluene-D8 | 100 | 88-110 | | | |
| 4-bromofluorobenzene | 99 | 86-115 | | | |

U = Below quantitation limit

Lab Number: 10171-53
Sample Designation: Lab Control Sample/Lab Control Sample Duplicate
File Name: X:\DATA\AWOA03\2006\APR06\041806W3041844.D
Date Analyzed: 4/19/06
SW 846 Method 5030B/8280B

| Compound | LCS | | LCSD | | RPD |
|-----------------------------|--------------|------------|--------------|------------|-----|
| | Amount Found | % Recovery | Amount Found | % Recovery | |
| dichlorodifluoromethane | 14 | 70% | 13 | 66% | 5% |
| chloromethane | 17 | 84% | 16 | 79% | 6% |
| vinyl chloride | 17 | 87% | 17 | 85% | 3% |
| bromomethane | 11 | 54% | 13 | 63% | 15% |
| chloroethane | 18 | 88% | 17 | 84% | 4% |
| trichlorofluoromethane | 19 | 93% | 18 | 88% | 5% |
| diethylether | 19 | 93% | 19 | 94% | 1% |
| acetone | 18 | 88% | 18 | 88% | 1% |
| 1,1-dichloroethane | 17 | 87% | 18 | 88% | 1% |
| methylene chloride | 20 | 99% | 19 | 94% | 5% |
| carbon disulfide | 18 | 88% | 17 | 84% | 5% |
| methyl-t-butyl ether (MTBE) | 38 | 95% | 37 | 93% | 3% |
| trans-1,2-dichloroethane | 18 | 91% | 18 | 90% | 2% |
| isopropyl ether (DIPE) | 18 | 95% | 18 | 90% | 5% |
| ethyl-t-butyl ether (ETBE) | 19 | 97% | 19 | 94% | 4% |
| 1,1-dichloroethane | 18 | 91% | 17 | 86% | 5% |
| 1-butanol (TBA) | 90 | 90% | 88 | 88% | 3% |
| 2-butanone (MEK) | 19 | 93% | 18 | 89% | 4% |
| 2,2-dichloropropane | 13 | 87% | 13 | 66% | 2% |
| cis-1,2-dichloroethane | 21 | 104% | 20 | 99% | 5% |
| chloroform | 19 | 95% | 19 | 98% | 1% |
| bromochloromethane | 21 | 106% | 21 | 104% | 2% |
| tetrahydrofuran (THF) | 19 | 96% | 19 | 95% | 1% |
| 1,1,1-trichloroethane | 20 | 98% | 18 | 91% | 7% |
| 1,1-dichloropropene | 19 | 97% | 19 | 96% | 1% |
| 1-amyl-methyl ether (TAME) | 20 | 100% | 19 | 94% | 6% |
| carbon tetrachloride | 17 | 87% | 17 | 83% | 4% |
| 1,2-dichloroethane | 18 | 88% | 18 | 89% | 2% |
| benzene | 20 | 88% | 20 | 99% | 1% |
| trichloroethane | 21 | 103% | 20 | 101% | 2% |
| 1,2-dichloropropane | 20 | 102% | 20 | 100% | 2% |
| bromodichloromethane | 18 | 88% | 17 | 87% | 1% |
| dibromomethane | 21 | 103% | 21 | 104% | 1% |
| 4-methyl-2-pentanone (MIBK) | 21 | 104% | 20 | 100% | 4% |
| cis-1,3-dichloropropene | 18 | 92% | 18 | 89% | 3% |
| toluene | 20 | 102% | 21 | 103% | 1% |
| trans-1,3-dichloropropene | 17 | 85% | 18 | 81% | 4% |
| 2-hexanone | 16 | 82% | 17 | 85% | 4% |
| 1,1,2-trichloroethane | 21 | 106% | 20 | 101% | 5% |
| 1,3-dichloropropane | 20 | 102% | 21 | 104% | 3% |
| tetrachloroethene | 22 | 108% | 22 | 111% | 3% |
| dibromochloromethane | 19 | 96% | 20 | 98% | 2% |
| 1,2-dibromoethane (EDB) | 20 | 99% | 21 | 104% | 4% |
| chlorobenzene | 20 | 98% | 20 | 102% | 4% |
| 1,1,1,2-tetrachloroethane | 20 | 99% | 20 | 100% | 2% |
| ethylbenzene | 20 | 98% | 20 | 100% | 3% |
| m,p-xylenes | 41 | 102% | 43 | 106% | 4% |
| o-xylene | 20 | 102% | 21 | 106% | 4% |
| styrene | 8 | 42% | 14 | 72% | 53% |
| bromoform | 19 | 97% | 20 | 101% | 3% |
| isopropylbenzene | 21 | 104% | 22 | 108% | 4% |
| 1,1,2,2-tetrachloroethane | 18 | 90% | 19 | 94% | 4% |
| 1,2,3-trichloropropane | 18 | 88% | 18 | 89% | 1% |
| n-propylbenzene | 19 | 93% | 20 | 100% | 7% |
| bromobenzene | 19 | 96% | 21 | 103% | 7% |
| 1,3,5-trimethylbenzene | 17 | 87% | 18 | 97% | 10% |
| 2-chlorotoluene | 17 | 86% | 19 | 94% | 9% |
| 4-chlorotoluene | 18 | 88% | 19 | 95% | 8% |
| tert-butylbenzene | 17 | 83% | 17 | 87% | 4% |
| 1,2,4-trimethylbenzene | 18 | 90% | 19 | 97% | 7% |
| sec-butylbenzene | 18 | 91% | 18 | 92% | 0% |
| 1,3-dichlorobenzene | 19 | 94% | 20 | 99% | 5% |
| 4-isopropyltoluene | 19 | 94% | 20 | 102% | 8% |
| 1,4-dichlorobenzene | 18 | 92% | 19 | 96% | 5% |
| 1,2-dichlorobenzene | 18 | 96% | 20 | 100% | 4% |
| n-butylbenzene | 19 | 95% | 20 | 100% | 5% |
| 1,2-dibromo-3-chloropropane | 17 | 85% | 19 | 94% | 10% |
| 1,2,4-trichlorobenzene | 17 | 86% | 19 | 94% | 8% |
| hexachlorobutadiene | 18 | 90% | 19 | 95% | 5% |
| naphthalene | 15 | 74% | 16 | 80% | 8% |
| 1,2,3-trichlorobenzene | 18 | 91% | 20 | 100% | 9% |
| 1,4-dioxane | 35 | 89% | 38 | 96% | 8% |

SURROGATE STANDARDS

| | | |
|-------------------------|------|------|
| SS dibromofluoromethane | 99% | 99% |
| SS toluene-D8 | 108% | 108% |
| SS 4-bromofluorobenzene | 107% | 101% |

Preliminary Acceptance Criteria: Recovery 70% - 130% RPD 20%

* Indicates compounds known to be problematic. These analytes are more likely to show recovery outside the QC limits.

= This analyte showed recovery outside the acceptance limits.

+ = The RPD was above the acceptance limit.

Lab Number: 10171-54
 Sample Designation: Method Blank 041806
 Date Sampled: N/A
 Date Received: N/A
 Date Analyzed: 4/18/06
 Matrix: Water
 Containers: Satisfactory
 Sample Preservation: N/A
 Temperature: N/A
 Instrument Dilution Factor: 1
 Analyst: LMM

VPH ANALYTICAL RESULTS
 MADEP VPH 2004-1.1

| | Elution Range | Concentration ug/L | Quantitation Limit ug/L |
|----------------------------------|------------------|-----------------------|----------------------------|
| Unadjusted C5-C8 Aliphatics (1) | N/A | U | 100 |
| Unadjusted C9-C12 Aliphatics (1) | N/A | U | 100 |
| methyl-t-butyl ether (MTBE) | C5-C8 | U | 2 |
| benzene | C5-C8 | U | 1 |
| toluene | C5-C8 | U | 2 |
| ethylbenzene | C9-C12 | U | 2 |
| m&p-xylenes | C9-C12 | U | 2 |
| o-xylene | C9-C12 | U | 2 |
| naphthalene | N/A | U | 5 |
| C5-C8 Aliphatics (1,2) | N/A | U | 100 |
| C9-C12 Aliphatics (1,3) | N/A | U | 100 |
| C9-C10 Aromatics (1) | N/A | U | 100 |
| Surrogate Recovery | | | |
| 2,5-dibromotoluene as aromatic | | 92% | |
| 2,5-dibromotoluene as aliphatic | | 93% | |
| Surrogate Acceptance Range | | 70-130% | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C5-C8 Aliphatic Hydrocarbons exclude the concentration of Target Analytes eluting in that range.

3 C9-C12 Aliphatic Hydrocarbons exclude conc of Target Analytes eluting in that range AND concentration of C9-C10 Aromatic Hydrocarbons.

U = Below quantitation limit

Lab Number: 10171-55
 Sample Designation: LCS/LCSD
 Date Sampled: N/A
 Date Received: N/A
 Date Analyzed: 4/19/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: LMM

VOLATILE ORGANICS MS/MSD LCS/LCSD Report
 MADEP VPH 2004-1.1

| Compound | Spike Added (ug/L) | Sample Concentration (ug/L) | MS/LCS Concentration (ug/L) | % Recovery | QC Lower Limit | QC Upper Limit |
|---------------------------------|--------------------|-----------------------------|-----------------------------|------------|----------------|----------------|
| methyl-t-butyl ether (MTBE) | 150 | | 138 | 92% | 70% | 130% |
| benzene | 50 | | 47 | 93% | 70% | 130% |
| toluene | 150 | | 133 | 89% | 70% | 130% |
| ethylbenzene | 50 | | 45 | 89% | 70% | 130% |
| m&p-xylenes | 200 | | 186 | 93% | 70% | 130% |
| o-xylene | 100 | | 89 | 89% | 70% | 130% |
| naphthalene | 100 | | 91 | 91% | 70% | 130% |
| C5-C8 Aliphatics | 750 | | 638 | 85% | 70% | 130% |
| C9-C12 Aliphatics | 550 | | 445 | 81% | 70% | 130% |
| C9-C10 Aromatics | 100 | | 94 | 94% | 70% | 130% |
| 2,5-dibromotoluene as aromatic | | | | 101% | 70% | 130% |
| 2,5-dibromotoluene as aliphatic | | | | 102% | 70% | 130% |

| Compound | Spike Added (ug/L) | Sample Concentration (ug/L) | MSD/LCSD Concentration (ug/L) | % Recovery | QC Lower Limit | QC Upper Limit | RPD | RPD Limit |
|---------------------------------|--------------------|-----------------------------|-------------------------------|------------|----------------|----------------|-----|-----------|
| methyl-t-butyl ether (MTBE) | 150 | | 140 | 93% | 70% | 130% | 1% | 20 |
| benzene | 50 | | 47 | 94% | 70% | 130% | 1% | 20 |
| toluene | 150 | | 137 | 92% | 70% | 130% | 3% | 20 |
| ethylbenzene | 50 | | 47 | 93% | 70% | 130% | 4% | 20 |
| m&p-xylenes | 200 | | 191 | 96% | 70% | 130% | 3% | 20 |
| o-xylene | 100 | | 93 | 93% | 70% | 130% | 4% | 20 |
| naphthalene | 100 | | 92 | 92% | 70% | 130% | 2% | 20 |
| C5-C8 Aliphatics | 750 | | 653 | 87% | 70% | 130% | 2% | 20 |
| C9-C12 Aliphatics | 550 | | 459 | 84% | 70% | 130% | 3% | 20 |
| C9-C10 Aromatics | 100 | | 101 | 101% | 70% | 130% | 7% | 20 |
| 2,5-dibromotoluene as aromatic | | | | 101% | 70% | 130% | 0% | 20 |
| 2,5-dibromotoluene as aliphatic | | | | 104% | 70% | 130% | 1% | 20 |

Lab Number: 10171-50
 Sample Designation: Preparation Blank 181 W
 Date Sampled: N/A
 Date Received: N/A
 Date Extracted: 4/18/06
 Matrix: Water
 Containers: N/A
 Sample Preservation: N/A
 Temperature: N/A
 Dilution Factor: 1
 Analyst: AJD

EPH ANALYTICAL RESULTS

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | Concentration ug/L | Quantitation Limit ug/L | Date of Analysis |
|-------------------------------------|-----------------------|----------------------------|------------------|
| Diesel PAH Analytes: | | | |
| naphthalene | U | 0.5 | 4/18/06 |
| 2-methylnaphthalene | U | 0.5 | 4/18/06 |
| phenanthrene | U | 0.5 | 4/18/06 |
| acenaphthene | U | 0.5 | 4/18/06 |
| Other PAH Analytes: | | | |
| acenaphthylene | U | 0.5 | 4/18/06 |
| fluorene | U | 0.5 | 4/18/06 |
| anthracene | U | 0.5 | 4/18/06 |
| fluoranthene | U | 0.5 | 4/18/06 |
| pyrene | U | 0.5 | 4/18/06 |
| benzo(a)anthracene | U | 0.5 | 4/18/06 |
| chrysene | U | 0.5 | 4/18/06 |
| benzo(b)fluoranthene | U | 0.5 | 4/18/06 |
| benzo(k)fluoranthene | U | 0.5 | 4/18/06 |
| benzo(a)pyrene | U | 0.2 | 4/18/06 |
| indeno(1,2,3-cd)pyrene | U | 0.5 | 4/18/06 |
| dibenzo(a,h)anthracene | U | 0.5 | 4/18/06 |
| benzo(g,h,i)perylene | U | 0.5 | 4/18/06 |
| Ranges: | | | |
| Unadjusted C11-C22 Aromatics (1) | U | 200 | 4/19/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | U | 200 | 4/19/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | U | 200 | 4/19/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | U | 200 | 4/19/06 |
| Extraction Surrogate Recoveries: | | | |
| 1-chloro-octadecane | 46% | | |
| o-terphenyl | 53% | | |
| Fractionation Surrogate Recoveries: | | | |
| 2-fluorobiphenyl | 98% | | |
| 2-bromonaphthalene | 90% | | |
| Surrogate Acceptance Range | 40-140% | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

Lab Number: 10171-51
 Sample Designation: Lab Control Sample/Duplicate 181 W
 Date Sampled: N/A
 Date Received: N/A
 Date Extracted: 4/18/06
 Matrix: Water
 Containers: N/A
 Sample Preservation: N/A
 Temperature: N/A
 Dilution Factor: 1
 Analyst: A.JD

EPH ANALYTICAL RESULTS

| Method for Ranges: MADEP EPH 2004-1.1 | Amount Added ug/L | Amount Found in LCS ug/L | LCS Recovery (%) | Amount Found in LCSD ug/L | LCSD Recovery (%) | RPD (%) | Acceptance Criteria (%) | Date of Analysis |
|---------------------------------------|-------------------------|-----------------------------------|------------------------|------------------------------------|-------------------------|------------|-------------------------------|------------------|
| Method for Target Analytes: EPA 8270C | | | | | | | | |
| Diesel PAH Analytes: | | | | | | | | |
| naphthalene | 60 | 32.97 | 55% | 29.46 | 49% | -11% | <25% | 4/19/06 |
| 2-methylnaphthalene | 60 | 34.12 | 57% | 29.66 | 49% | -14% | <25% | 4/19/06 |
| phenanthrene | 60 | 42.98 | 72% | 38.40 | 64% | -11% | <25% | 4/19/06 |
| acenaphthene | 60 | 41.06 | 68% | 36.16 | 60% | -13% | <25% | 4/19/06 |
| Other PAH Analytes: | | | | | | | | |
| acenaphthylene | 60 | 38.41 | 64% | 33.49 | 56% | -14% | <25% | 4/19/06 |
| fluorene | 60 | 40.28 | 67% | 35.86 | 60% | -12% | <25% | 4/19/06 |
| anthracene | 60 | 45.02 | 75% | 40.95 | 68% | -9% | <25% | 4/19/06 |
| fluoranthene | 60 | 48.34 | 81% | 42.15 | 70% | -14% | <25% | 4/19/06 |
| pyrene | 60 | 45.14 | 75% | 40.17 | 67% | -12% | <25% | 4/19/06 |
| benzo(a)anthracene | 60 | 47.06 | 78% | 41.83 | 70% | -12% | <25% | 4/19/06 |
| chrysene | 60 | 48.65 | 81% | 42.52 | 71% | -13% | <25% | 4/19/06 |
| benzo(b)fluoranthene | 60 | 38.30 | 64% | 36.83 | 61% | -4% | <25% | 4/19/06 |
| benzo(k)fluoranthene | 60 | 59.95 | 100% | 47.64 | 79% | -23% | <25% | 4/19/06 |
| benzo(a)pyrene | 60 | 51.65 | 86% | 45.99 | 77% | -12% | <25% | 4/19/06 |
| indeno(1,2,3-cd)pyrene | 60 | 49.36 | 82% | 43.65 | 73% | -12% | <25% | 4/19/06 |
| dibenzo(a,h)anthracene | 60 | 49.35 | 82% | 43.30 | 72% | -13% | <25% | 4/19/06 |
| benzo(g,h,i)perylene | 60 | 49.46 | 82% | 43.35 | 72% | -13% | <25% | 4/19/06 |
| Ranges: | | | | | | | | |
| Unadjusted C11-C22 Aromatics (1) | 1020 | 685.66 | 67% | 688.84 | 68% | 1% | <25% | 4/19/06 |
| C9-C18 Aliphatic Hydrocarbons (1) | 360 | 261.88 | 73% | 233.25 | 65% | -12% | <25% | 4/19/06 |
| C19-C36 Aliphatic Hydrocarbons (1) | 480 | 510.83 | 106% | 456.23 | 95% | -11% | <25% | 4/19/06 |
| C11-C22 Aromatic Hydrocarbons (1,2) | | | | | | | | 4/19/06 |
| Extraction Surrogate Recoveries: | | | | | | | | |
| 1-chloro-octadecane | | 46% | | 42% | | | | |
| o-terphenyl | | 59% | | 57% | | | | |
| Fractionation Surrogate Recoveries: | | | | | | | | |
| 2-fluorobiphenyl | | 102% | | 87% | | | | |
| 2-bromonaphthalene | | 92% | | 81% | | | | |
| Acceptance Range | | 40-140% | 40-140% | 40-140% | 40-140% | | | |

1 Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

2 C11-C22 Aromatic Hydrocarbons exclude the concentration of target PAH Analytes.

U = Below quantitation limit

Lab Number: 10171-52
 Sample Designation: LCS 181 W
 Date Sampled: N/A
 Date Received: N/A
 Date Extracted: 4/18/06
 Matrix: Water
 Containers: N/A
 Sample Preservation: N/A
 Temperature: N/A
 Dilution Factor: 1
 Analyst: AJD

NAPHTHALENE BREAKTHROUGH CALCULATION

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | LCS Aliphatic Breakthrough (%) | Acceptance Criteria | Date of Analysis |
|---------------------|--------------------------------------|------------------------|------------------|
| naphthalene | 2.5% | <5.0% | 4/19/06 |
| 2-methylnaphthalene | 1.5% | <5.0% | 4/19/06 |

Lab Number: 10171-53
 Sample Designation: LCSD 181 W
 Date Sampled: N/A
 Date Received: N/A
 Date Extracted: 4/18/06
 Matrix: Water
 Containers: N/A
 Sample Preservation: N/A
 Temperature: N/A
 Dilution Factor: 1
 Analyst: AJD

NAPHTHALENE BREAKTHROUGH CALCULATION

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

| | LCSD | | |
|---------------------|------------------------|------------|------------------|
| | Aliphatic Breakthrough | Acceptance | Date of Analysis |
| | (%) | Criteria | |
| naphthalene | 3.4% | <5.0% | 4/19/06 |
| 2-methylnaphthalene | 2.0% | <5.0% | 4/19/06 |

Lab Number: 10171-50
 Sample Designation: Preparation Blank 209 W
 Date Sampled: N/A
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 74 | 30-150 |
| Decachlorobiphenyl | 86 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-51
 Sample Designation: Laboratory Control Sample/Duplicate 209 W
 Date Sampled: N/A
 Date Extracted: 4/12/06
 Date Analyzed: 4/13/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: A.J.D

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Amount Added ug/L | Amount Found in LCS ug/L | LCS Recovery (%) | Amount Found in LCSD ug/L | LCSD Recovery (%) | RPD (%) | Date of Analysis |
|----------------------|-------------------------|-----------------------------------|------------------------|------------------------------------|-------------------------|------------|------------------|
| PCB-1016 | 2 | 1.5 | 75% | 1.5 | 75% | 0.0% | 4/13/06 |
| PCB-1242 | | | | | | | 4/13/06 |
| PCB-1221 | | | | | | | 4/13/06 |
| PCB-1232 | | | | | | | 4/13/06 |
| PCB-1248 | | | | | | | 4/13/06 |
| PCB-1254 | | | | | | | 4/13/06 |
| PCB-1260 | 2 | 1.4 | 70% | 1.5 | 75% | 6.9% | 4/13/06 |
| PCB-1262 | | | | | | | 4/13/06 |
| PCB-1268 | | | | | | | 4/13/06 |
| Acceptance Criteria: | | | 40-140% | | 40-140% | 20% | |

| SURROGATE STANDARDS | Recovery (%) | Acceptance Limits (%) | Recovery (%) | Acceptance Limits (%) |
|----------------------|-----------------|--------------------------|-----------------|--------------------------|
| Tetrachloro-m-xylene | 83 | 30-150 | 78 | 30-150 |
| Decachlorobiphenyl | 88 | 30-150 | 92 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-60
 Sample Designation: Preparation Blank 210 W
 Date Sampled: N/A
 Date Extracted: 4/14/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: AJD

POLYCHLORINATED BIPHENYLS

SW 846 Method 3510C/8082A.

| | Concentration | Quantitation Limit |
|----------|---------------|--------------------|
| | ug/L | ug/L |
| PCB-1016 | U | 0.3 |
| PCB-1242 | U | 0.3 |
| PCB-1221 | U | 0.3 |
| PCB-1232 | U | 0.3 |
| PCB-1248 | U | 0.3 |
| PCB-1254 | U | 0.3 |
| PCB-1260 | U | 0.3 |
| PCB-1262 | U | 0.3 |
| PCB-1268 | U | 0.3 |

| SURROGATE STANDARDS | Recovery | Acceptance Limits |
|----------------------|----------|-------------------|
| | (%) | (%) |
| Tetrachloro-m-xylene | 80 | 30-150 |
| Decachlorobiphenyl | 96 | 30-150 |

U = Below quantitation limit

Lab Number: 10171-61
 Sample Designation: Laboratory Control Sample/Duplicate 210 W
 Date Sampled: N/A
 Date Extracted: 4/12/06
 Date Analyzed: 4/18/06
 Matrix: Water
 Dilution Factor: 1
 Analyst: A.JD

POLYCHLORINATED BIPHENYLS
 SW 846 Method 3510C/8082A.

| | Amount Added ug/L | Amount Found in LCS ug/L | LCS Recovery (%) | Amount Found in LCSD ug/L | LCSD Recovery (%) | RPD (%) | Date of Analysis |
|----------------------|-------------------------|-----------------------------------|------------------------|------------------------------------|-------------------------|------------|------------------|
| PCB-1016 | 2 | 1.5 | 75% | 1.8 | 90% | 18% | 4/18/06 |
| PCB-1242 | | | | | | | 4/18/06 |
| PCB-1221 | | | | | | | 4/18/06 |
| PCB-1232 | | | | | | | 4/18/06 |
| PCB-1248 | | | | | | | 4/18/06 |
| PCB-1254 | | | | | | | 4/18/06 |
| PCB-1260 | 2 | 1.7 | 85% | 1.8 | 90% | 6% | 4/18/06 |
| PCB-1262 | | | | | | | 4/18/06 |
| PCB-1268 | | | | | | | 4/18/06 |
| Acceptance Criteria: | | | 40-140% | | 40-140% | 20% | |

| SURROGATE STANDARDS | Recovery (%) | Acceptance Limits (%) | Recovery (%) | Acceptance Limits (%) |
|----------------------|-----------------|--------------------------|-----------------|--------------------------|
| Tetrachloro-m-xylene | 79 | 30-150 | 90 | 30-150 |
| Decachlorobiphenyl | 101 | 30-150 | 102 | 30-150 |

U = Below quantitation limit

Lab Number 10171 METALS QC REPORT
Batch QC Results

Prep Blank

| Analyte | Sample ID # | Result (mg/L) | Reporting Limit (mg/L) |
|----------------|--------------------|--------------------------|-----------------------------------|
| Silver | ICB041306 | < 0.007 | 0.007 |
| Arsenic | ICB041306 | < 0.01 | 0.01 |
| Barium | ICB041306 | < 0.05 | 0.05 |
| Cadmium | ICB041306 | < 0.005 | 0.005 |
| Chromium | ICB041306 | < 0.05 | 0.05 |
| Lead | ICB041306 | < 0.01 | 0.01 |
| Selenium | ICB041306 | < 0.05 | 0.05 |

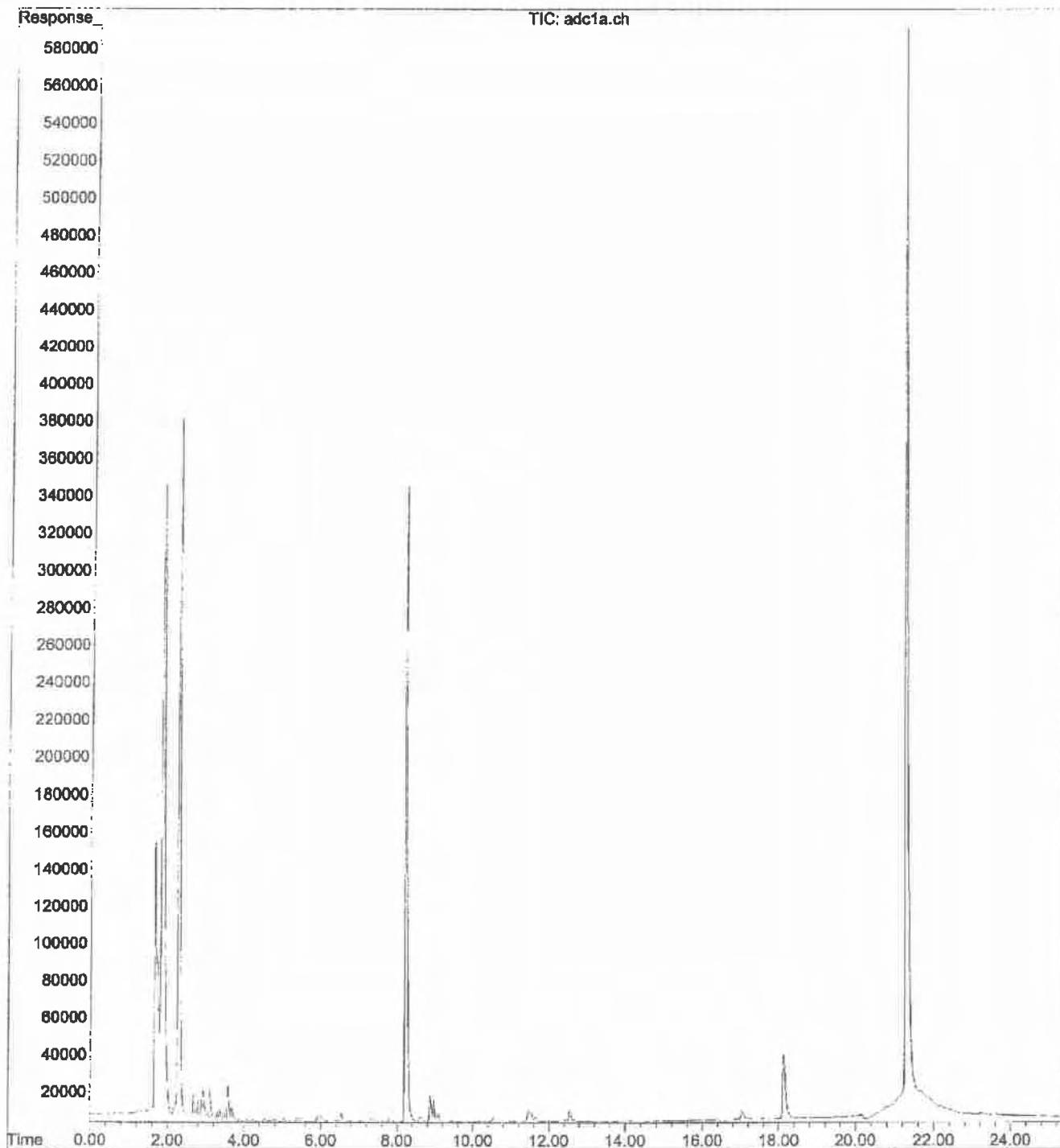
Laboratory Control Sample

| Analyte | Sample ID # | Result (mg/L) | True Value (mg/L) | %Recovery Control Limits (85-115%) |
|----------------|--------------------|--------------------------|------------------------------|---|
| Silver | ICV041306 | 0.26 | 0.25 | 103 |
| Arsenic | ICV041306 | 0.49 | 0.50 | 99 |
| Barium | ICV041306 | 0.51 | 0.50 | 101 |
| Cadmium | ICV041306 | 0.51 | 0.50 | 102 |
| Chromium | ICV041306 | 0.50 | 0.50 | 100 |
| Lead | ICV041306 | 0.49 | 0.50 | 97 |
| Selenium | ICV041306 | 0.50 | 0.50 | 99 |

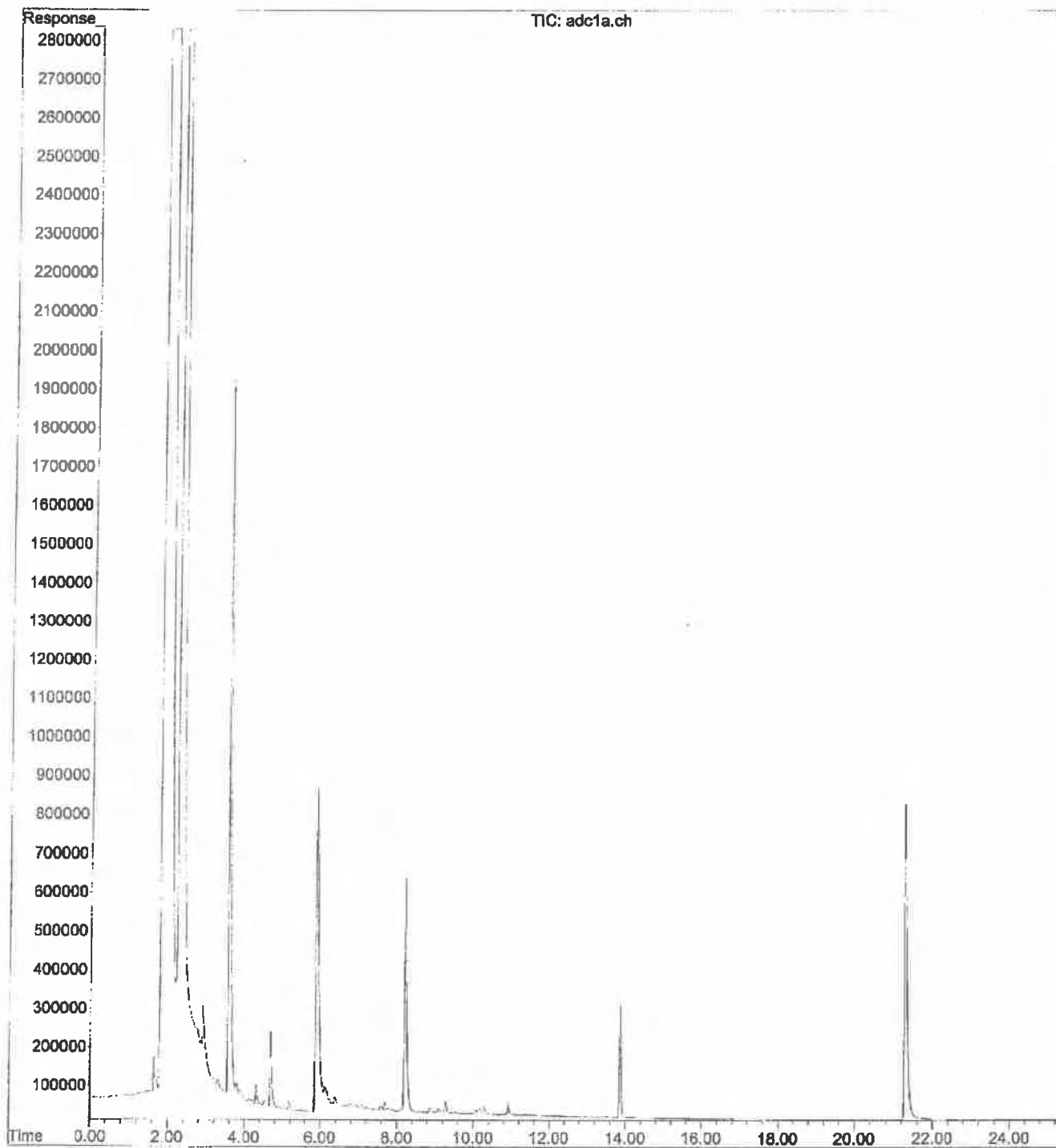
Sample Spike Data (MS)

| Analyte | Sample ID # | Sample Result (mg/L) | Spike Sample Result (mg/L) | Spike Amount (mg/L) | %Recovery Control Limits (75-125%) |
|----------------|--------------------|-------------------------------------|---|------------------------------------|---|
| Silver | 10171-04MS | 0 | 0.24 | 0.3 | 97 |
| Arsenic | 10171-04MS | 0.0032 | 0.53 | 0.5 | 105 |
| Barium | 10171-04MS | 0.1773 | 0.67 | 0.5 | 99 |
| Cadmium | 10171-04MS | 0 | 0.51 | 0.5 | 103 |
| Chromium | 10171-04MS | 0 | 0.50 | 0.5 | 101 |
| Lead | 10171-04MS | 0.0027 | 0.47 | 0.5 | 94 |
| Selenium | 10171-04MS | 0.0152 | 0.53 | 0.5 | 103 |

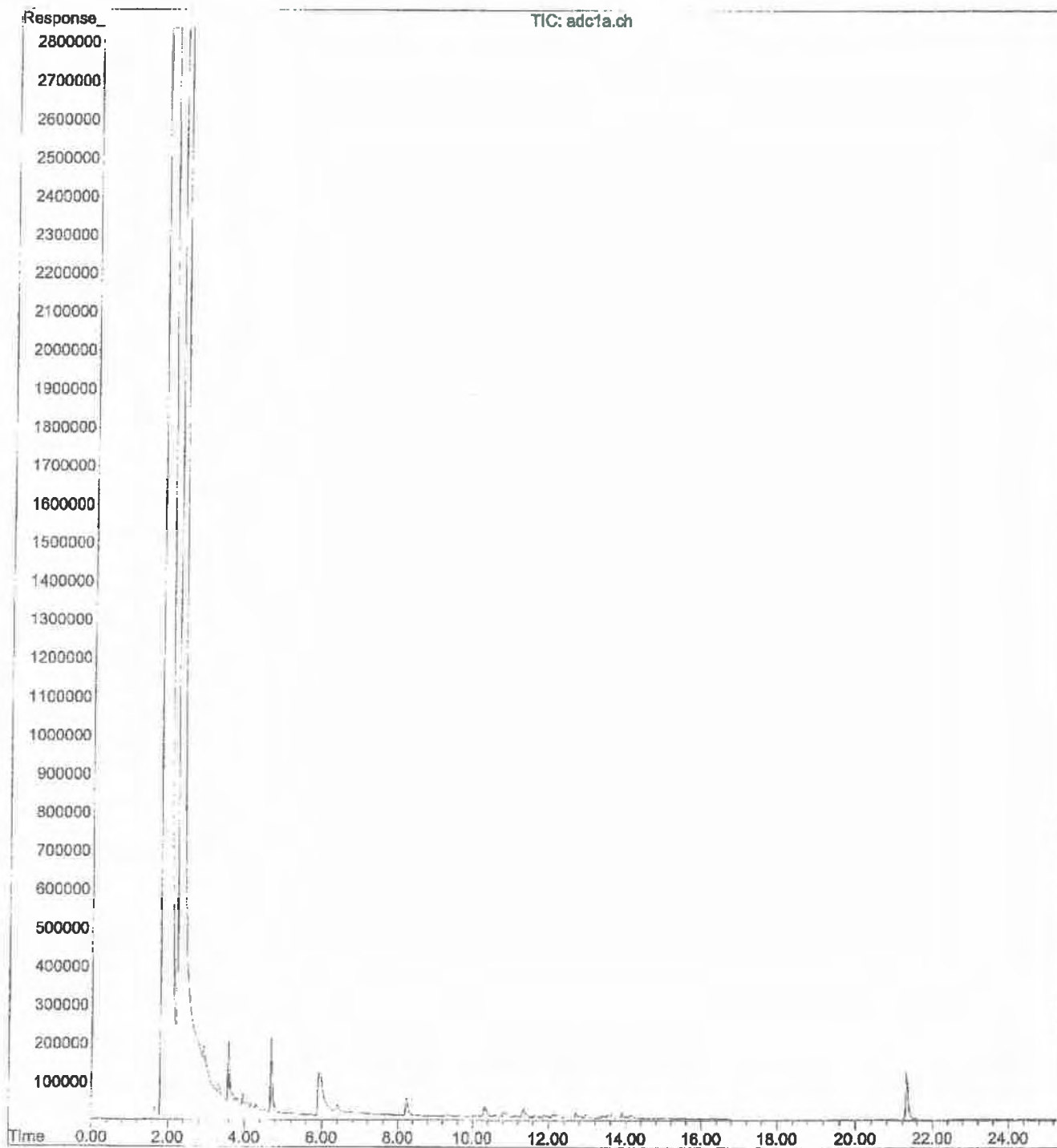
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Operator : AJD
Acquired : 18 Apr 2006 12:46 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10171-01 pcb x1
Misc Info :
Vial Number: 10



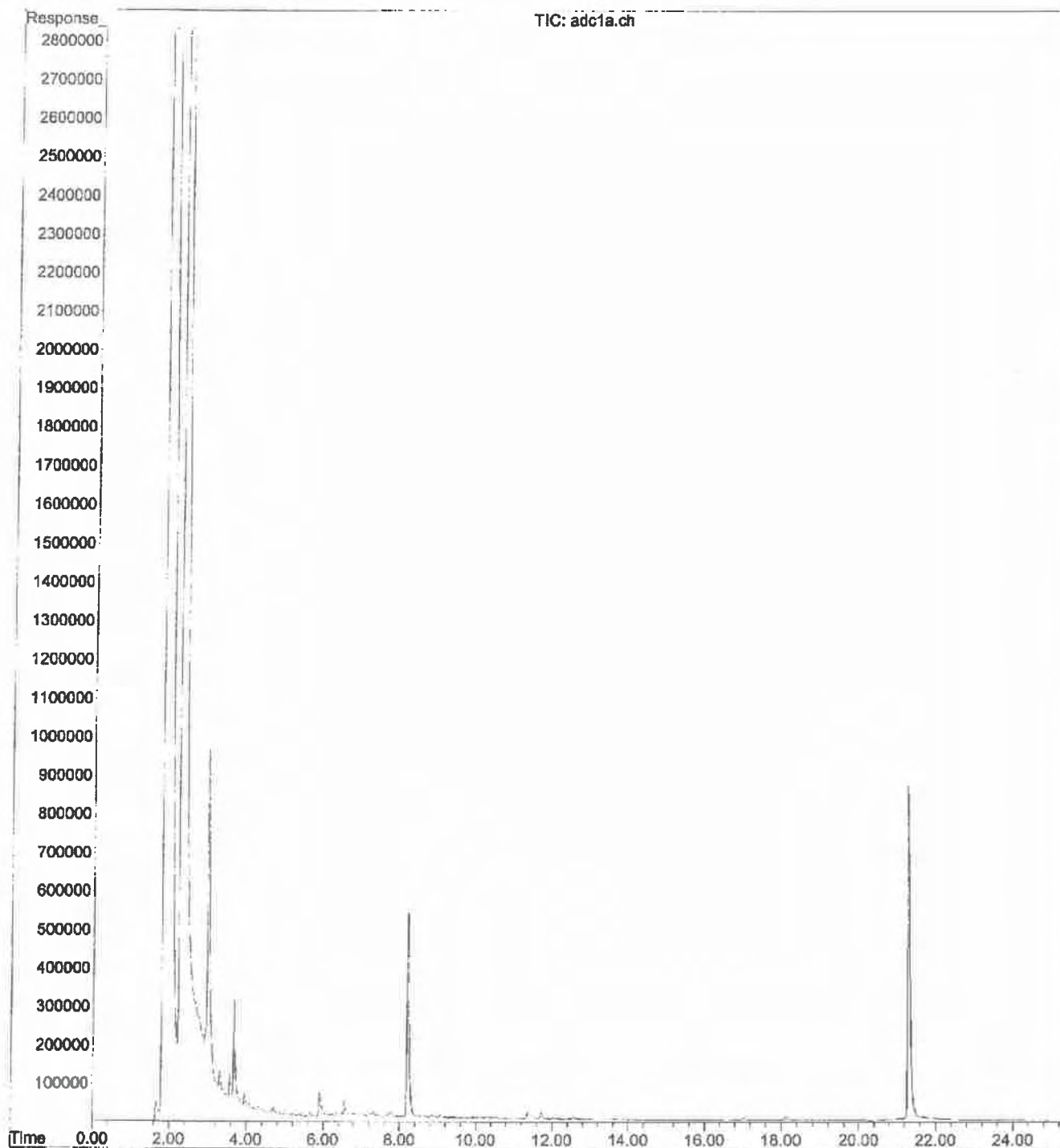
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Operator : AJD
Acquired : 18 Apr 2006 06:59 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10171-02 pcb x1
Misc Info :
Vial Number: 23



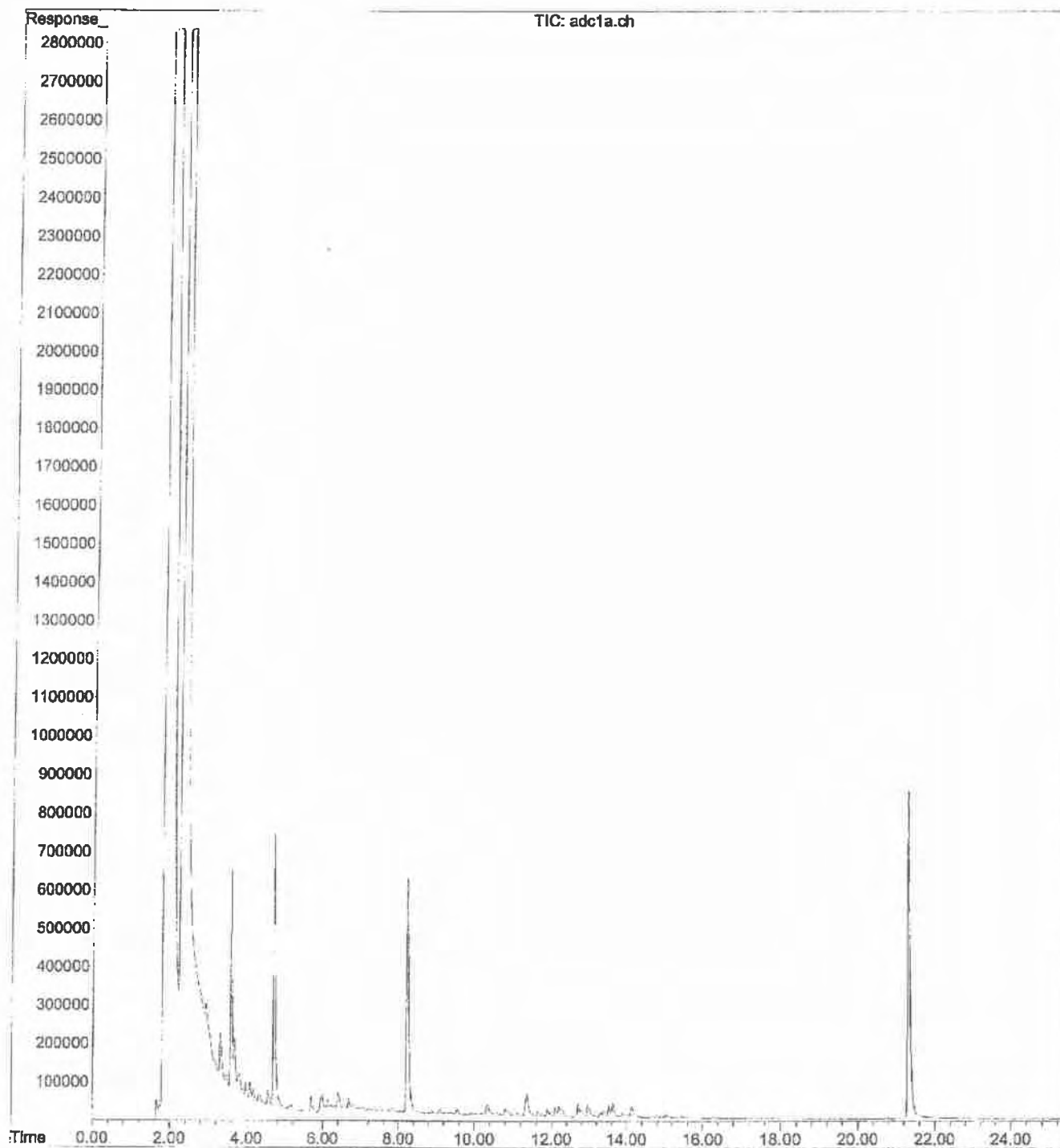
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Operator : AJD
Acquired : 19 Apr 2006 06:37 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10171-04rrpcbx20
Misc Info :
Vial Number: 11



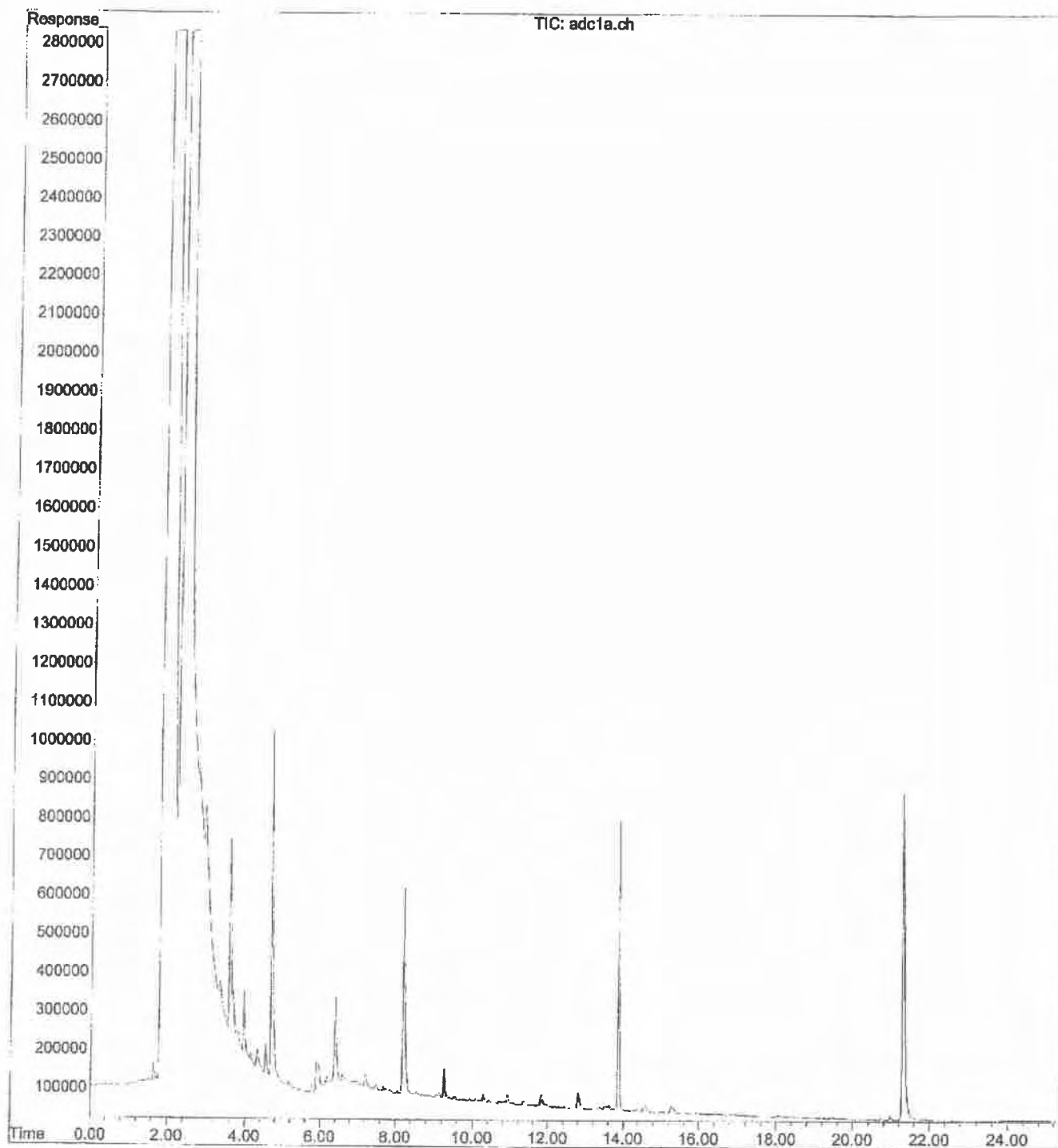
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Operator : AJD
Acquired : 18 Apr 2006 01:14 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10171-05 pcb x1
Misc Info :
Vial Number: 11



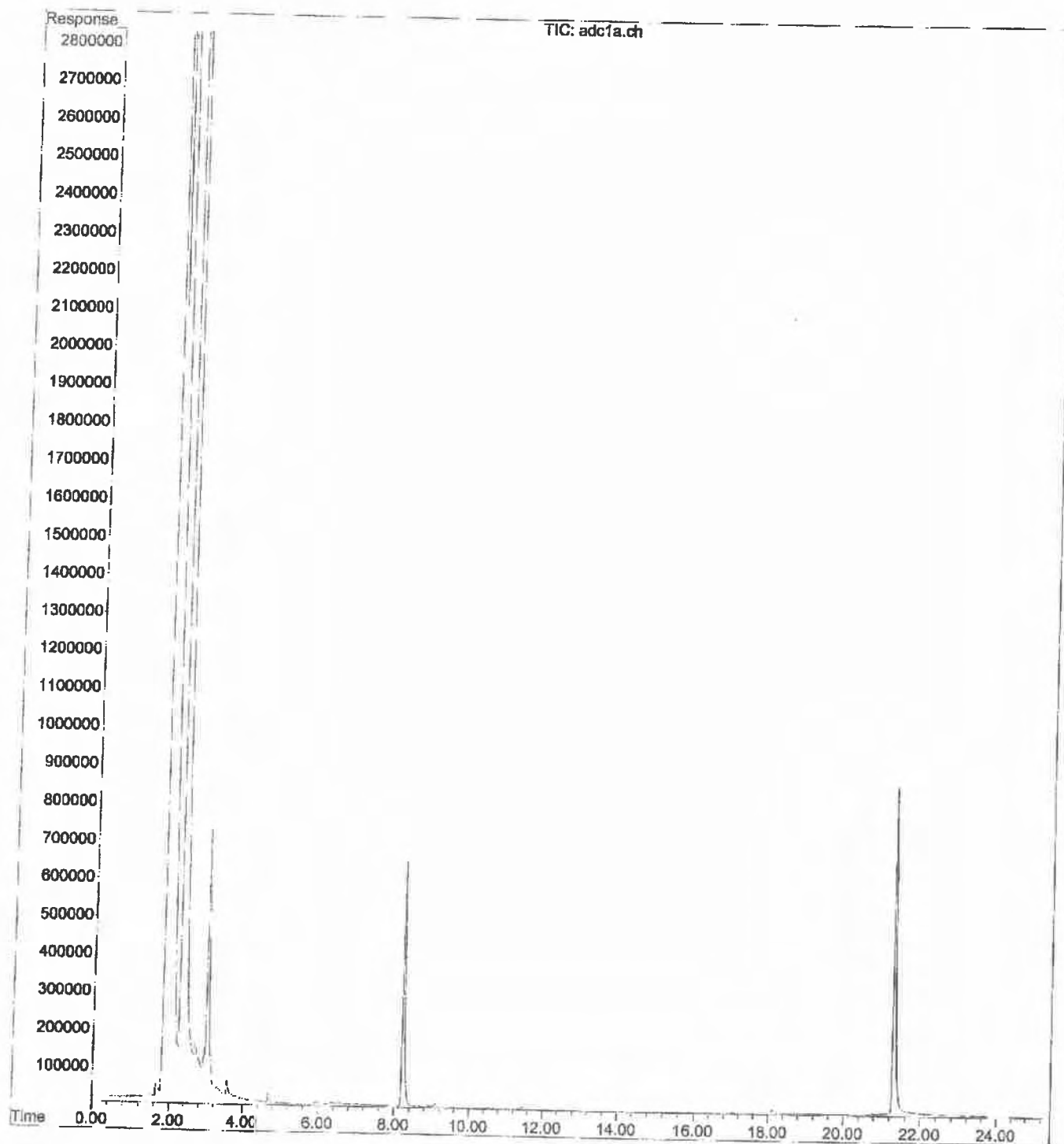
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Operator : AJD
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Instrument : SVOA03
Sample Name: 10171-06rr pcbx1
Misc Info :
Vial Number: 6



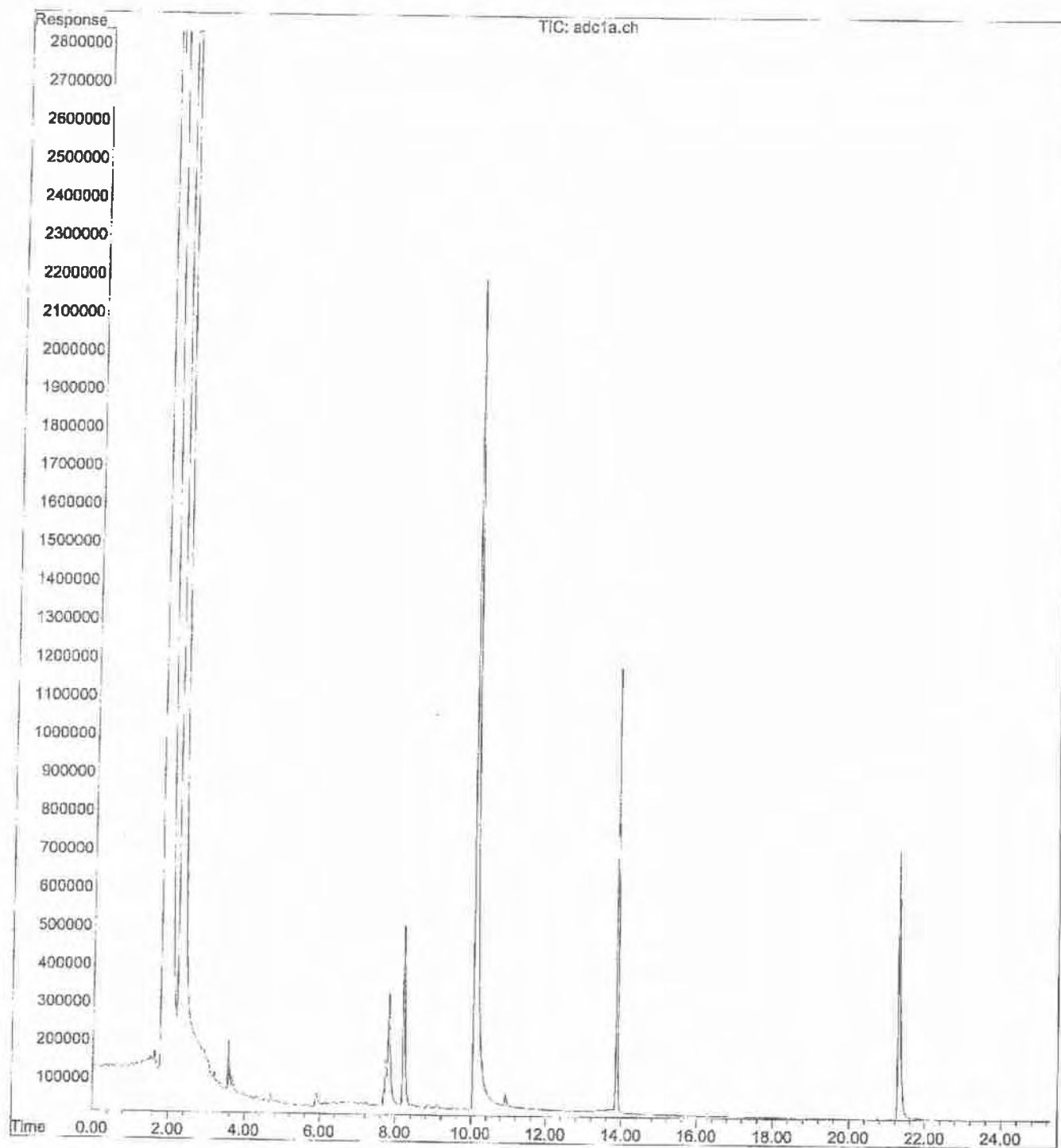
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Operator : AJD
Acquired : 18 Apr 2006 06:02 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10171-07 pcb x1
Misc Info :
Vial Number: 21



File : X:\DATA\SVOA03\2006\APR06\041806\012F0101.D
Operator : AJD
Acquired : 18 Apr 2006 01:43 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10171-08 pcb x1
Misc Info :
Vial Number: 12



File : X:\DATA\SVOA03\2006\APR06\041806\022F0101.D
Operator : AJD
Acquired : 18 Apr 2006 06:31 pm using AcqMethod S3PP10B.MTH
Instrument : SVOA03
Sample Name: 10171-09 pcb x1
Misc Info :
Vial Number: 22



Resource Laboratories, LLC

124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

10171

PAGE 1 OF 1

Company Name:

Phone #: 603-285-9700

Company Address:

FAX #: 603-430-9457

Project Manager:

Site Location (City, State):

154 West Main St, Portsmouth, NH

14702 Park, MA

Project Manager:

Project ID / Name:

Invoice To:

Protocol: RCRA SDWA NPDES
MCP NHDES OTHER

ES-M

2006-05-12 / Leach Chemical

| Lab Sample ID (Lab Use Only) | Field ID | # CONTAINERS | Matrix | | | Preservation Method | | | | | | Sampling | | |
|---------------------------------|-----------|--------------|--------|-------|-------|---------------------|------------------|--------------------------------|------|------|-----------------|----------|------|---------|
| | | | WATER | SOLID | OTHER | HCl | HNO ₃ | H ₂ SO ₄ | NaOH | MeOH | OTHER (Specify) | DATE | TIME | SAMPLER |
| 10171-01 | ES-M-04 | 4 | Y | | | 3 | | | | | | 4/10/06 | 0830 | LM |
| 02 | ES-M-04 | 4 | X | | | 3 | | | | | | | 0900 | LM |
| 03 | DUP-2 | 3 | X | | | 3 | | | | | | | 0915 | LM |
| 04 | ES-M-05 | 5 | X | | | 3 | 1 | | | | | | 0930 | LM |
| 05 | ES-M-05B | 5 | X | | | 3 | 1 | | | | | | 1000 | LM |
| 06 | ES-M-06 | 4 | X | | | 3 | | | | | | | 1040 | LM |
| 07 | ES-M-15 | 4 | X | | | 3 | | | | | | | 1115 | LM |
| 08 | ES-M-07 | 4 | X | | | 3 | | | | | | | 1145 | LM |
| 09 | ES-M-10 | 8 | Y | | | 6 | | | | | | | 1230 | LM |
| 10 | Imp Blade | 1 | X | | | 1 | | | | | | | | |

| | | |
|--|--|--|
| <input type="checkbox"/> VOC B260-NH List | <input checked="" type="checkbox"/> MADEP VPH | <input type="checkbox"/> MEGRO |
| <input checked="" type="checkbox"/> VOC B260 | <input type="checkbox"/> VOC B015GRO | <input type="checkbox"/> VOC B24 (8260R) |
| <input type="checkbox"/> VOC B260 BTEX, MIBE, Naphthalene only | | |
| <input type="checkbox"/> VOC 524.2 | <input type="checkbox"/> VOC 524.2 NH List | |
| <input type="checkbox"/> TPH Fingerprint | <input type="checkbox"/> MEDRO | <input checked="" type="checkbox"/> DRO 8015 |
| <input type="checkbox"/> 8270PAH | <input type="checkbox"/> 8270ABN | <input type="checkbox"/> 625 |
| <input checked="" type="checkbox"/> 8082 PCB | <input type="checkbox"/> 8081 Pesticides | <input type="checkbox"/> 606 |
| <input type="checkbox"/> O&G 1064 | <input type="checkbox"/> O&G SM5520F | |
| <input type="checkbox"/> pH | <input type="checkbox"/> BOD | <input type="checkbox"/> Conductivity |
| <input type="checkbox"/> TSS | <input type="checkbox"/> TDS | <input type="checkbox"/> TS |
| <input checked="" type="checkbox"/> RCRA Metals | <input type="checkbox"/> Priority Pollutant Metals | <input type="checkbox"/> TAL Metals |
| <input type="checkbox"/> Total Metals-list | <input type="checkbox"/> Dissolved Metals-list | |
| <input type="checkbox"/> Ammonia | <input type="checkbox"/> COD | |
| <input type="checkbox"/> T-Phosphate | <input type="checkbox"/> Phenol | |
| <input type="checkbox"/> Cyanide | <input type="checkbox"/> Sulfide | |
| <input type="checkbox"/> Nitrate | <input type="checkbox"/> Nitrite | <input type="checkbox"/> Ortho P |
| <input type="checkbox"/> Sulfate | <input type="checkbox"/> Bromide | <input type="checkbox"/> Chloride |
| <input type="checkbox"/> Corrosivity | <input type="checkbox"/> Reactive CN | <input type="checkbox"/> Reactive S- |
| <input type="checkbox"/> Ignitibility/FP | | |
| <input type="checkbox"/> TCLP Metals | <input type="checkbox"/> TCLP VOC | <input type="checkbox"/> TCLP SVOC |
| <input type="checkbox"/> TCLP Pesticide | <input type="checkbox"/> TCLP Herbicides (subcontract) | |
| <input type="checkbox"/> Standard Drinking Water Test | <input type="checkbox"/> Bacteria P/A | |

Grab (G) or Composite (C)

TEST REQUESTED

SPECIAL INSTRUCTIONS

Priority (24 hr) ☐

Expected (48 hr) ☐

10 Business Days ☒

Other ☐

E-Mail Address: jallen@resource.com

Queue #

PO #

REPORTING INSTRUCTIONS

☐ FAX ☐ OTHER (specify)

☐ PDF ☐ Excel Spreadsheet

RECEIVED ON ICE

YES ☒ NO ☐

TEMPERATURE

64 °C

Lab Use Only

CUSTODY RECORD

Relinquished by: [Signature]

Date: 4/10/06 Time: 1145

Received by: [Signature]

Date: 4/11/06 Time: 1145

Date: 4/11/06 Time: 1145

Relinquished by:

Date: Time:

Received by Laboratory:

Date: Time:

Way Bill #:

Laboratory Report

Joe Callahan
Environmental Strategies & Management
184 West Main Street
Norton, MA 02766

PO Number: None
LabID: 11371
Date Received: 11/16/06

Project: 2006-056 DND Lewis Chem

Attached please find results for the analysis of the samples received on the date referenced above.

Enclosed please find the revised report. The report now includes the MADEP MCP Report Certification Form.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Resource Laboratories, LLC Quality Assurance Plan. The Standard Operating Procedures (SOP) are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Resource Laboratories, LLC maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Resource Laboratories, LLC



Susan Sylvester
Principal, General Manager

12-7-06

Date

Total number of pages 21

Resource Laboratories, LLC Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Lab ID: 11371

Lab Number: 11371-001

Sample ID: PZ-04

Matrix: Water

Sampled: 11/15/06 11:30

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| dichlorodifluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| chloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| vinyl chloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| bromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| chloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| trichlorofluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| diethyl ether | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| acetone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,1-dichloroethene | < 1 | 1 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| methylene chloride | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| carbon disulfide | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| methyl t-butyl ether (MTBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| trans-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| isopropyl ether (DIPE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| ethyl t-butyl ether (ETBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,1-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| t-butanol (TBA) | < 40 | 40 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 2-butanone (MEK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 2,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| cis-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| chloroform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| bromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| tetrahydrofuran (THF) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,1,1-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,1-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| t-amyl-methyl ether (TAME) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| carbon tetrachloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,2-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| benzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| trichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| bromodichloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,4-dioxane | < 50 | 50 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| dibromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 4-methyl-2-pentanone (MIBK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| cis-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| toluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| trans-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 2-hexanone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,1,2-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,3-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| tetrachloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| dibromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-001

Sample ID: PZ-04

Matrix: Water

Sampled: 11/15/06 11:30

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|---------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| 1,2-dibromoethane (EDB) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| chlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,1,1,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| ethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| m&p-xylenes | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| o-xylene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| styrene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| bromoform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| isopropylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,1,2,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,2,3-trichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| n-propylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| bromobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,3,5-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 2-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 4-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| tert-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,2,4-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| sec-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,3-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 4-isopropyltoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,4-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,2-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| n-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,2-dibromo-3-chloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,2,4-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| hexachlorobutadiene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| naphthalene | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 1,2,3-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| Surrogate Recovery: | | Limits | | | | | | | |
| dibromofluoromethane SUR | 98 | 78-114 | % | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| toluene-D8 SUR | 102 | 88-110 | % | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |
| 4-bromofluorobenzene SUR | 93 | 86-115 | % | 1 | LMM | | 11/22/06 | 5:10 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-002

Sample ID: PZ-05

Matrix: Water

Sampled: 11/15/06 12:20

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| dichlorodifluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| chloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| vinyl chloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| bromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| chloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| trichlorofluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| diethyl ether | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| acetone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,1-dichloroethene | < 1 | 1 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| methylene chloride | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| carbon disulfide | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| methyl t-butyl ether (MTBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| trans-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| isopropyl ether (DIPE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| ethyl t-butyl ether (ETBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,1-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| t-butanol (TBA) | < 40 | 40 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 2-butanone (MEK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 2,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| cis-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| chloroform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| bromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| tetrahydrofuran (THF) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,1,1-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,1-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| t-amyl-methyl ether (TAME) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| carbon tetrachloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,2-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| benzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| trichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| bromodichloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,4-dioxane | < 50 | 50 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| dibromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 4-methyl-2-pentanone (MIBK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| cis-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| toluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| trans-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 2-hexanone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,1,2-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,3-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| tetrachloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| dibromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-002

Sample ID: PZ-05

Matrix: Water

Sampled: 11/15/06 12:20

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|---------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| 1,2-dibromoethane (EDB) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| chlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,1,1,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| ethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| m&p-xylenes | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| o-xylene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| styrene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| bromoform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| isopropylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,1,2,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,2,3-trichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| n-propylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| bromobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,3,5-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 2-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 4-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| tert-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,2,4-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| sec-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,3-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 4-isopropyltoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,4-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,2-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| n-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,2-dibromo-3-chloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,2,4-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| hexachlorobutadiene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| naphthalene | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 1,2,3-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| Surrogate Recovery: | | Limits | | | | | | | |
| dibromofluoromethane SUR | 98 | 78-114 | % | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| toluene-D8 SUR | 101 | 88-110 | % | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |
| 4-bromofluorobenzene SUR | 91 | 86-115 | % | 1 | LMM | | 11/22/06 | 5:41 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-003

Sample ID: PZ-06

Matrix: Water

Sampled: 11/15/06 13:00

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| dichlorodifluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| chloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| vinyl chloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| bromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| chloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| trichlorofluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| diethyl ether | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| acetone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,1-dichloroethene | < 1 | 1 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| methylene chloride | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| carbon disulfide | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| methyl t-butyl ether (MTBE) | 74 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| trans-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| isopropyl ether (DIPE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| ethyl t-butyl ether (ETBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,1-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| t-butanol (TBA) | < 40 | 40 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 2-butanone (MEK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 2,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| cis-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| chloroform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| bromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| tetrahydrofuran (THF) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,1,1-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,1-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| t-amyl-methyl ether (TAME) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| carbon tetrachloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,2-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| benzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| trichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| bromodichloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,4-dioxane | < 50 | 50 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| dibromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 4-methyl-2-pentanone (MIBK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| cis-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| toluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| trans-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 2-hexanone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,1,2-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,3-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| tetrachloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| dibromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-003

Sample ID: PZ-06

Matrix: Water

Sampled: 11/15/06 13:00

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|---------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| 1,2-dibromoethane (EDB) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| chlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,1,1,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| ethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| m&p-xylenes | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| o-xylene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| styrene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| bromoform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| isopropylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,1,2,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,2,3-trichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| n-propylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| bromobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,3,5-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 2-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 4-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| tert-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,2,4-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| sec-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,3-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 4-isopropyltoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,4-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,2-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| n-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,2-dibromo-3-chloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,2,4-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| hexachlorobutadiene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| naphthalene | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 1,2,3-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| Surrogate Recovery: | | Limits | | | | | | | |
| dibromofluoromethane SUR | 98 | 78-114 | % | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| toluene-D8 SUR | 101 | 88-110 | % | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |
| 4-bromofluorobenzene SUR | 97 | 86-115 | % | 1 | LMM | | 11/22/06 | 6:11 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-004

Sample ID: PZ-07

Matrix: Water

Sampled: 11/15/06 15:00

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| dichlorodifluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| chloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| vinyl chloride | 4 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| bromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| chloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| trichlorofluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| diethyl ether | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| acetone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,1-dichloroethene | < 1 | 1 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| methylene chloride | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| carbon disulfide | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| methyl t-butyl ether (MTBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| trans-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| isopropyl ether (DIPE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| ethyl t-butyl ether (ETBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,1-dichloroethane | 10 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| t-butanol (TBA) | < 40 | 40 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 2-butanone (MEK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 2,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| cis-1,2-dichloroethene | 15 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| chloroform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| bromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| tetrahydrofuran (THF) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,1,1-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,1-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| t-amyl-methyl ether (TAME) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| carbon tetrachloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,2-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| benzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| trichloroethene | 11 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| bromodichloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,4-dioxane | < 50 | 50 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| dibromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 4-methyl-2-pentanone (MIBK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| cis-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| toluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| trans-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 2-hexanone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,1,2-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,3-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| tetrachloroethene | 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| dibromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-004

Sample ID: PZ-07

Matrix: Water

Sampled: 11/15/06 15:00

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|---------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| 1,2-dibromoethane (EDB) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| chlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,1,1,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| ethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| m&p-xylenes | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| o-xylene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| styrene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| bromoform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| isopropylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,1,2,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,2,3-trichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| n-propylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| bromobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,3,5-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 2-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 4-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| tert-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,2,4-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| sec-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,3-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 4-isopropyltoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,4-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,2-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| n-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,2-dibromo-3-chloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,2,4-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| hexachlorobutadiene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| naphthalene | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 1,2,3-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| Surrogate Recovery: | | Limits | | | | | | | |
| dibromofluoromethane SUR | 99 | 78-114 | % | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| toluene-D8 SUR | 103 | 88-110 | % | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |
| 4-bromofluorobenzene SUR | 95 | 86-115 | % | 1 | LMM | | 11/22/06 | 6:42 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-005

Sample ID: Trip Blank

Matrix: Water

Sampled: 11/15/06

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|-------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| dichlorodifluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| chloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| vinyl chloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| bromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| chloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| trichlorofluoromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| diethyl ether | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| acetone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,1-dichloroethene | < 1 | 1 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| methylene chloride | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| carbon disulfide | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| methyl t-butyl ether (MTBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| trans-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| isopropyl ether (DIPE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| ethyl t-butyl ether (ETBE) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,1-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| t-butanol (TBA) | < 40 | 40 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 2-butanone (MEK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 2,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| cis-1,2-dichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| chloroform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| bromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| tetrahydrofuran (THF) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,1,1-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,1-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| t-amyl-methyl ether (TAME) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| carbon tetrachloride | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,2-dichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| benzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| trichloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,2-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| bromodichloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,4-dioxane | < 50 | 50 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| dibromomethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 4-methyl-2-pentanone (MIBK) | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| cis-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| toluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| trans-1,3-dichloropropene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 2-hexanone | < 10 | 10 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,1,2-trichloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,3-dichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| tetrachloroethene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| dibromochloromethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |

Lab ID: 11371

Lab Number: 11371-005

Sample ID: Trip Blank

Matrix: Water

Sampled: 11/15/06

| Parameter: | Result | Quant Limit | Units | Instr Dil'n Factor | Analyst | Prep Date | Analysis Date | Analysis Time | Reference |
|-----------------------------|--------|---------------|-------|--------------------|---------|-----------|---------------|---------------|--------------|
| 1,2-dibromoethane (EDB) | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| chlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,1,1,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| ethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| m&p-xylenes | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| o-xylene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| styrene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| bromoform | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| isopropylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,1,2,2-tetrachloroethane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,2,3-trichloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| n-propylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| bromobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,3,5-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 2-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 4-chlorotoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| tert-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,2,4-trimethylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| sec-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,3-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 4-isopropyltoluene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,4-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,2-dichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| n-butylbenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,2-dibromo-3-chloropropane | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,2,4-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| hexachlorobutadiene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| naphthalene | < 5 | 5 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 1,2,3-trichlorobenzene | < 2 | 2 | ug/L | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| Surrogate Recovery: | | Limits | | | | | | | |
| dibromofluoromethane SUR | 98 | 78-114 | % | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| toluene-D8 SUR | 104 | 88-110 | % | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |
| 4-bromofluorobenzene SUR | 93 | 86-115 | % | 1 | LMM | | 11/22/06 | 3:08 | SW5030B8260B |

Quality Control Report

RESOURCE LABORATORIES, LLC.

Case Narrative

Lab # 11371

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 4 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

VOC 8260: LCS 0403790 did not meet acceptance criteria for bromoform and 1,2 dibromo-3-chloropropane. LCSD 0403790 did not meet acceptance criteria for bromoform and 1,2 dibromo-3-chloropropane. These compounds are known to be problematic in the method.

No other exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

MADEP MCP Analytical Method Report Certification Form

Laboratory Name: Resource Laboratories, LLC

Lab # 11371

Project Location Hyde Park

Project #

MADEP RTN (if available)

This form provides certifications for the following data set in the Lab # referenced above (see Chain of Custody for samples numbers)

Sample Matrices: Groundwater (x) Soil/Sediment () Drinking Water () Other:

| | | | | | |
|--------------------------------|----------|----------|---------------|-------------|-----------|
| MCP SW- 846 Methods Used | 8260 (x) | 8081 () | 6010 () | Cyanide () | Other () |
| | 8270 () | VPH () | 7470/7471 () | Other () | Other () |
| | 8082 () | EPH () | Other () | Other () | Other () |

| | | |
|---|--|-------------------|
| A | Were all samples received by the laboratory in a condition consistent with that described on the Chain of Custody documentation for the data set? | Yes (x) No () |
| B | Were all QA/QC procedures required for the specified analytical method(s) included in this report followed, including the requirement to note and discuss in a narrative QC data that did not meet appropriate performance standards or guidelines? | Yes (x) No () |
| C | Does the data included in this report meet all the analytical requirements for "Presumptive Certainty" as described in Section 2.0 (a), (b), (c) and (d) of the MADEP document CAM VII A, "Quality Assurance and Quality Control Guidelines for the Acquisition and Reporting of Analytical Data?" | Yes (x) No () |
| D | VPH and EPH methods only: Was the VPH or EPH method run without significant modifications? (see Section 11.3 of respective Methods) | Yes () No () NA |
| E | Were all QC performance standards and recommendations for the specified methods achieved? | Yes () No (x) |
| F | Were results for all analyte-list compounds/elements for the specified method(s) reported? | Yes (x) No () |

I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.

Signature: Susan C. SylvesterPosition: Lab DirectorPrinted Name: Susan C. SylvesterDate: 12-6-06

- QC Report -

| Method | QC ID | Parameter | Associated Sample | Result | Units | Amt Added | %R | Limit | RPD | RPD Limit |
|--------------|------------|-----------------------------|-------------------|--------|---------|-----------|----|-------|-----|-----------|
| SW5030B8260B | BLK0403790 | dichlorodifluoromethane | | < | 2 ug/L | | | | | |
| | | chloromethane | | < | 2 ug/L | | | | | |
| | | vinyl chloride | | < | 2 ug/L | | | | | |
| | | bromomethane | | < | 2 ug/L | | | | | |
| | | chloroethane | | < | 2 ug/L | | | | | |
| | | trichlorofluoromethane | | < | 2 ug/L | | | | | |
| | | diethyl ether | | < | 10 ug/L | | | | | |
| | | acetone | | < | 10 ug/L | | | | | |
| | | 1,1-dichloroethene | | < | 1 ug/L | | | | | |
| | | methylene chloride | | < | 5 ug/L | | | | | |
| | | carbon disulfide | | < | 2 ug/L | | | | | |
| | | methyl t-butyl ether (MTBE) | | < | 2 ug/L | | | | | |
| | | trans-1,2-dichloroethene | | < | 2 ug/L | | | | | |
| | | isopropyl ether (DIPE) | | < | 2 ug/L | | | | | |
| | | ethyl t-butyl ether (ETBE) | | < | 2 ug/L | | | | | |
| | | 1,1-dichloroethane | | < | 2 ug/L | | | | | |
| | | t-butanol (TBA) | | < | 40 ug/L | | | | | |
| | | 2-butanone (MEK) | | < | 10 ug/L | | | | | |
| | | 2,2-dichloropropane | | < | 2 ug/L | | | | | |
| | | cis-1,2-dichloroethene | | < | 2 ug/L | | | | | |
| | | chloroform | | < | 2 ug/L | | | | | |
| | | bromochloromethane | | < | 2 ug/L | | | | | |
| | | tetrahydrofuran (THF) | | < | 10 ug/L | | | | | |
| | | 1,1,1-trichloroethane | | < | 2 ug/L | | | | | |
| | | 1,1-dichloropropene | | < | 2 ug/L | | | | | |
| | | t-amyl-methyl ether (TAME) | | < | 2 ug/L | | | | | |
| | | carbon tetrachloride | | < | 2 ug/L | | | | | |
| | | 1,2-dichloroethane | | < | 2 ug/L | | | | | |
| | | benzene | | < | 2 ug/L | | | | | |
| | | trichloroethene | | < | 2 ug/L | | | | | |
| | | 1,2-dichloropropane | | < | 2 ug/L | | | | | |
| | | bromodichloromethane | | < | 2 ug/L | | | | | |
| | | 1,4-dioxane | | < | 50 ug/L | | | | | |
| | | dibromomethane | | < | 2 ug/L | | | | | |
| | | 4-methyl-2-pentanone (MIBK) | | < | 10 ug/L | | | | | |
| | | cis-1,3-dichloropropene | | < | 2 ug/L | | | | | |
| | | toluene | | < | 2 ug/L | | | | | |
| | | trans-1,3-dichloropropene | | < | 2 ug/L | | | | | |
| | | 2-hexanone | | < | 10 ug/L | | | | | |
| | | 1,1,2-trichloroethane | | < | 2 ug/L | | | | | |
| | | 1,3-dichloropropane | | < | 2 ug/L | | | | | |
| | | tetrachloroethene | | < | 2 ug/L | | | | | |
| | | dibromochloromethane | | < | 2 ug/L | | | | | |
| | | 1,2-dibromoethane (EDB) | | < | 2 ug/L | | | | | |
| | | chlorobenzene | | < | 2 ug/L | | | | | |
| | | 1,1,1,2-tetrachloroethane | | < | 2 ug/L | | | | | |
| | | ethylbenzene | | < | 2 ug/L | | | | | |
| | | m&p-xylenes | | < | 2 ug/L | | | | | |
| | | o-xylene | | < | 2 ug/L | | | | | |
| | | styrene | | < | 2 ug/L | | | | | |
| | | bromoform | | < | 2 ug/L | | | | | |

| Method | QC ID | Parameter | Associated Sample | Result | Units | Amt Added | %R | Limit | RPD | RPD Limit |
|--------------|------------|-----------------------------|-------------------|--------|--------|-----------|----|-------|-----|-----------|
| SW5030B8260B | BLK0403790 | isopropylbenzene | | < | 2 ug/L | | | | | |
| | | 1,1,2,2-tetrachloroethane | | < | 2 ug/L | | | | | |
| | | 1,2,3-trichloropropane | | < | 2 ug/L | | | | | |
| | | n-propylbenzene | | < | 2 ug/L | | | | | |
| | | bromobenzene | | < | 2 ug/L | | | | | |
| | | 1,3,5-trimethylbenzene | | < | 2 ug/L | | | | | |
| | | 2-chlorotoluene | | < | 2 ug/L | | | | | |
| | | 4-chlorotoluene | | < | 2 ug/L | | | | | |
| | | tert-butylbenzene | | < | 2 ug/L | | | | | |
| | | 1,2,4-trimethylbenzene | | < | 2 ug/L | | | | | |
| | | sec-butylbenzene | | < | 2 ug/L | | | | | |
| | | 1,3-dichlorobenzene | | < | 2 ug/L | | | | | |
| | | 4-isopropyltoluene | | < | 2 ug/L | | | | | |
| | | 1,4-dichlorobenzene | | < | 2 ug/L | | | | | |
| | | 1,2-dichlorobenzene | | < | 2 ug/L | | | | | |
| | | n-butylbenzene | | < | 2 ug/L | | | | | |
| | | 1,2-dibromo-3-chloropropane | | < | 2 ug/L | | | | | |
| | | 1,2,4-trichlorobenzene | | < | 2 ug/L | | | | | |
| | | hexachlorobutadiene | | < | 2 ug/L | | | | | |
| | | naphthalene | | < | 5 ug/L | | | | | |
| | | 1,2,3-trichlorobenzene | | < | 2 ug/L | | | | | |
| | | dibromofluoromethane SUR | | 100 | % | | | 78 | 114 | |
| | | toluene-D8 SUR | | 100 | % | | | 88 | 110 | |
| | | 4-bromofluorobenzene SUR | | 92 | % | | | 86 | 115 | |

| Method | QC ID | Parameter | Associated Sample | Result | Units | Amt Added | %R | Limit | RPD | RPD Limit |
|--------------|------------|-----------------------------|-------------------|---------|-------|-----------|-----|-------|-----|-----------|
| SW5030B8260B | LCS0403790 | dichlorodifluoromethane | | 17 ug/L | | 20 | 86 | 70 | 130 | |
| | | chloromethane | | 21 ug/L | | 20 | 103 | 70 | 130 | |
| | | vinyl chloride | | 19 ug/L | | 20 | 95 | 70 | 130 | |
| | | bromomethane | | 17 ug/L | | 20 | 87 | 70 | 130 | |
| | | chloroethane | | 20 ug/L | | 20 | 101 | 70 | 130 | |
| | | trichlorofluoromethane | | 18 ug/L | | 20 | 92 | 70 | 130 | |
| | | diethyl ether | | 22 ug/L | | 20 | 110 | 70 | 130 | |
| | | acetone | | 21 ug/L | | 20 | 104 | 70 | 130 | |
| | | 1,1-dichloroethene | | 20 ug/L | | 20 | 102 | 70 | 130 | |
| | | methylene chloride | | 22 ug/L | | 20 | 111 | 70 | 130 | |
| | | carbon disulfide | | 17 ug/L | | 20 | 83 | 70 | 130 | |
| | | methyl t-butyl ether (MTBE) | | 21 ug/L | | 20 | 103 | 70 | 130 | |
| | | trans-1,2-dichloroethene | | 21 ug/L | | 20 | 106 | 70 | 130 | |
| | | isopropyl ether (DIPE) | | 20 ug/L | | 20 | 99 | 70 | 130 | |
| | | ethyl t-butyl ether (ETBE) | | 21 ug/L | | 20 | 105 | 70 | 130 | |
| | | 1,1-dichloroethane | | 21 ug/L | | 20 | 103 | 70 | 130 | |
| | | t-butanol (TBA) | | 91 ug/L | | 100 | 91 | 70 | 130 | |
| | | 2-butanone (MEK) | | 17 ug/L | | 20 | 87 | 70 | 130 | |
| | | 2,2-dichloropropane | | 15 ug/L | | 20 | 75 | 70 | 130 | |
| | | cis-1,2-dichloroethene | | 21 ug/L | | 20 | 107 | 70 | 130 | |
| | | chloroform | | 21 ug/L | | 20 | 107 | 70 | 130 | |
| | | bromochloromethane | | 22 ug/L | | 20 | 108 | 70 | 130 | |
| | | tetrahydrofuran (THF) | | 18 ug/L | | 20 | 91 | 70 | 130 | |
| | | 1,1,1-trichloroethane | | 20 ug/L | | 20 | 100 | 70 | 130 | |
| | | 1,1-dichloropropene | | 21 ug/L | | 20 | 107 | 70 | 130 | |
| | | t-amyl-methyl ether (TAME) | | 21 ug/L | | 20 | 104 | 70 | 130 | |
| | | carbon tetrachloride | | 17 ug/L | | 20 | 86 | 70 | 130 | |
| | | 1,2-dichloroethane | | 21 ug/L | | 20 | 107 | 70 | 130 | |
| | | benzene | | 21 ug/L | | 20 | 105 | 70 | 130 | |
| | | trichloroethene | | 21 ug/L | | 20 | 106 | 70 | 130 | |
| | | 1,2-dichloropropane | | 21 ug/L | | 20 | 105 | 70 | 130 | |
| | | bromodichloromethane | | 18 ug/L | | 20 | 89 | 70 | 130 | |
| | | 1,4-dioxane | < | 50 ug/L | | 40 | 101 | | | |
| | | dibromomethane | | 21 ug/L | | 20 | 105 | 70 | 130 | |
| | | 4-methyl-2-pentanone (MIBK) | | 18 ug/L | | 20 | 92 | 70 | 130 | |
| | | cis-1,3-dichloropropene | | 18 ug/L | | 20 | 92 | 70 | 130 | |
| | | toluene | | 22 ug/L | | 20 | 111 | 70 | 130 | |
| | | trans-1,3-dichloropropene | | 18 ug/L | | 20 | 89 | 70 | 130 | |
| | | 2-hexanone | | 17 ug/L | | 20 | 85 | 70 | 130 | |
| | | 1,1,2-trichloroethane | | 21 ug/L | | 20 | 106 | 70 | 130 | |
| | | 1,3-dichloropropane | | 21 ug/L | | 20 | 103 | 70 | 130 | |
| | | tetrachloroethene | | 21 ug/L | | 20 | 105 | 70 | 130 | |
| | | dibromochloromethane | | 16 ug/L | | 20 | 78 | 70 | 130 | |
| | | 1,2-dibromoethane (EDB) | | 20 ug/L | | 20 | 100 | 70 | 130 | |
| | | chlorobenzene | | 21 ug/L | | 20 | 104 | 70 | 130 | |
| | | 1,1,1,2-tetrachloroethane | | 17 ug/L | | 20 | 87 | 70 | 130 | |
| | | ethylbenzene | | 22 ug/L | | 20 | 109 | 70 | 130 | |
| | | m&p-xylenes | | 45 ug/L | | 40 | 113 | 70 | 130 | |
| | | o-xylene | | 23 ug/L | | 20 | 115 | 70 | 130 | |
| | | styrene | | 22 ug/L | | 20 | 109 | 70 | 130 | |
| | | bromoform | | 12 ug/L | | 20 | 61 | # 70 | 130 | |
| | | isopropylbenzene | | 22 ug/L | | 20 | 112 | 70 | 130 | |
| | | 1,1,2,2-tetrachloroethane | | 17 ug/L | | 20 | 84 | 70 | 130 | |

| Method | QC ID | Parameter | Associated Sample | Result | Units | Amt Added | %R | Limit | RPD | RPD Limit |
|--------------|------------|-----------------------------|-------------------|---------|-------|-----------|-----|-------|-----|-----------|
| SW503088260B | LCS0403790 | 1,2,3-trichloropropane | | 18 ug/L | | 20 | 90 | 70 | 130 | |
| | | n-propylbenzene | | 21 ug/L | | 20 | 105 | 70 | 130 | |
| | | bromobenzene | | 20 ug/L | | 20 | 99 | 70 | 130 | |
| | | 1,3,5-trimethylbenzene | | 22 ug/L | | 20 | 109 | 70 | 130 | |
| | | 2-chlorotoluene | | 21 ug/L | | 20 | 105 | 70 | 130 | |
| | | 4-chlorotoluene | | 20 ug/L | | 20 | 102 | 70 | 130 | |
| | | tert-butylbenzene | | 22 ug/L | | 20 | 111 | 70 | 130 | |
| | | 1,2,4-trimethylbenzene | | 22 ug/L | | 20 | 109 | 70 | 130 | |
| | | sec-butylbenzene | | 21 ug/L | | 20 | 106 | 70 | 130 | |
| | | 1,3-dichlorobenzene | | 19 ug/L | | 20 | 95 | 70 | 130 | |
| | | 4-isopropyltoluene | | 22 ug/L | | 20 | 111 | 70 | 130 | |
| | | 1,4-dichlorobenzene | | 19 ug/L | | 20 | 94 | 70 | 130 | |
| | | 1,2-dichlorobenzene | | 21 ug/L | | 20 | 104 | 70 | 130 | |
| | | n-butylbenzene | | 22 ug/L | | 20 | 108 | 70 | 130 | |
| | | 1,2-dibromo-3-chloropropane | | 13 ug/L | | 20 | 67 | # 70 | 130 | |
| | | 1,2,4-trichlorobenzene | | 19 ug/L | | 20 | 96 | 70 | 130 | |
| | | hexachlorobutadiene | | 21 ug/L | | 20 | 104 | 70 | 130 | |
| | | naphthalene | | 17 ug/L | | 20 | 84 | 70 | 130 | |
| | | 1,2,3-trichlorobenzene | | 19 ug/L | | 20 | 95 | 70 | 130 | |
| | | dibromofluoromethane SUR | | 98 | % | | | 78 | 114 | |
| | | toluene-D8 SUR | | 103 | % | | | 88 | 110 | |
| | | 4-bromofluorobenzene SUR | | 106 | % | | | 86 | 115 | |

| Method | QC ID | Parameter | Associated Sample | Result | Units | Amt Added | %R | Limit | RPD | RPD Limit |
|--------------|-------------|-----------------------------|-------------------|---------|-------|-----------|------|-------|-----|-----------|
| SW5030B8260B | LCSD0403790 | dichlorodifluoromethane | | 17 ug/L | 20 | 87 | 70 | 130 | 1 | 20 |
| | | chloromethane | | 21 ug/L | 20 | 103 | 70 | 130 | 0.2 | 20 |
| | | vinyl chloride | | 19 ug/L | 20 | 95 | 70 | 130 | 0.4 | 20 |
| | | bromomethane | | 18 ug/L | 20 | 89 | 70 | 130 | 2.2 | 20 |
| | | chloroethane | | 20 ug/L | 20 | 100 | 70 | 130 | 1.2 | 20 |
| | | trichlorofluoromethane | | 19 ug/L | 20 | 95 | 70 | 130 | 2.9 | 20 |
| | | diethyl ether | | 21 ug/L | 20 | 105 | 70 | 130 | 4.7 | 20 |
| | | acetone | | 21 ug/L | 20 | 103 | 70 | 130 | 0.9 | 20 |
| | | 1,1-dichloroethene | | 20 ug/L | 20 | 100 | 70 | 130 | 1.8 | 20 |
| | | methylene chloride | | 20 ug/L | 20 | 99 | 70 | 130 | 12 | 20 |
| | | carbon disulfide | | 16 ug/L | 20 | 82 | 70 | 130 | 0.4 | 20 |
| | | methyl t-butyl ether (MTBE) | | 20 ug/L | 20 | 101 | 70 | 130 | 1.9 | 20 |
| | | trans-1,2-dichloroethene | | 21 ug/L | 20 | 103 | 70 | 130 | 3 | 20 |
| | | Isopropyl ether (DIPE) | | 20 ug/L | 20 | 102 | 70 | 130 | 2.6 | 20 |
| | | ethyl t-butyl ether (ETBE) | | 20 ug/L | 20 | 102 | 70 | 130 | 3.1 | 20 |
| | | 1,1-dichloroethane | | 20 ug/L | 20 | 102 | 70 | 130 | 0.9 | 20 |
| | | t-butanol (TBA) | | 88 ug/L | 100 | 88 | 70 | 130 | 3.6 | 20 |
| | | 2-butanone (MEK) | | 18 ug/L | 20 | 92 | 70 | 130 | 4.9 | 20 |
| | | 2,2-dichloropropane | | 15 ug/L | 20 | 77 | 70 | 130 | 2.2 | 20 |
| | | cis-1,2-dichloroethene | | 22 ug/L | 20 | 109 | 70 | 130 | 1.9 | 20 |
| | | chloroform | | 21 ug/L | 20 | 105 | 70 | 130 | 1.4 | 20 |
| | | bromochloromethane | | 21 ug/L | 20 | 105 | 70 | 130 | 2.1 | 20 |
| | | tetrahydrofuran (THF) | | 19 ug/L | 20 | 96 | 70 | 130 | 4.9 | 20 |
| | | 1,1,1-trichloroethane | | 20 ug/L | 20 | 102 | 70 | 130 | 1.7 | 20 |
| | | 1,1-dichloropropene | | 21 ug/L | 20 | 104 | 70 | 130 | 3.4 | 20 |
| | | t-amyl-methyl ether (TAME) | | 20 ug/L | 20 | 100 | 70 | 130 | 4.1 | 20 |
| | | carbon tetrachloride | | 17 ug/L | 20 | 85 | 70 | 130 | 1 | 20 |
| | | 1,2-dichloroethane | | 21 ug/L | 20 | 107 | 70 | 130 | 0.3 | 20 |
| | | benzene | | 21 ug/L | 20 | 106 | 70 | 130 | 0.6 | 20 |
| | | trichloroethene | | 21 ug/L | 20 | 104 | 70 | 130 | 2.3 | 20 |
| | | 1,2-dichloropropane | | 21 ug/L | 20 | 104 | 70 | 130 | 1 | 20 |
| | | bromodichloromethane | | 18 ug/L | 20 | 91 | 70 | 130 | 1.8 | 20 |
| | | 1,4-dioxane | | 50 ug/L | 40 | 95 | 70 | 130 | 6.7 | 20 |
| | | dibromomethane | | 21 ug/L | 20 | 105 | 70 | 130 | 0.6 | 20 |
| | | 4-methyl-2-pentanone (MIBK) | | 18 ug/L | 20 | 90 | 70 | 130 | 2.7 | 20 |
| | | cis-1,3-dichloropropene | | 18 ug/L | 20 | 90 | 70 | 130 | 2 | 20 |
| | | toluene | | 22 ug/L | 20 | 109 | 70 | 130 | 2.4 | 20 |
| | | trans-1,3-dichloropropene | | 18 ug/L | 20 | 88 | 70 | 130 | 2.2 | 20 |
| | | 2-hexanone | | 17 ug/L | 20 | 83 | 70 | 130 | 2.5 | 20 |
| | | 1,1,2-trichloroethane | | 21 ug/L | 20 | 103 | 70 | 130 | 2.6 | 20 |
| | | 1,3-dichloropropane | | 20 ug/L | 20 | 101 | 70 | 130 | 1.8 | 20 |
| | | tetrachloroethene | | 21 ug/L | 20 | 103 | 70 | 130 | 2 | 20 |
| | | dibromochloromethane | | 16 ug/L | 20 | 79 | 70 | 130 | 0.8 | 20 |
| | | 1,2-dibromoethane (EDB) | | 20 ug/L | 20 | 98 | 70 | 130 | 1.2 | 20 |
| | | chlorobenzene | | 21 ug/L | 20 | 103 | 70 | 130 | 1.1 | 20 |
| | | 1,1,1,2-tetrachloroethane | | 18 ug/L | 20 | 88 | 70 | 130 | 1.7 | 20 |
| | | ethylbenzene | | 22 ug/L | 20 | 108 | 70 | 130 | 1 | 20 |
| | | m&p-xylenes | | 44 ug/L | 40 | 110 | 70 | 130 | 2.5 | 20 |
| | | o-xylene | | 22 ug/L | 20 | 110 | 70 | 130 | 5 | 20 |
| | | styrene | | 21 ug/L | 20 | 106 | 70 | 130 | 2.2 | 20 |
| | | bromoform | | 13 ug/L | 20 | 63 | # 70 | 130 | 3.2 | 20 |
| | | isopropylbenzene | | 22 ug/L | 20 | 110 | 70 | 130 | 1.1 | 20 |
| | | 1,1,2,2-tetrachloroethane | | 17 ug/L | 20 | 86 | 70 | 130 | 2.4 | 20 |

| Method | QC ID | Parameter | Associated Sample | Result | Units | Amt Added | %R | Limit | RPD | RPD Limit |
|--------------|-------------|-----------------------------|-------------------|---------|-------|-----------|-----|----------|-----|-----------|
| SW5030B8260B | LCSD0403790 | 1,2,3-trichloropropane | | 18 ug/L | | 20 | 90 | 70 130 | 0.2 | 20 |
| | | n-propylbenzene | | 22 ug/L | | 20 | 109 | 70 130 | 3.5 | 20 |
| | | bromobenzene | | 20 ug/L | | 20 | 102 | 70 130 | 3.1 | 20 |
| | | 1,3,5-trimethylbenzene | | 22 ug/L | | 20 | 111 | 70 130 | 1.8 | 20 |
| | | 2-chlorotoluene | | 22 ug/L | | 20 | 108 | 70 130 | 2.3 | 20 |
| | | 4-chlorotoluene | | 21 ug/L | | 20 | 105 | 70 130 | 2.4 | 20 |
| | | tert-butylbenzene | | 21 ug/L | | 20 | 105 | 70 130 | 5.8 | 20 |
| | | 1,2,4-trimethylbenzene | | 22 ug/L | | 20 | 110 | 70 130 | 1 | 20 |
| | | sec-butylbenzene | | 21 ug/L | | 20 | 103 | 70 130 | 2.7 | 20 |
| | | 1,3-dichlorobenzene | | 19 ug/L | | 20 | 97 | 70 130 | 1.6 | 20 |
| | | 4-isopropyltoluene | | 22 ug/L | | 20 | 109 | 70 130 | 2 | 20 |
| | | 1,4-dichlorobenzene | | 19 ug/L | | 20 | 97 | 70 130 | 2.8 | 20 |
| | | 1,2-dichlorobenzene | | 21 ug/L | | 20 | 105 | 70 130 | 0.6 | 20 |
| | | n-butylbenzene | | 21 ug/L | | 20 | 107 | 70 130 | 0.5 | 20 |
| | | 1,2-dibromo-3-chloropropane | | 14 ug/L | | 20 | 68 | # 70 130 | 1.8 | 20 |
| | | 1,2,4-trichlorobenzene | | 19 ug/L | | 20 | 93 | 70 130 | 3.3 | 20 |
| | | hexachlorobutadiene | | 20 ug/L | | 20 | 102 | 70 130 | 1.5 | 20 |
| | | naphthalene | < | 17 ug/L | | 20 | 86 | | 2.2 | 20 |
| | | 1,2,3-trichlorobenzene | | 19 ug/L | | 20 | 95 | 70 130 | 0.3 | 20 |
| | | dibromofluoromethane SUR | | 102 | % | | | 78 114 | | |
| | | toluene-D8 SUR | | 105 | % | | | 88 110 | | |
| | | 4-bromofluorobenzene SUR | | 105 | % | | | 86 115 | | |

Resource Laboratories, LLC

124 Heritage Avenue • Portsmouth, NH 03801
Phone: 603-436-2001 • Fax: 603-430-2100

Company Name: ESM Phone #: 603 435 4750

Company Address: 184 West Main St Northfield FAX #: 603 485 9957

Project Manager: Joe Callahan Project ID / Name: 2006-056 / DVD Lewis Chem

Invoice To: ESM Protocol: RCRA SDWA NPDES
AMCP NHDES OTHER

| Lab Sample ID (Lab Use Only) | Field ID | # CONTAINERS | Matrix | | | Preservation Method | | | | | | Sampling | | SAMPLER |
|---------------------------------|-----------|--------------|--------|-------|-------|---------------------|------------------|--------------------------------|------|------|-----------------|----------|-------|---------|
| | | | WATER | SOLID | OTHER | HCl | HNO ₃ | H ₂ SO ₄ | NaOH | MeOH | OTHER (Specify) | DATE | TIME | |
| 11371-01 | P2-04 | 2 | ✓ | | | ✓ | | | | | | 11/15/06 | 11:30 | AF |
| 11371-02 | P2-05 | 2 | ✓ | | | ✓ | | | | | | 11/15/06 | 12:20 | AF |
| 11371-03 | P2-06 | 2 | ✓ | | | ✓ | | | | | | 11/15/06 | 1:30 | AF |
| 11371-04 | P2-07 | 2 | ✓ | | | ✓ | | | | | | 11/15/06 | 1:50 | AF |
| 11371-05 | TRIP BUNK | 1 | | | | | | | | | | | | BY |

- ☐ VOC 8260-NH List ☐ MADEP VPH ☐ MEGRO
- ☒ VOC 8260 ☐ VOC8015GRO ☐ VOC 624
- ☐ VOC 8260 BTEX, MIBE, Naphthalene only
- ☐ VOC 524.2 ☐ VOC 524.2 NH List
- ☐ TPH Fingerprint ☐ MEDRO ☐ DRO 8015 ☐ EPH
- ☐ 8270PAH ☐ 8270ABN ☐ 625
- ☐ 8082 PCB ☐ 8081 Pesticides ☐ 608
- ☐ O&G 1664 ☐ O&G SM5520F
- ☐ pH ☐ BOD ☐ Conductivity
- ☐ TSS ☐ TDS ☐ TS
- ☐ RCRA Metals ☐ Priority Pollutant Metals ☐ TAL Metals
- ☐ Total Metals-list ☐ Dissolved Metals-list
- ☐ Ammonia ☐ COD
- ☐ T-Phosphorus ☐ Phenol
- ☐ Cyanide ☐ Sulfide
- ☐ Nitrate ☐ Nitrite ☐ Ortho P ☐ Sulfate ☐ Bromide ☐ Chloride
- ☐ Corrosivity ☐ Reactive CN ☐ Reactive S- ☐ Ignitability/FP
- ☐ TCLP Metals ☐ TCLP VOC ☐ TCLP SVOC
- ☐ TCLP Pesticide ☐ TCLP Herbicides (subcontract)
- ☐ Standard Drinking Water Test ☐ Bacteria P/A

Grab (G) or Composite (C)

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

ANALYSIS REQUEST

11371

| | | | | | |
|---|-----------------------------|---|---|-------------------------|---------------------|
| TAI REQUESTED | SPECIAL INSTRUCTIONS | REPORTING INSTRUCTIONS | RECEIVED ON ICE | TEMPERATURE | Lab Use Only |
| Priority (24 hr) <input type="checkbox"/> Expected (48 hr) <input type="checkbox"/> 10 Business Days <input type="checkbox"/> Other <input type="checkbox"/> | | <input type="checkbox"/> FAX <input type="checkbox"/> OTHER (specify) _____ <input checked="" type="checkbox"/> PDF <input type="checkbox"/> Excel Spreadsheet | RECEIVED ON ICE <input checked="" type="checkbox"/> YES <input type="checkbox"/> NO | TEMPERATURE <u>4</u> °C | |
| PO # _____ | | | | | |

| | | | | |
|-----------------------|---|--|---|--|
| CUSTODY RECORD | Relinquished by Sampler: <u>Gina-Ann Mearns</u> | Date: <u>11/16/06</u> Time: <u>12:00</u> | Received by: | Date: _____ Time: _____ |
| | Relinquished by: | Date: _____ Time: _____ | Received by: | Date: _____ Time: _____ |
| | Relinquished by: | Date: _____ Time: _____ | Received by Laboratory: <u>Thomas M. McLaughlin</u> | Date: <u>11-16-06</u> Time: <u>12:00</u> |