

## 1. INTRODUCTION

Task Order Title: NPDES Monitoring  
Contract Task Order #: 313150010  
SDG#: IOD0948  
Project Manager: B. McIlvaine  
Matrix: Water  
Analysis: Volatiles  
QC Level: Level IV  
No. of Samples: 2  
No. of Reanalyses/Dilutions: 0  
Reviewer: K. Shadowlight  
Date of Review: May 13, 2005

The samples listed in Table 1 were validated based on the guidelines outlined in the *AMEC Data Validation Procedure for Levels C and D Volatile Organics (DVP-2, Rev. 2)*, *EPA Method 624, SW846 Method 8260B*, and the *National Functional Guidelines For Organic Data Review (2/94)*. Any deviations from these procedures are documented herein. Qualifiers were applied in cases where the data did not meet the required QC criteria or where special consideration by the data user is required. Data qualifiers were placed on Form Is with the associated qualification codes. Analytes that were rejected for any reason are denoted on the summary forms as having only the "R" data qualifier and associated qualification code(s) denoting the reason for rejection. Any additional problems with the data that may have resulted in an estimated value were not denoted by a qualification code since the data had already been rejected.

**Table 1. Sample identification**

| Client ID   | EPA ID      | Lab No.    | Matrix | Method |
|-------------|-------------|------------|--------|--------|
| Outfall 012 | Outfall 012 | IOD0948-01 | water  | 624    |
| Trip Blank  | Trip Blank  | IOD0948-02 | water  | 624    |

## 2. DATA VALIDATION FINDINGS

### 2.1 SAMPLE MANAGEMENT

The following are findings associated with sample management:

#### 2.1.1 Sample Preservation, Handling, and Transport

The samples in this SDG were received at the laboratory within the temperature limits of 4°C ±2°C. The samples were properly preserved. The COC noted that the samples were received intact; however, information regarding absence of headspace was not provided. No qualifications were required.

#### 2.1.2 Chain of Custody

The COC was signed and dated by both field and laboratory personnel. The COC accounted for the analyses presented in this SDG. As the samples were couriered directly to the laboratory, custody seals were not required. No qualifications were required.

#### 2.1.3 Holding Times

The samples were analyzed within 14 days of collection. No qualifications were required.

### 2.2 GC/MS TUNING

The ion abundance windows shown on the quantitation reports were consistent with those specified in EPA Method 624, and all ion abundances were within the established windows. The samples and associated QC were analyzed within 12 hours of the BFB injection time. The BFB summary report was verified from the raw data and no discrepancies between the summary report and the raw data were noted. No qualifications were required.

### 2.3 CALIBRATION

One initial calibration dated 04/07/05 was associated with this SDG. The average RRFs were ≥0.05 for the target compounds listed on the sample result summaries. The %RSDs were ≤35% for all applicable target compounds. One continuing calibration dated 04/24/05 was associated with the sample analyses in this SDG. The %D for diisopropyl ether (DIPE) exceeded 20% in the continuing calibration dated 04/24/05; therefore, the nondetect result for diisopropyl ether was qualified as estimated, "UJ," in sample Outfall 012. No qualifications were required for the Trip Blank. The RRFs were ≥0.05 for the target compounds listed on the sample result summaries. A representative number of %RSDs and average RRFs from the initial calibration, and %Ds and RRFs from the continuing calibration were recalculated from the raw data, and no calculation or transcription errors were found. No further qualifications were required.

## 2.4 BLANKS

One water method blank (5D24007-BLK1) was associated with the sample analyses. There were no detects above the MDLs for the target compounds listed on the sample result summaries. The method blank raw data showed no evidence of false negatives. No qualifications were required.

## 2.5 BLANK SPIKES AND LABORATORY CONTROL SAMPLES

One water blank spike (5D24007-BS1) was associated with the sample analyses. All recoveries were within the laboratory-established QC limits. A representative number of recoveries were recalculated from the raw data and no calculation or transcription errors were found. No qualifications were required.

## 2.6 SURROGATE RECOVERY

The surrogates were recovered within the QC limits of 80-120% in the samples and associated QC. A representative number of surrogate recoveries were recalculated from the raw data and no calculation or transcription errors were found. No qualifications were required.

## 2.7 MATRIX SPIKE/MATRIX SPIKE DUPLICATE

MS/MSD analyses were not performed for this SDG. Evaluation of method accuracy was based on blank spike results. No qualifications were required.

## 2.8 FIELD QC SAMPLES

Field QC samples were evaluated, and if necessary, qualified based on method blanks and other laboratory QC results affecting the usability of the field QC data. Any remaining detects were used to evaluate the associated site sample. Following are findings associated with field QC samples:

### 2.8.1 Trip Blanks

Sample Trip Blank was the trip blank associated with this SDG. There were no target compounds detected above the MDLs in the trip blank. No qualifications were required.

### 2.8.2 Field Blanks and Equipment Rinsates

There were no field QC samples associated with this SDG. No qualifications were required.

### 2.8.3 Field Duplicates

There were no field duplicate samples associated with this SDG.

## 2.9 INTERNAL STANDARDS PERFORMANCE

Internal standard area counts and retention times for the samples in this SDG were within the control limits established by the continuing calibration standards: +100%/-50% for internal standard areas and  $\pm 0.50$  minutes for retention times. A representative number of internal standard areas and retention times were verified from the raw data, and no calculation or transcription errors were noted. No qualifications were required.

## 2.10 COMPOUND IDENTIFICATION

Target compound identification was verified at a Level IV data validation. The laboratory analyzed the volatile target compounds by EPA Method 624. Chromatograms, retention times, and spectra for the samples and QC were examined and no target compound identification problems were noted. No qualifications were required.

## 2.11 COMPOUND QUANTIFICATION AND REPORTED DETECTION LIMITS

Compound quantification is verified at a Level IV data validation. The reporting limits were supported by the lowest concentrations of the initial calibration standard and by the MDL study. As there were no sample detects in this SDG, compound quantitation was verified by recalculating a representative number of blank spike and surrogate recoveries from the raw data. Results were reported in  $\mu\text{g/L}$  (ppb). No calculation or transcription errors were noted. No qualifications were required.

## 2.12 TENTATIVELY IDENTIFIED COMPOUNDS

The laboratory did not provide TICs for this SDG. No qualifications were required.

## 2.13 SYSTEM PERFORMANCE

A review of the chromatograms and other raw data showed no identifiable problems with system performance. No qualifications were required.



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MWH-Pasadena/Boeing  
 300 North Lake Avenue, Suite 1200  
 Pasadena, CA 91101  
 Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD0948

Sampled: 04/13/05  
 Received: 04/13/05

## DRAFT: PURGEABLES BY GC/MS (EPA 624)

| Analyte   | Method  | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|---|---------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| <b>Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water)</b> |         |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                                     |         |         |           |                 |               |                 |                |               |                 |
| 1,2-Dibromoethane (EDB)                                   | EPA 624 | 5D24007 | 0.32      | 2.0             | ND            | 1               | 04/24/05       | 04/24/05      | U               |
| Methyl-tert-butyl Ether (MTBE)                            | EPA 624 | 5D24007 | 0.32      | 5.0             | ND            | 1               | 04/24/05       | 04/24/05      | ↓               |
| 1,2,3-Trichloropropane                                    | EPA 624 | 5D24007 | 0.85      | 10              | ND            | 1               | 04/24/05       | 04/24/05      | ↓               |
| Di-isopropyl Ether (DIPE)                                 | EPA 624 | 5D24007 | 0.25      | 5.0             | ND            | 1               | 04/24/05       | 04/24/05      | U.S             |
| tert-Butanol (TBA)  | EPA 624 | 5D24007 | 3.1       | 25              | ND            | 1               | 04/24/05       | 04/24/05      | U               |
| Surrogate: Dibromofluoromethane (80-120%)                 |         |         |           |                 | 113 %         |                 |                |               |                 |
| Surrogate: Toluene-d8 (80-120%)                           |         |         |           |                 | 106 %         |                 |                |               |                 |
| Surrogate: 4-Bromofluorobenzene (80-120%)                 |         |         |           |                 | 106 %         |                 |                |               |                 |
| <b>Sample ID: IOD0948-02 (DRAFT: Trip Blank - Water)</b>  |         |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                                     |         |         |           |                 |               |                 |                |               |                 |
| 1,2-Dibromoethane (EDB)                                   | EPA 624 | 5D24007 | 0.32      | 2.0             | ND            | 1               | 04/24/05       | 04/24/05      | U               |
| Methyl-tert-butyl Ether (MTBE)                            | EPA 624 | 5D24007 | 0.32      | 5.0             | ND            | 1               | 04/24/05       | 04/24/05      | ↓               |
| 1,2,3-Trichloropropane                                    | EPA 624 | 5D24007 | 0.85      | 10              | ND            | 1               | 04/24/05       | 04/24/05      | ↓               |
| Di-isopropyl Ether (DIPE)                                 | EPA 624 | 5D24007 | 0.25      | 5.0             | ND            | 1               | 04/24/05       | 04/24/05      | ↓               |
| tert-Butanol (TBA)  | EPA 624 | 5D24007 | 3.1       | 25              | ND            | 1               | 04/24/05       | 04/24/05      | ↓               |
| Surrogate: Dibromofluoromethane (80-120%)                 |         |         |           |                 | 107 %         |                 |                |               |                 |
| Surrogate: Toluene-d8 (80-120%)                           |         |         |           |                 | 106 %         |                 |                |               |                 |
| Surrogate: 4-Bromofluorobenzene (80-120%)                 |         |         |           |                 | 104 %         |                 |                |               |                 |

**AMEC VALIDATED  
 LEVEL IV**

DRAFT REPORT  
 DRAFT REPORT  
 DATA SUBJECT TO CHANGE

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**CONTRACT COMPLIANCE SCREENING FORM FOR HARDCOPY DATA**

AMEC Earth & Environmental  
 550 South Wadsworth Boulevard  
 Suite 500  
 Lakewood, CO 80226

Package ID T711VO102  
 Task Order 313150010  
 SDG No. IOD0948

No. of Analyses 1

Laboratory Del Mar Analytical  
 Reviewer K. Shadowlight  
 Analysis/Method 1,4-Dioxane by 8260

Date May 13, 2005  
 Reviewer's Signature K. Shadowlight

| ACTION ITEMS*   |  |
|---|--|
| <b>1. Case Narrative</b>  |  |
| Deficiencies  |  |
| <b>2. Out of Scope</b>  |  |
| Analyses  |  |
| <b>3. Analyses Not Conducted</b>  |  |
| <b>4. Missing Hardcopy Deliverables</b>   |  |
| <b>5. Incorrect Hardcopy Deliverables</b>   |  |
| <b>6. Deviations from Analysis</b>  | Qualification was assigned for a %D continuing calibration outlier |
| GC/MS Tune/Inst. Perform  |  |
| Calibrations  |  |
| Blanks  |  |
| Surrogates  |  |
| Matrix Spike/Dup LCS  |  |
| Field QC  |  |
| Internal Standard Performance   |  |
| Compound Identification and Quantitation  |  |
| System Performance  |  |
| <b>COMMENTS<sup>b</sup></b>   |  |
|   |  |
|   |  |
|   |  |
| <p>* Subcontracted analytical laboratory is not meeting contract and/or method requirements.<br/>                     * Differences in protocol have been adopted by the laboratory but no action against the laboratory is required.</p> |  |

### Data Qualifier Reference Table

| Qualifier | Organics  | Inorganics   |
|-----------|---|--|
| U         | The analyte was analyzed for, but was not detected above the reported sample quantitation limit.  | The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit. |
| J         | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.  | The associated value is an estimated quantity.   |
| N         | The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification."   | Not applicable.  |
| NJ        | The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.   | Not applicable.  |
| UJ        | The analyte was not deemed above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. | The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.   |
| R         | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and to meet quality control criteria. The presence or absence of the analyte cannot be verified.   | The data are unusable. (Note: Analyte may or may not be present).  |



## Qualification Code Reference Table

| Qualifier | Organics   | Inorganics   |
|-----------|--|--|
| H         | Holding times were exceeded.   | Holding times were exceeded.   |
| S         | Surrogate recovery was outside QC limits.  | The sequence or number of standards used for the calibration was incorrect   |
| C         | Calibration %RSD or %D were noncompliant.  | Correlation coefficient is <0.995.   |
| R         | Calibration RRF was <0.05.   | %R for calibration is not within control limits.   |
| B         | Presumed contamination from preparation (method) blank.  | Presumed contamination from preparation (method) or calibration blank.   |
| L         | Laboratory Blank Spike/Blank Spike Duplicate %R was not within control limits.   | Laboratory Control Sample %R was not within control limits.  |
| Q         | MS/MSD recovery was poor or RPD high.  | MS recovery was poor.  |
| E         | Not applicable.  | Duplicates showed poor agreement.  |
| I         | Internal standard performance was unsatisfactory.  | ICP ICS results were unsatisfactory.   |
| A         | Not applicable.  | ICP Serial Dilution %D were not within control limits.   |
| M         | Tuning (BFB or DFTPP) was noncompliant.  | Not applicable.  |
| T         | Presumed contamination from trip blank.  | Not applicable.  |
| +         | False positive – reported compound was not present. Not applicable.  |  |
| -         | False negative – compound was present but not reported.  | Not applicable.  |
| F         | Presumed contamination from FB, or ER.   | Presumed contamination from FB or ER.  |
| \$        | Reported result or other information was incorrect.  | Reported result or other information was incorrect.  |
| ?         | TIC identity or reported retention time has been changed.  | Not applicable.  |
| D         | The analysis with this flag should not be used because another more technically sound analysis is available.   | The analysis with this flag should not be used because another more technically sound analysis is available.   |
| P         | Instrument performance for pesticides was poor.  | Post Digestion Spike recovery was not within control limits.   |
| DNQ       | The compound was detected between the MDL and the RL and, by definition, is considered an estimated value.   | The compound was detected between the MDL and the RL and, by definition, is considered an estimated value.   |
| *#        | Unusual problems found with the data that have been described in Section 2.#, "Data Validation Findings." The number following the asterisk (*) will indicate the subsection where a description of the problem can be found (eg. *1 would indicate a sample was not within temperature limits). | Unusual problems found with the data that have been described in Section 2.#, "Data Validation Findings." The number following the asterisk (*) will indicate the subsection where a description of the problem can be found (eg. *1 would indicate a sample was not within temperature limits). |



# DATA VALIDATION REPORT

NPDES Monitoring

ANALYSIS: VOLATILES

SAMPLE DELIVERY GROUP: IOD0948

Prepared by

AMEC—Denver Operations  
550 South Wadsworth Boulevard, Suite 500  
Lakewood, Colorado 80226

## 1. INTRODUCTION

Task Order Title: NPDES Monitoring  
Contract Task Order #: 313150010  
Sample Delivery Group #: IOD0948  
Project Manager: B. McIlvaine  
Matrix: Water  
Analysis: Volatiles (1,4-dioxane)  
QC Level: Level IV  
No. of Samples: 1  
No. of Reanalyses/Dilutions: 0  
Reviewer: K. Shadowlight  
Date of Review: May 13, 2005

The samples listed in Table 1 were validated based on the guidelines outlined in *the AMEC Data Validation Procedure for Levels C and D Volatile Organics (DVP-2, Rev. 2)*, *EPA Method SW-846 8260B* and the *National Functional Guidelines For Organic Data Review (2/94)*. Any deviations from these procedures and guidelines are documented herein. Qualifiers were applied in cases where the data did not meet the required QC criteria or where special consideration by the data user is required. Data qualifiers were placed on Form Is with the associated qualification codes. Analytes that were rejected for any reason are denoted on the Form I as having only the "R" data qualifier and associated qualification code(s) denoting the reason for rejection. Any additional problems with the data that may have resulted in an estimated value were not denoted by a qualification code since the data had already been rejected.

**Table 1. Sample identification**

| Client ID   | EPA ID      | Lab No.<br>Del Mar, CA | Lab No.<br>Del Mar, AZ | Matrix | Method |
|-------------|-------------|------------------------|------------------------|--------|--------|
| Outfall 012 | Outfall 012 | IOD0948-01             | POD0411-01             | water  | 8260B  |

## 2. DATA VALIDATION FINDINGS

### 2.1 SAMPLE MANAGEMENT

Following are findings associated with sample management:

#### 2.1.1 Sample Preservation, Handling, and Transport

The sample in this SDG was received at the Del Mar within the temperature limits of  $4^{\circ}\text{C} \pm 2^{\circ}\text{C}$ . The sample was subcontracted to Del Mar (Phoenix) for 1,4-dioxane analysis, and the sample was received within the temperature limits of  $4^{\circ}\text{C} \pm 2^{\circ}\text{C}$ . The sample was properly preserved. The COC and transfer COC noted that the sample was received intact; however, information regarding absence of headspace was not provided. No qualifications were required.

#### 2.1.2 Chain of Custody

The COC and transfer COC were signed by field and laboratory personnel. As the sample was couriered directly to the laboratory from the field, custody seals were not required. According to the transfer COC, there were no custody seals present on the cooler received by Del Mar Analytical in Arizona. No qualifications were required.

#### 2.1.3 Holding Times

The sample was analyzed within 14 days of collection. No qualifications were required.

### 2.2 GC/MS TUNING

The ion abundance windows were consistent with those specified in EPA Method 8260B. All ion abundances were within the established windows, and the sample was analyzed within 12 hours of the BFB injection time. No qualifications were required.

### 2.3 CALIBRATION

One initial calibration, dated 03/19/05, was associated with this SDG. The average RRF for 1,4-dioxane was  $\geq 0.05$  and the  $r^2$  value was  $\geq 0.995$ . The laboratory reported the continuing calibration and the blank spike (P5D1803-BS1) from the same analysis. As the analysis cannot be reported as both a CCV and a blank spike, the reviewer reported P5D1803-BS1 as the continuing calibration. The RRF for 1,4-dioxane was  $\geq 0.05$ ; however, the %D exceeded 20%. The nondetect result for 1,4-dioxane was qualified as estimated, "UJ," in the site sample. The  $r^2$  value and average RRF for 1,4-dioxane in the initial calibration, and the %D and RRF for 1,4-dioxane in the continuing calibration were recalculated from the raw data, and no calculation or transcription errors were found. No further qualifications were required.

## 2.4 BLANKS

One water method blank (P5D1803-BLK1) was associated with this SDG. Target compound 1,4-dioxane was not detected above the MDL in the method blank. The method blank raw data showed no evidence of a false negative. No qualifications were required.

## 2.5 BLANK SPIKES AND LABORATORY CONTROL SAMPLES

The laboratory analyzed a blank spike/blank spike duplicate pair (P5D1803-BS1/BS1D) with this SDG; however, P5D1803-BS1 was reported as the CCV (see section 2.3); therefore, P5D1803-BS1D was evaluated as a single blank spike. The recovery for 1,4-dioxane was within the QC limits of 70-130%. The recovery was recalculated from the raw data and no calculation or transcription errors were found. No qualifications were required.

## 2.6 SURROGATE RECOVERY

The sample and QC were fortified with dibromofluoromethane. The surrogate was recovered within the laboratory QC limits of 80-125%. The surrogate recovery for the sample was recalculated from the raw data and no calculation or transcription errors were found. No qualifications were required.

## 2.7 MATRIX SPIKE/MATRIX SPIKE DUPLICATE

No MS/MSD analyses were associated with this SDG. Evaluation of method accuracy was based on blank spike results. No qualifications were required.

## 2.8 FIELD QC SAMPLES

Field QC samples were evaluated, and if necessary, qualified based on method blanks and other laboratory QC results affecting the usability of the field QC data. Any remaining detects were used to evaluate the associated site sample. Following are findings associated with field QC samples:

### 2.8.1 Trip Blanks

The sample in this SDG had no associated trip blank. No qualifications were required.

#### 2.8.1.1 Field Blanks and Equipment Rinsates

The site sample in this SDG had no associated field QC samples. No qualifications were required.

### 2.8.2 Field Duplicates

There were no field duplicate samples associated with this SDG.

## 2.9 INTERNAL STANDARDS PERFORMANCE

Internal standard area counts and retention times for the sample were within the control limits established by the continuing calibration standard: +100%/-50% for internal standard areas and  $\pm 0.50$  minutes for retention times. Internal standard areas and retention times were verified from the raw data, and no calculation or transcription errors were noted. No qualifications were required.

## 2.10 COMPOUND IDENTIFICATION

Target compound identification was verified at a Level IV data validation. The laboratory analyzed for 1,4-dioxane by Method 8260B/SIM. Chromatograms, retention times, and spectra for the sample and QC were examined and no target compound identification problems were noted. No qualifications were required.

## 2.11 COMPOUND QUANTIFICATION AND REPORTED DETECTION LIMITS

Compound quantification is verified at a Level IV data validation. The reporting limit was supported by the lowest concentration of the initial calibration standards and by the undated MDL supplied by the laboratory. Compound quantitation was verified by recalculating blank spike and surrogate recoveries from the raw data. No calculation or transcription errors were noted. No qualifications were required.

## 2.12 TENTATIVELY IDENTIFIED COMPOUNDS

TICs are not typically reported for SIM methods.

## 2.13 SYSTEM PERFORMANCE

A review of the chromatograms and other raw data showed no identifiable problems with system performance. No qualifications were required.



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 9830 South 51st St., Suite B-120, Phoenix, AZ 85044 (480) 785-0044 FAX (480) 785-0871  
 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

MWH-Pasadena/Boeing  
 300 North Lake Avenue, Suite 1200  
 Pasadena, CA 91101  
 Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD0948

Sampled: 04/13/05  
 Received: 04/13/05

## DRAFT: 1,4-DIOXANE BY GC/MS (EPA 5030B/8260B)

| Analyte  | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed |           | Data Qualifiers |  |
|--|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------|-----------------|--|
|  |           |         |           |                 |               |                 |                | Key Qual      | Qual Lock |                 |  |
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water) - cont. |           |         |           |                 |               |                 |                |               |           |                 |  |
| Reporting Units: ug/l                                      |           |         |           |                 |               |                 |                |               |           |                 |  |
| 1,4-Dioxane  | EPA 8260B | P5D1803 | 0.49      | 1.0             | ND            | 1               | 04/18/05       | 04/18/05      | US        | C               |  |
| Surrogate: Dibromofluoromethane (80-125%)                  |           |         |           |                 | 117%          |                 |                |               |           |                 |  |

**AMEC VALIDATED**

**LEVEL IV**

DRAFT REPORT  
 DRAFT REPORT  
 DATA SUBJECT TO CHANGE

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### Data Qualifier Reference Table

| Qualifier | Organics  | Inorganics   |
|-----------|---|--|
| U         | The analyte was analyzed for, but was not detected above the reported sample quantitation limit.  | The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit. |
| J         | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.  | The associated value is an estimated quantity.   |
| N         | The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification."   | Not applicable.  |
| NJ        | The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.   | Not applicable.  |
| UJ        | The analyte was not deemed above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. | The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.   |
| R         | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and to meet quality control criteria. The presence or absence of the analyte cannot be verified.   | The data are unusable. (Note: Analyte may or may not be present).  |

## Qualification Code Reference Table

| Qualifier | Organics   | Inorganics   |
|-----------|--|--|
| H         | Holding times were exceeded.   | Holding times were exceeded.   |
| S         | Surrogate recovery was outside QC limits.  | The sequence or number of standards used for the calibration was incorrect   |
| C         | Calibration %RSD or %D were noncompliant.  | Correlation coefficient is <0.995.   |
| R         | Calibration RRF was <0.05.   | %R for calibration is not within control limits.   |
| B         | Presumed contamination from preparation (method) blank.  | Presumed contamination from preparation (method) or calibration blank.   |
| L         | Laboratory Blank Spike/Blank Spike Duplicate %R was not within control limits.   | Laboratory Control Sample %R was not within control limits.  |
| Q         | MS/MSD recovery was poor or RPD high.  | MS recovery was poor.  |
| E         | Not applicable.  | Duplicates showed poor agreement.  |
| I         | Internal standard performance was unsatisfactory.  | ICP ICS results were unsatisfactory.   |
| A         | Not applicable.  | ICP Serial Dilution %D were not within control limits.   |
| M         | Tuning (BFB or DFTPP) was noncompliant.  | Not applicable.  |
| T         | Presumed contamination from trip blank.  | Not applicable.  |
| +         | False positive – reported compound was not present. Not applicable.  |  |
| -         | False negative – compound was present but not reported.  | Not applicable.  |
| F         | Presumed contamination from FB, or ER.   | Presumed contamination from FB or ER.  |
| \$        | Reported result or other information was incorrect.  | Reported result or other information was incorrect.  |
| ?         | TIC identity or reported retention time has been changed.  | Not applicable.  |
| D         | The analysis with this flag should not be used because another more technically sound analysis is available.   | The analysis with this flag should not be used because another more technically sound analysis is available.   |
| P         | Instrument performance for pesticides was poor.  | Post Digestion Spike recovery was not within control limits.   |
| DNQ       | The compound was detected between the MDL and the RL and, by definition, is considered an estimated value.   | The compound was detected between the MDL and the RL and, by definition, is considered an estimated value.   |
| *#        | Unusual problems found with the data that have been described in Section 2.#, "Data Validation Findings." The number following the asterisk (*) will indicate the subsection where a description of the problem can be found (eg. *1 would indicate a sample was not within temperature limits). | Unusual problems found with the data that have been described in Section 2.#, "Data Validation Findings." The number following the asterisk (*) will indicate the subsection where a description of the problem can be found (eg. *1 would indicate a sample was not within temperature limits). |



# DATA VALIDATION REPORT

NPDES Monitoring

ANALYSIS: GENERAL MINERALS

SAMPLE DELIVERY GROUP: IOD0948

Prepared by

AMEC—Denver Operations  
550 South Wadsworth Boulevard, Suite 500  
Lakewood, Colorado 80226

## 1. INTRODUCTION

Task Order Title: NPDES Monitoring  
Contract Task Order #: 313150010  
Sample Delivery Group #: IOD0948  
Project Manager: B. McIlvaine  
Matrix: Water  
Analysis: General Minerals  
QC Level: Level IV  
No. of Samples: 1  
Reviewer: L. Jarusewic  
Date of Review: March 10, 2005

The sample listed in Table 1 was validated based on the guidelines outlined in the AMEC *Data Validation Procedures SOP DVP-6, Rev. 2, USEPA Methods for Chemical Analysis of Water and Wastes Method 350.2, 405.1, 413.1, 418.1, 160.2, 160.5, and 180.1, Standard Methods for the Examination of Water and Wastewater Method SM2540C*, and validation guidelines outlined in the USEPA *Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (2/94)*. Any deviations from these procedures and guidelines are documented herein. Qualifiers were applied in cases where the data did not meet the required QC criteria or where special consideration by the data user is required. Data qualifiers were placed on Form Is with the associated qualification codes. Analytes that were rejected for any reason are denoted on the Form I as having only the "R" data qualifier and associated qualification code(s) denoting the reason for rejection. Any additional problems with the data that may have resulted in an estimated value were not denoted by a qualification code since the data had already been rejected.

**Table 1. Sample identification**

| Client ID   | EPA ID      | Laboratory ID | Matrix | COC Method       |
|-------------|-------------|---------------|--------|------------------|
| Outfall 012 | Outfall 012 | IOD0948-01    | Water  | General Minerals |

## 2. DATA VALIDATION FINDINGS

### 2.1 SAMPLE MANAGEMENT

Following are findings associated with sample management:

#### 2.1.1 Sample Preservation, Handling, and Transport

The sample in this SDG was received at the laboratory within the temperature limits of  $4^{\circ}\text{C} \pm 2^{\circ}\text{C}$ . No preservation problems were noted by the laboratory. No qualifications were required.

#### 2.1.2 Chain of Custody

The COC was signed and dated by field and laboratory personnel. The COC accounted for all analyses present in this SDG. No sample qualifications were required.

#### 2.1.3 Holding Times

The holding times were assessed by comparing the date of collection with the dates of analyses. The 28-day analytical holding time for ammonia, total recoverable hydrocarbons, and oil and grease, the seven-day holding time for total suspended solids and total dissolved solids, and the 48-hour holding time for turbidity, biological oxygen demand, and total settleable solids were met. No qualifications were required.

### 2.2 CALIBRATION

For the applicable analyses, the initial calibration correlation coefficients were  $\geq 0.995$ . Initial and continuing calibration information was acceptable with recoveries within the control limits of 90-110%. For ammonia, no information regarding the standardization of the titrant was provided; however, as the LCS recovery was within the CCV control limits, no qualifications were required. For BOD, no information regarding the calibration of the oxygen meter was provided; however, as the LCS recovery was within the CCV control limits, no qualifications were required. Calibration is not applicable to total settleable solids or oil and grease. No qualifications were required.

### 2.3 BLANKS

Turbidity was detected in a bracketing CCB at 0.040NTU; however, the turbidity CCB result was insufficient to qualify the Outfall 012 result. The remaining method blank and CCB results reported on the summary forms and in the raw data for blank analyses associated with the sample were nondetects at the reporting limit. No qualifications were required.

## **2.4 BLANK SPIKES AND LABORATORY CONTROL SAMPLES**

The laboratory control sample and laboratory control sample duplicate (BOD, oil and grease, and total recoverable hydrocarbons only) recoveries and RPDs were within the laboratory-established control limits. The LCS is not applicable to turbidity or settleable solids. No qualifications were required.

## **2.5 SURROGATES RECOVERY**

Surrogate recovery is not applicable to the analyses presented in this SDG.

## **2.6 LABORATORY DUPLICATES**

No MS/MSD analyses were performed in association with the sample in this SDG; therefore, no assessment was made with respect to this criterion.

## **2.7 MATRIX SPIKE/MATRIX SPIKE DUPLICATE**

No MS/MSD analyses were performed in association with the sample in this SDG; therefore, no assessment was made with respect to this criterion. Method accuracy was based on LCS results.

## **2.8 FURNACE ATOMIC ABSORPTION QC**

Furnace atomic absorption was not utilized for the analyses of this sample; therefore, furnace atomic absorption QC is not applicable.

## **2.9 ICP SERIAL DILUTION**

ICP serial dilution is not applicable to the analyses presented in this data validation report.

## **2.10 SAMPLE RESULT VERIFICATION**

A Level IV review was performed for the samples in these data packages. Calculations were verified, and the sample results reported on the Form Is were verified against the raw data. No transcription errors or calculation errors were noted. Oil and grease detected below the reporting limit was qualified as estimated, "J." No further qualifications were required.



## **2.11 FIELD QC SAMPLES**

Field QC samples are evaluated, and if necessary, qualified based only on laboratory blanks. Any remaining detects are used to evaluate the associated sample. The following are findings associated with field QC samples:

### **2.11.1 Field Blanks and Equipment Rinsates**

The sample in this SDG had no associated field QC samples. No qualifications were required.

### **2.11.2 Field Duplicates**

There were no field duplicate pairs associated with this SDG.



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MWH-Pasadena/Boeing  
 300 North Lake Avenue, Suite 1200  
 Pasadena, CA 91101  
 Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD0948

Sampled: 04/13/05  
 Received: 04/13/05

## DRAFT: INORGANICS

| Analyte  | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|--|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water) - cont. |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: mg/l                                      |           |         |           |                 |               |                 |                |               |                 |
| Ammonia-N (Distilled)                                      | EPA 350.2 | 5D14077 | 0.30      | 0.50            | ND            | 1               | 04/14/05       | 04/14/05      | U               |
| Biochemical Oxygen Demand                                  | EPA 405.1 | 5D14060 | 0.59      | 2.0             | 3.8           | 1               | 04/14/05       | 04/19/05      | J               |
| Oil & Grease   | EPA 413.1 | 5D14081 | 0.94      | 5.0             | 4.2           | 1               | 04/14/05       | 04/14/05      | J               |
| Total Dissolved Solids                                     | SM2540C   | 5D18095 | 10        | 10              | 240           | 1               | 04/18/05       | 04/18/05      | J               |
| Total Suspended Solids                                     | EPA 160.2 | 5D18089 | 10        | 10              | 12            | 1               | 04/18/05       | 04/18/05      | J               |
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water)         |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ml/hr                                     |           |         |           |                 |               |                 |                |               |                 |
| Total Settleable Solids                                    | EPA 160.5 | 5D15069 | 0.10      | 0.10            | ND            | 1               | 04/15/05       | 04/15/05      | U               |
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water)         |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: NTU                                       |           |         |           |                 |               |                 |                |               |                 |
| Turbidity  | EPA 180.1 | 5D15076 | 0.040     | 1.0             | 25            | 1               | 04/15/05       | 04/15/05      | J               |
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water)         |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                                      |           |         |           |                 |               |                 |                |               |                 |
| Perchlorate  | EPA 314.0 | 5D14043 | 0.80      | 4.0             | ND            | 1               | 04/14/05       | 04/14/05      | *               |

REV  
QUAL  
CODE

DNQ

**AMEC VALIDATED**

**LEVEL IV**

\*Analysis Not Validated

DRAFT REPORT  
 DRAFT REPORT  
 DATA SUBJECT TO CHANGE

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MWH-Pasadena/Boeing  
 300 North Lake Avenue, Suite 1200  
 Pasadena, CA 91101  
 Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD0948

Sampled: 04/13/05  
 Received: 04/13/05

**DRAFT: TOTAL RECOVERABLE PETROLEUM HYDROCARBONS (EPA 418.1)**

| Analyte  | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers  |
|--|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|------------------|
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water) |           |         |           |                 |               |                 |                |               |                  |
| Reporting Units: mg/l                              |           |         |           |                 |               |                 |                |               |                  |
| Total Recoverable Hydrocarbons                     | EPA 418.1 | 5D21036 | 0.31      | 1.0             | 9.8           | 1               | 04/21/05       | 04/21/05      | REV<br>QUAL CODE |

**AMEC VALIDATED**

**LEVEL III**

DRAFT REPORT  
 DRAFT REPORT  
 DATA SUBJECT TO CHANGE

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### Data Qualifier Reference Table

| Qualifier | Organics  | Inorganics   |
|-----------|---|--|
| U         | The analyte was analyzed for, but was not detected above the reported sample quantitation limit.  | The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit. |
| J         | The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.  | The associated value is an estimated quantity.   |
| N         | The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification."   | Not applicable.  |
| NJ        | The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.   | Not applicable.  |
| UJ        | The analyte was not deemed above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample. | The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.   |
| R         | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and to meet quality control criteria. The presence or absence of the analyte cannot be verified.   | The data are unusable. (Note: Analyte may or may not be present).  |

## Qualification Code Reference Table

| Qualifier | Organics   | Inorganics   |
|-----------|--|--|
| H         | Holding times were exceeded.   | Holding times were exceeded.   |
| S         | Surrogate recovery was outside QC limits.  | The sequence or number of standards used for the calibration was incorrect   |
| C         | Calibration %RSD or %D were noncompliant.  | Correlation coefficient is <0.995.   |
| R         | Calibration RRF was <0.05.   | %R for calibration is not within control limits.   |
| B         | Presumed contamination from preparation (method) blank.  | Presumed contamination from preparation (method) or calibration blank.   |
| L         | Laboratory Blank Spike/Blank Spike Duplicate %R was not within control limits.   | Laboratory Control Sample %R was not within control limits.  |
| Q         | MS/MSD recovery was poor or RPD high.  | MS recovery was poor.  |
| E         | Not applicable.  | Duplicates showed poor agreement.  |
| I         | Internal standard performance was unsatisfactory.  | ICP ICS results were unsatisfactory.   |
| A         | Not applicable.  | ICP Serial Dilution %D were not within control limits.   |
| M         | Tuning (BFB or DFTPP) was noncompliant.  | Not applicable.  |
| T         | Presumed contamination from trip blank.  | Not applicable.  |
| +         | False positive - reported compound was not present. Not applicable.  |  |
| -         | False negative - compound was present but not reported.  | Not applicable.  |
| F         | Presumed contamination from FB, or ER.   | Presumed contamination from FB or ER.  |
| S         | Reported result or other information was incorrect.  | Reported result or other information was incorrect.  |
| ?         | TIC identity or reported retention time has been changed.  | Not applicable.  |
| D         | The analysis with this flag should not be used because another more technically sound analysis is available.   | The analysis with this flag should not be used because another more technically sound analysis is available.   |
| P         | Instrument performance for pesticides was poor.  | Post Digestion Spike recovery was not within control limits.   |
| DNQ       | The compound was detected between the MDL and the RL and, by definition, is considered an estimated value.   | The compound was detected between the MDL and the RL and, by definition, is considered an estimated value.   |
| *#        | Unusual problems found with the data that have been described in Section 2.#, "Data Validation Findings." The number following the asterisk (*) will indicate the subsection where a description of the problem can be found (eg. *1 would indicate a sample was not within temperature limits). | Unusual problems found with the data that have been described in Section 2.#, "Data Validation Findings." The number following the asterisk (*) will indicate the subsection where a description of the problem can be found (eg. *1 would indicate a sample was not within temperature limits). |



# DATA VALIDATION REPORT

## NPDES Monitoring

ANALYSIS: PERCHLORATE

SAMPLE DELIVERY GROUP: IOD0948

Prepared by

AMEC—Denver Operations  
550 South Wadsworth Boulevard, Suite 500  
Lakewood, Colorado 80226

## 1. INTRODUCTION

Task Order Title: NPDES Monitoring  
Contract Task Order #: 313150010  
Sample Delivery Group #: IOD0948  
Project Manager: B. McIlvaine  
Matrix: Water  
Analysis: Perchlorate  
QC Level: Level IV  
No. of Samples: 1  
Reviewer: L. Jarusewic  
Date of Review: May 18, 2005

The samples listed in Table 1 was validated based on the guidelines outlined in the AMEC *Data Validation Procedures SOP DVP-6, Rev. 2*, *USEPA Methods for Chemical Analysis of Water and Wastes Method 314.0*, and validation guidelines outlined in the *USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (2/94)*. Any deviations from these procedures and guidelines are documented herein. Qualifiers were applied in cases where the data did not meet the required QC criteria or where special consideration by the data user is required. Data qualifiers were placed on Form Is with the associated qualification codes. Analytes that were rejected for any reason are denoted on the Form I as having only the "R" data qualifier and associated qualification code(s) denoting the reason for rejection. Any additional problems with the data that may have resulted in an estimated value were not denoted by a qualification code since the data had already been rejected.



**Table 1. Sample identification**

| Client ID   | EPA ID      | Laboratory ID | Matrix | COC Method  |
|-------------|-------------|---------------|--------|-------------|
| Outfall 012 | Outfall 012 | IOD0948-01    | Water  | Perchlorate |

## 2. DATA VALIDATION FINDINGS

### 2.1 SAMPLE MANAGEMENT

Following are findings associated with sample management:

#### 2.1.1 Sample Preservation, Handling, and Transport

The sample in this SDG was received at the laboratory within the temperature limits of  $4^{\circ}\text{C} \pm 2^{\circ}\text{C}$ . The analysis did not require preservation and no preservation was noted in the field. No qualifications were required.

#### 2.1.2 Chain of Custody

The COC was signed and dated by field and laboratory personnel, and accounted for the sample and analysis presented in this SDG. No qualifications were required.

#### 2.1.3 Holding Times

The holding time was assessed by comparing the date of collection with the date of analysis. The 28-day analytical holding time for perchlorate was met, and no qualifications were required.

### 2.2 CALIBRATION

The initial calibration correlation coefficient was  $\geq 0.995$ . The IPC-MA recovery was within the control limits of 80-120%. The ICV, CCV, ICCS, and IPC recoveries were within the control limits of 90-110%. No qualifications were required.

### 2.3 BLANKS

The method blank and CCB results reported on the summary forms and in the raw data for the blank analyses associated with the sample were nondetects at the reporting limit. No qualifications were required.

### 2.4 BLANK SPIKES AND LABORATORY CONTROL SAMPLES

The laboratory control sample recovery was within the method control limits of 85-115%. No qualifications were required.

### 2.5 SURROGATES RECOVERY

Surrogate recovery is not applicable to the analysis presented in this SDG.

## 2.6 LABORATORY DUPLICATES

No MS/MSD or duplicate analyses were performed in association with the sample in this SDG; therefore, no assessment was made with respect to this criterion.

## 2.7 MATRIX SPIKE/MATRIX SPIKE DUPLICATE

No MS/MSD analyses were performed in association with the sample in this SDG; therefore, no assessment was made with respect to this criterion. Method accuracy was assessed based on LCS results.

## 2.8 FURNACE ATOMIC ABSORPTION QC

Furnace atomic absorption was not utilized for the analysis of this sample; therefore, furnace atomic absorption QC is not applicable.

## 2.9 ICP SERIAL DILUTION

ICP serial dilution is not applicable to the analysis presented in this data validation report.

## 2.10 SAMPLE RESULT VERIFICATION

A Level IV review was performed for the sample in this data package. Calculations were verified, and the sample result reported on the Form I was verified against the raw data. No transcription errors or calculation errors were noted. No qualifications were required.

## 2.11 FIELD QC SAMPLES

Field QC samples are evaluated, and if necessary, qualified based only on laboratory blanks. Any remaining detects are used to evaluate the associated sample. The following are findings associated with field QC samples:

### 2.11.1 Field Blanks and Equipment Rinsates

The sample in this SDG had no associated field QC samples. No qualifications were required.

### 2.11.2 Field Duplicates

There were no field duplicate pairs associated with this SDG.



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MWH-Pasadena/Boeing  
 300 North Lake Avenue, Suite 1200  
 Pasadena, CA 91101  
 Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD0948

Sampled: 04/13/05  
 Received: 04/13/05

## DRAFT: INORGANICS

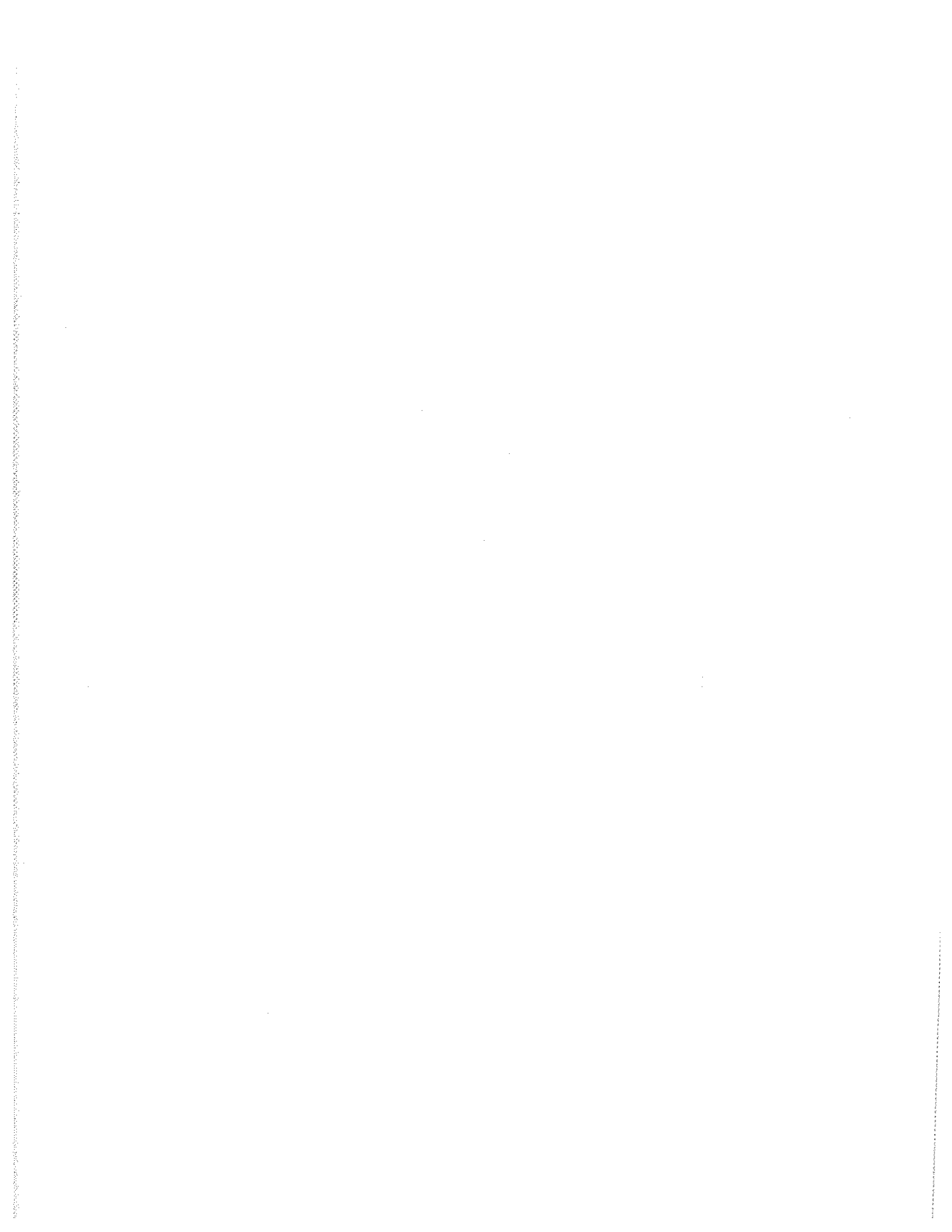
| Analyte  | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers          |
|--|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|--------------------------|
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water) - cont. |           |         |           |                 |               |                 |                |               |                          |
| Reporting Units: mg/l                                      |           |         |           |                 |               |                 |                |               |                          |
| Ammonia-N (Distilled)                                      | EPA 350.2 | 5D14077 | 0.30      | 0.50            | ND            | 1               | 04/14/05       | 04/14/05      | * <u>REV</u> <u>QUAL</u> |
| Biochemical Oxygen Demand                                  | EPA 405.1 | 5D14060 | 0.59      | 2.0             | 3.8           | 1               | 04/14/05       | 04/19/05      |                          |
| Oil & Grease   | EPA 413.1 | 5D14081 | 0.94      | 5.0             | 4.2           | 1               | 04/14/05       | 04/14/05      |                          |
| Total Dissolved Solids                                     | SM2540C   | 5D18095 | 10        | 10              | 240           | 1               | 04/18/05       | 04/18/05      |                          |
| Total Suspended Solids                                     | EPA 160.2 | 5D18089 | 10        | 10              | 12            | 1               | 04/18/05       | 04/18/05      |                          |
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water)         |           |         |           |                 |               |                 |                |               |                          |
| Reporting Units: ml/hr                                     |           |         |           |                 |               |                 |                |               |                          |
| Total Settleable Solids                                    | EPA 160.5 | 5D15069 | 0.10      | 0.10            | ND            | 1               | 04/15/05       | 04/15/05      |                          |
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water)         |           |         |           |                 |               |                 |                |               |                          |
| Reporting Units: NTU                                       |           |         |           |                 |               |                 |                |               |                          |
| Turbidity  | EPA 180.1 | 5D15076 | 0.040     | 1.0             | 25            | 1               | 04/15/05       | 04/15/05      |                          |
| Sample ID: IOD0948-01 (DRAFT: Outfall 012 - Water)         |           |         |           |                 |               |                 |                |               |                          |
| Reporting Units: ug/l                                      |           |         |           |                 |               |                 |                |               |                          |
| Perchlorate  | EPA 314.0 | 5D14043 | 0.80      | 4.0             | ND            | 1               | 04/14/05       | 04/14/05      | u                        |

**AMEC VALIDATED**  
**LEVEL IV**

\* Analytic Not Valid

DRAFT REPORT  
 DRAFT REPORT  
 DATA SUBJECT TO CHANGE

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LABORATORY REPORT

Prepared For: MWH-Pasadena/Boeing  
300 North Lake Avenue, Suite 1200  
Pasadena, CA 91101  
Attention: Bronwyn Kelly

Project: Alfa Outfall 012 - During Test

Sampled: 03/30/05  
Received: 03/30/05  
Issued: 04/13/05 09:21

NELAP #01108CA California ELAP#1197 CSDLAC #10117

*The results listed within this Laboratory Report pertain only to the samples tested in the laboratory. The analyses contained in this report were performed in accordance with the applicable certifications as noted. All soil samples are reported on a wet weight basis unless otherwise noted in the report. This Laboratory Report is confidential and is intended for the sole use of Del Mar Analytical and its client. This report shall not be reproduced, except in full, without written permission from Del Mar Analytical. The Chain(s) of Custody, 2 pages, are included and are an integral part of this report. This entire report was reviewed and approved for release.*

SAMPLE CROSS REFERENCE

SUBCONTRACTED: Refer to the last page for specific subcontract laboratory information included in this report.

| LABORATORY ID | CLIENT ID   | MATRIX |
|---------------|-------------|--------|
| IOC2360-01    | Outfall 012 | Water  |
| IOC2360-02    | Trip Blank  | Water  |

Reviewed By:

Del Mar Analytical, Irvine  
Michele Harper  
Project Manager



LABORATORY REPORT

Prepared For: MWH-Pasadena/Boeing  
300 North Lake Avenue, Suite 1200  
Pasadena, CA 91101  
Attention: Bronwyn Kelly

Project: Alfa Outfall 012 - During Test

Sampled: 04/28/05  
Received: 04/28/05  
Issued: 06/22/05 09:53

NELAP #01108CA California ELAP#1197 CSDLAC #10117

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*This entire report was reviewed and approved for release.*

SAMPLE CROSS REFERENCE

SUBCONTRACTED: Refer to the last page for specific subcontract laboratory information included in this report.

| LABORATORY ID | CLIENT ID   | MATRIX |
|---------------|-------------|--------|
| IOD2047-01    | Outfall 012 | Water  |
| IOD2047-02    | Trip Blank  | Water  |

Reviewed By:

Del Mar Analytical, Irvine  
Michele Harper  
Project Manager



MWH-Pasadena/Boeing  
300 North Lake Avenue, Suite 1200  
Pasadena, CA 91101  
Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

**TOTAL RECOVERABLE PETROLEUM HYDROCARBONS (EPA 418.1)**

| Analyte                                     | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|---|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| Sample ID: IOD2047-01 (Outfall 012 - Water) |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: mg/l                       |           |         |           |                 |               |                 |                |               |                 |
| Total Recoverable Hydrocarbons              | EPA 418.1 | 5D30026 | 0.31      | 1.0             | 5.6           | 1               | 04/30/05       | 04/30/05      |                 |

Del Mar Analytical, Irvine  
Michele Harper  
Project Manager





MWH-Pasadena/Boeing  
300 North Lake Avenue, Suite 1200  
Pasadena, CA 91101  
Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

**EXTRACTABLE FUEL HYDROCARBONS (CADHS/8015 Modified)**

| Analyte   | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|---|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| Sample ID: IOD2047-01 (Outfall 012 - Water) - cont. |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: mg/l                               |           |         |           |                 |               |                 |                |               |                 |
| EFH (C13 - C22)                                     | EPA 8015B | 5E03053 | 0.082     | 0.50            | 1.2           | 0.962           | 05/03/05       | 05/03/05      |                 |
| Surrogate: n-Octacosane (40-125%)                   |           |         |           |                 | 74 %          |                 |                |               |                 |

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Received: 04/28/05

**VOLATILE FUEL HYDROCARBONS (EPA 5030/CADHS Mod. 8015)**

| Analyte  | Method        | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|--|---------------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| <b>Sample ID: IOD2047-01 (Outfall 012 - Water) - cont.</b> |               |         |           |                 |               |                 |                |               |                 |
| Reporting Units: mg/l                                      |               |         |           |                 |               |                 |                |               |                 |
| GRO (C4 - C12)   | EPA 8015 Mod. | 5E04040 | 0.50      | 1.0             | 2.7           | 10              | 05/04/05       | 05/04/05      |                 |
| Surrogate: 4-BFB (FID) (65-140%)                           |               |         |           |                 | 100 %         |                 |                |               |                 |
| <b>Sample ID: IOD2047-02 (Trip Blank - Water)</b>          |               |         |           |                 |               |                 |                |               |                 |
| Reporting Units: mg/l                                      |               |         |           |                 |               |                 |                |               |                 |
| GRO (C4 - C12)   | EPA 8015 Mod. | 5E04040 | 0.050     | 0.10            | ND            | 1               | 05/04/05       | 05/04/05      |                 |
| Surrogate: 4-BFB (FID) (65-140%)                           |               |         |           |                 | 89 %          |                 |                |               |                 |

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Sampled: 04/28/05  
Received: 04/28/05

PURGEABLES BY GC/MS (EPA 624)

| Analyte  | Method  | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|--|---------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| <b>Sample ID: IOD2047-01 (Outfall 012 - Water)</b> |         |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                              |         |         |           |                 |               |                 |                |               |                 |
| 1,2-Dibromoethane (EDB)                            | EPA 624 | 5E05024 | 0.32      | 2.0             | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| Methyl-tert-butyl Ether (MTBE)                     | EPA 624 | 5E05024 | 0.32      | 5.0             | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| 1,2,3-Trichloropropane                             | EPA 624 | 5E05024 | 0.85      | 10              | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| Di-isopropyl Ether (DIPE)                          | EPA 624 | 5E05024 | 0.25      | 5.0             | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| tert-Butanol (TBA)                                 | EPA 624 | 5E05024 | 3.1       | 25              | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| Surrogate: Dibromofluoromethane (80-120%)          |         |         |           |                 |               |                 |                |               | 114 %           |
| Surrogate: Toluene-d8 (80-120%)                    |         |         |           |                 |               |                 |                |               | 112 %           |
| Surrogate: 4-Bromofluorobenzene (80-120%)          |         |         |           |                 |               |                 |                |               | 110 %           |
| <b>Sample ID: IOD2047-02 (Trip Blank - Water)</b>  |         |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                              |         |         |           |                 |               |                 |                |               |                 |
| 1,2-Dibromoethane (EDB)                            | EPA 624 | 5E05024 | 0.32      | 2.0             | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| Methyl-tert-butyl Ether (MTBE)                     | EPA 624 | 5E05024 | 0.32      | 5.0             | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| 1,2,3-Trichloropropane                             | EPA 624 | 5E05024 | 0.85      | 10              | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| Di-isopropyl Ether (DIPE)                          | EPA 624 | 5E05024 | 0.25      | 5.0             | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| tert-Butanol (TBA)                                 | EPA 624 | 5E05024 | 3.1       | 25              | ND            | 1               | 05/05/05       | 05/05/05      |                 |
| Surrogate: Dibromofluoromethane (80-120%)          |         |         |           |                 |               |                 |                |               | 111 %           |
| Surrogate: Toluene-d8 (80-120%)                    |         |         |           |                 |               |                 |                |               | 112 %           |
| Surrogate: 4-Bromofluorobenzene (80-120%)          |         |         |           |                 |               |                 |                |               | 107 %           |

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**ACID & BASE/NEUTRALS BY GC/MS (EPA 625)**

| Analyte  | Method  | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|--|---------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| <b>Sample ID: IOD2047-01 (Outfall 012 - Water)</b> |         |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                              |         |         |           |                 |               |                 |                |               |                 |
| Naphthalene  | EPA 625 | 5E01019 | 4.5       | 10              | 24            | 0.99            | 05/01/05       | 05/04/05      |                 |
| N-Nitrosodimethylamine                             | EPA 625 | 5E01019 | 3.7       | 20              | ND            | 0.99            | 05/01/05       | 05/04/05      |                 |
| Surrogate: 2-Fluorophenol (30-120%)                |         |         |           |                 | 53 %          |                 |                |               |                 |
| Surrogate: Phenol-d6 (35-120%)                     |         |         |           |                 | 68 %          |                 |                |               |                 |
| Surrogate: 2,4,6-Tribromophenol (45-120%)          |         |         |           |                 | 69 %          |                 |                |               |                 |
| Surrogate: Nitrobenzene-d5 (45-120%)               |         |         |           |                 | 85 %          |                 |                |               |                 |
| Surrogate: 2-Fluorobiphenyl (45-120%)              |         |         |           |                 | 74 %          |                 |                |               |                 |
| Surrogate: Terphenyl-d14 (45-120%)                 |         |         |           |                 | 132 %         |                 |                |               | ZX              |

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Sampled: 04/28/05  
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INORGANICS

| Analyte  | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|--|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| <b>Sample ID: IOD2047-01 (Outfall 012 - Water) - cont.</b> |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: mg/l                                      |           |         |           |                 |               |                 |                |               |                 |
| Ammonia-N (Distilled)                                      | EPA 350.2 | 5E02067 | 0.30      | 0.50            | ND            | 1               | 05/02/05       | 05/02/05      |                 |
| Biochemical Oxygen Demand                                  | EPA 405.1 | 5D29091 | 0.59      | 2.0             | 3.2           | 1               | 04/29/05       | 05/04/05      |                 |
| Oil & Grease   | EPA 413.1 | 5E04036 | 0.94      | 5.0             | ND            | 1               | 05/04/05       | 05/04/05      |                 |
| Total Dissolved Solids                                     | SM2540C   | 5D29129 | 10        | 10              | 250           | 1               | 04/29/05       | 04/29/05      |                 |
| Total Suspended Solids                                     | EPA 160.2 | 5E04071 | 10        | 10              | 21            | 1               | 05/04/05       | 05/04/05      |                 |
| <b>Sample ID: IOD2047-01 (Outfall 012 - Water)</b>         |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ml/l/hr                                   |           |         |           |                 |               |                 |                |               |                 |
| Total Settleable Solids                                    | EPA 160.5 | 5D29094 | 0.10      | 0.10            | 0.10          | 1               | 04/29/05       | 04/29/05      |                 |
| <b>Sample ID: IOD2047-01 (Outfall 012 - Water)</b>         |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: NTU                                       |           |         |           |                 |               |                 |                |               |                 |
| Turbidity  | EPA 180.1 | 5D29110 | 0.040     | 1.0             | 23            | 1               | 04/29/05       | 04/29/05      |                 |
| <b>Sample ID: IOD2047-01 (Outfall 012 - Water)</b>         |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                                      |           |         |           |                 |               |                 |                |               |                 |
| Perchlorate  | EPA 314.0 | 5D29065 | 0.80      | 4.0             | ND            | 1               | 04/29/05       | 04/30/05      |                 |

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Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

**1,4-DIOXANE BY GC/MS (EPA 5030B/8260B)**

| Analyte  | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|--|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| <b>Sample ID: IOD2047-01 (Outfall 012 - Water) - cont.</b> |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                                      |           |         |           |                 |               |                 |                |               |                 |
| 1,4-Dioxane  | EPA 8260B | P5E0214 | 0.49      | 1.0             | ND            | 1               | 05/02/05       | 05/02/05      |                 |
| Surrogate: Dibromofluoromethane (80-125%)                  |           |         |           |                 | 97 %          |                 |                |               |                 |

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Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

**SHORT HOLD TIME DETAIL REPORT**

| Sample ID: Outfall 012 (IOD2047-01) - Water | Hold Time<br>(in days) | Date/Time<br>Sampled | Date/Time<br>Received | Date/Time<br>Extracted | Date/Time<br>Analyzed |
|---|------------------------|----------------------|-----------------------|------------------------|-----------------------|
| EPA 160.5                                   | 2                      | 04/28/2005 14:35     | 04/28/2005 18:15      | 04/29/2005 13:29       | 04/29/2005 15:00      |
| EPA 180.1                                   | 2                      | 04/28/2005 14:35     | 04/28/2005 18:15      | 04/29/2005 15:00       | 04/29/2005 16:00      |
| EPA 405.1                                   | 2                      | 04/28/2005 14:35     | 04/28/2005 18:15      | 04/29/2005 13:10       | 05/04/2005 10:00      |

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Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

**METHOD BLANK/QC DATA**

**TOTAL RECOVERABLE PETROLEUM HYDROCARBONS (EPA 418.1)**

| Analyte  | Result | Reporting Limit | MDL  | Units | Spike Level | Source Result | %REC %REC | RPD Limits | RPD RPD | RPD Limit | Data Qualifiers |
|--|--------|-----------------|------|-------|-------------|---------------|-----------|------------|---------|-----------|-----------------|
| <b>Batch: 5D30026 Extracted: 04/30/05</b>          |        |                 |      |       |             |               |           |            |         |           |                 |
| <b>Blank Analyzed: 04/30/2005 (5D30026-BLK1)</b>   |        |                 |      |       |             |               |           |            |         |           |                 |
| Total Recoverable Hydrocarbons                     | ND     | 1.0             | 0.31 | mg/l  |             |               |           |            |         |           |                 |
| <b>LCS Analyzed: 04/30/2005 (5D30026-BS1)</b>      |        |                 |      |       |             |               |           |            |         |           |                 |
| Total Recoverable Hydrocarbons                     | 4.84   | 1.0             | 0.31 | mg/l  | 5.00        |               | 97        | 65-120     |         |           | M-NRI           |
| <b>LCS Dup Analyzed: 04/30/2005 (5D30026-BSD1)</b> |        |                 |      |       |             |               |           |            |         |           |                 |
| Total Recoverable Hydrocarbons                     | 4.72   | 1.0             | 0.31 | mg/l  | 5.00        |               | 94        | 65-120     | 3       | 20        |                 |

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**METHOD BLANK/QC DATA**

**EXTRACTABLE FUEL HYDROCARBONS (CADHS/8015 Modified)**

| Analyte  | Result | Reporting Limit | MDL   | Units | Spike Level | Source Result | %REC %REC | RPD Limits | RPD RPD | Data Qualifiers |
|--|--------|-----------------|-------|-------|-------------|---------------|-----------|------------|---------|-----------------|
| <b>Batch: 5E03053 Extracted: 05/03/05</b>          |        |                 |       |       |             |               |           |            |         |                 |
| <b>Blank Analyzed: 05/03/2005 (5E03053-BLK1)</b>   |        |                 |       |       |             |               |           |            |         |                 |
| EFH (C13 - C22)                                    | ND     | 0.50            | 0.082 | mg/l  |             |               |           |            |         |                 |
| EFH (C13 - C40)                                    | ND     | 0.50            | 0.082 | mg/l  |             |               |           |            |         |                 |
| Surrogate: n-Octacosane                            | 0.161  |                 |       | mg/l  | 0.200       |               | 80        | 40-125     |         |                 |
| <b>LCS Analyzed: 05/03/2005 (5E03053-BS1)</b>      |        |                 |       |       |             |               |           |            |         |                 |
| EFH (C13 - C40)                                    | 0.509  | 0.50            | 0.082 | mg/l  | 0.775       |               | 66        | 40-120     |         | <b>M-NR1</b>    |
| Surrogate: n-Octacosane                            | 0.150  |                 |       | mg/l  | 0.200       |               | 75        | 40-125     |         |                 |
| <b>LCS Dup Analyzed: 05/03/2005 (5E03053-BSD1)</b> |        |                 |       |       |             |               |           |            |         |                 |
| EFH (C13 - C40)                                    | 0.503  | 0.50            | 0.082 | mg/l  | 0.775       |               | 65        | 40-120     | 1       | 25              |
| Surrogate: n-Octacosane                            | 0.150  |                 |       | mg/l  | 0.200       |               | 75        | 40-125     |         |                 |

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**METHOD BLANK/QC DATA**

**VOLATILE FUEL HYDROCARBONS (EPA 5030/CADHS Mod. 8015)**

| Analyte  | Result  | Reporting Limit | MDL   | Units | Spike Level | Source Result | %REC %REC | RPD Limits | RPD | RPD Limit | Data Qualifiers |
|--|---------|-----------------|-------|-------|-------------|---------------|-----------|------------|-----|-----------|-----------------|
| <b>Batch: 5E04040 Extracted: 05/04/05</b>                                      |         |                 |       |       |             |               |           |            |     |           |                 |
| <b>Blank Analyzed: 05/04/2005 (5E04040-BLK1)</b>                               |         |                 |       |       |             |               |           |            |     |           |                 |
| GRO (C4 - C12)   | ND      | 0.10            | 0.050 | mg/l  |             |               |           |            |     |           |                 |
| Surrogate: 4-BFB (FID)   | 0.00985 |                 |       | mg/l  | 0.0100      |               | 98        | 65-140     |     |           |                 |
| <b>LCS Analyzed: 05/04/2005 (5E04040-BS1)</b>                                  |         |                 |       |       |             |               |           |            |     |           |                 |
| GRO (C4 - C12)   | 0.878   | 0.10            | 0.050 | mg/l  | 0.800       |               | 110       | 70-140     |     |           |                 |
| Surrogate: 4-BFB (FID)   | 0.0309  |                 |       | mg/l  | 0.0300      |               | 103       | 65-140     |     |           |                 |
| <b>Matrix Spike Analyzed: 05/04/2005 (5E04040-MS1) Source: IOD1878-17</b>      |         |                 |       |       |             |               |           |            |     |           |                 |
| GRO (C4 - C12)   | 7.13    | 1.0             | 0.50  | mg/l  | 2.20        | 9.2           | -94       | 60-140     |     |           | M2              |
| Surrogate: 4-BFB (FID)   | 0.100   |                 |       | mg/l  | 0.100       |               | 100       | 65-140     |     |           |                 |
| <b>Matrix Spike Dup Analyzed: 05/04/2005 (5E04040-MSD1) Source: IOD1878-17</b> |         |                 |       |       |             |               |           |            |     |           |                 |
| GRO (C4 - C12)   | 7.32    | 1.0             | 0.50  | mg/l  | 2.20        | 9.2           | -85       | 60-140     | 3   | 20        | M2              |
| Surrogate: 4-BFB (FID)   | 0.102   |                 |       | mg/l  | 0.100       |               | 102       | 65-140     |     |           |                 |

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## METHOD BLANK/QC DATA

### PURGEABLES BY GC/MS (EPA 624)

| Analyte  | Result | Reporting Limit | MDL  | Units | Spike Level | Source Result | %REC %REC | RPD RPD | RPD Limit | Data Qualifiers |
|--|--------|-----------------|------|-------|-------------|---------------|-----------|---------|-----------|-----------------|
| <b>Batch: 5E05024 Extracted: 05/05/05</b>              |        |                 |      |       |             |               |           |         |           |                 |
| <b>Blank Analyzed: 05/05/2005 (5E05024-BLK1)</b>       |        |                 |      |       |             |               |           |         |           |                 |
| 1,2-Dibromoethane (EDB)                                | ND     | 2.0             | 0.32 | ug/l  |             |               |           |         |           |                 |
| Methyl-tert-butyl Ether (MTBE)                         | ND     | 5.0             | 0.32 | ug/l  |             |               |           |         |           |                 |
| 1,2,3-Trichloropropane                                 | ND     | 10              | 0.85 | ug/l  |             |               |           |         |           |                 |
| Di-isopropyl Ether (DIPE)                              | ND     | 5.0             | 0.25 | ug/l  |             |               |           |         |           |                 |
| tert-Butanol (TBA)                                     | ND     | 25              | 3.1  | ug/l  |             |               |           |         |           |                 |
| Surrogate: Dibromofluoromethane                        | 28.3   |                 |      | ug/l  | 25.0        |               | 113       |         | 80-120    |                 |
| Surrogate: Toluene-d8                                  | 27.9   |                 |      | ug/l  | 25.0        |               | 112       |         | 80-120    |                 |
| Surrogate: 4-Bromofluorobenzene                        | 26.2   |                 |      | ug/l  | 25.0        |               | 105       |         | 80-120    |                 |
| <b>LCS Analyzed: 05/05/2005 (5E05024-BS1)</b>          |        |                 |      |       |             |               |           |         |           |                 |
| 1,2-Dibromoethane (EDB)                                | 22.4   | 2.0             | 0.32 | ug/l  | 25.0        |               | 90        |         | 70-125    |                 |
| Methyl-tert-butyl Ether (MTBE)                         | 21.5   | 5.0             | 0.32 | ug/l  | 25.0        |               | 86        |         | 55-140    |                 |
| 1,2,3-Trichloropropane                                 | 21.1   | 10              | 0.85 | ug/l  | 25.0        |               | 84        |         | 55-130    |                 |
| Di-isopropyl Ether (DIPE)                              | 21.2   | 5.0             | 0.25 | ug/l  | 25.0        |               | 85        |         | 60-135    |                 |
| tert-Butanol (TBA)                                     | 127    | 25              | 3.1  | ug/l  | 125         |               | 102       |         | 65-135    |                 |
| Surrogate: Dibromofluoromethane                        | 28.9   |                 |      | ug/l  | 25.0        |               | 116       |         | 80-120    |                 |
| Surrogate: Toluene-d8                                  | 28.0   |                 |      | ug/l  | 25.0        |               | 112       |         | 80-120    |                 |
| Surrogate: 4-Bromofluorobenzene                        | 27.9   |                 |      | ug/l  | 25.0        |               | 112       |         | 80-120    |                 |
| <b>Matrix Spike Analyzed: 05/05/2005 (5E05024-MS1)</b> |        |                 |      |       |             |               |           |         |           |                 |
| <b>Source: IOD2064-02</b>                              |        |                 |      |       |             |               |           |         |           |                 |
| 1,2-Dibromoethane (EDB)                                | 20.6   | 2.0             | 0.32 | ug/l  | 25.0        | ND            | 82        |         | 65-130    |                 |
| Methyl-tert-butyl Ether (MTBE)                         | 20.8   | 5.0             | 0.32 | ug/l  | 25.0        | ND            | 83        |         | 50-150    |                 |
| 1,2,3-Trichloropropane                                 | 19.3   | 10              | 0.85 | ug/l  | 25.0        | ND            | 77        |         | 50-135    |                 |
| Di-isopropyl Ether (DIPE)                              | 20.4   | 5.0             | 0.25 | ug/l  | 25.0        | ND            | 82        |         | 60-140    |                 |
| tert-Butanol (TBA)                                     | 125    | 25              | 3.1  | ug/l  | 125         | ND            | 100       |         | 60-145    |                 |
| Surrogate: Dibromofluoromethane                        | 29.6   |                 |      | ug/l  | 25.0        |               | 118       |         | 80-120    |                 |
| Surrogate: Toluene-d8                                  | 28.8   |                 |      | ug/l  | 25.0        |               | 115       |         | 80-120    |                 |
| Surrogate: 4-Bromofluorobenzene                        | 27.6   |                 |      | ug/l  | 25.0        |               | 110       |         | 80-120    |                 |

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Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

**METHOD BLANK/QC DATA**

**PURGEABLES BY GC/MS (EPA 624)**

| Analyte   | Result | Reporting Limit | MDL  | Units | Spike Level | Source Result             | %REC | %REC Limits | RPD | RPD Limit | Data Qualifiers |
|---|--------|-----------------|------|-------|-------------|---------------------------|------|-------------|-----|-----------|-----------------|
| <b>Batch: 5E05024 Extracted: 05/05/05</b>                   |        |                 |      |       |             |                           |      |             |     |           |                 |
| <b>Matrix Spike Dup Analyzed: 05/05/2005 (5E05024-MSD1)</b> |        |                 |      |       |             | <b>Source: IOD2064-02</b> |      |             |     |           |                 |
| 1,2-Dibromoethane (EDB)                                     | 21.3   | 2.0             | 0.32 | ug/l  | 25.0        | ND                        | 85   | 65-130      | 3   | 25        |                 |
| Methyl-tert-butyl Ether (MTBE)                              | 21.2   | 5.0             | 0.32 | ug/l  | 25.0        | ND                        | 85   | 50-150      | 2   | 25        |                 |
| 1,2,3-Trichloropropane                                      | 20.4   | 10              | 0.85 | ug/l  | 25.0        | ND                        | 82   | 50-135      | 6   | 30        |                 |
| Di-isopropyl Ether (DIPE)                                   | 20.4   | 5.0             | 0.25 | ug/l  | 25.0        | ND                        | 82   | 60-140      | 0   | 25        |                 |
| tert-Butanol (TBA)  | 119    | 25              | 3.1  | ug/l  | 125         | ND                        | 95   | 60-145      | 5   | 25        |                 |
| Surrogate: Dibromofluoromethane                             | 29.3   |                 |      | ug/l  | 25.0        |                           | 117  | 80-120      |     |           |                 |
| Surrogate: Toluene-d8                                       | 28.7   |                 |      | ug/l  | 25.0        |                           | 115  | 80-120      |     |           |                 |
| Surrogate: 4-Bromofluorobenzene                             | 27.7   |                 |      | ug/l  | 25.0        |                           | 111  | 80-120      |     |           |                 |

Del Mar Analytical, Irvine  
Michele Harper  
Project Manager



MWH-Pasadena/Boeing  
300 North Lake Avenue, Suite 1200  
Pasadena, CA 91101  
Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

METHOD BLANK/QC DATA

ACID & BASE/NEUTRALS BY GC/MS (EPA 625)

| Analyte  | Result | Reporting Limit | MDL | Units | Spike Level | Source Result | %REC %REC | RPD Limits | RPD RPD | Data Limit | Qualifiers |
|--|--------|-----------------|-----|-------|-------------|---------------|-----------|------------|---------|------------|------------|
| <b>Batch: 5E01019 Extracted: 05/01/05</b>          |        |                 |     |       |             |               |           |            |         |            |            |
| <b>Blank Analyzed: 05/03/2005 (5E01019-BLK1)</b>   |        |                 |     |       |             |               |           |            |         |            |            |
| Naphthalene  | ND     | 10              | 4.5 | ug/l  |             |               |           |            |         |            |            |
| N-Nitrosodimethylamine                             | ND     | 20              | 3.7 | ug/l  |             |               |           |            |         |            |            |
| Surrogate: 2-Fluorophenol                          | 87.4   |                 |     | ug/l  | 200         |               | 44        | 30-120     |         |            |            |
| Surrogate: Phenol-d6                               | 125    |                 |     | ug/l  | 200         |               | 62        | 35-120     |         |            |            |
| Surrogate: 2,4,6-Tribromophenol                    | 132    |                 |     | ug/l  | 200         |               | 66        | 45-120     |         |            |            |
| Surrogate: Nitrobenzene-d5                         | 71.3   |                 |     | ug/l  | 100         |               | 71        | 45-120     |         |            |            |
| Surrogate: 2-Fluorobiphenyl                        | 73.8   |                 |     | ug/l  | 100         |               | 74        | 45-120     |         |            |            |
| Surrogate: Terphenyl-d14                           | 79.3   |                 |     | ug/l  | 100         |               | 79        | 45-120     |         |            |            |
| <b>LCS Analyzed: 05/03/2005 (5E01019-BS1)</b>      |        |                 |     |       |             |               |           |            |         |            |            |
| Naphthalene  | 66.5   | 10              | 4.5 | ug/l  | 100         |               | 66        | 50-120     |         |            | M-NRI      |
| N-Nitrosodimethylamine                             | 56.6   | 20              | 3.7 | ug/l  | 100         |               | 57        | 40-120     |         |            |            |
| Surrogate: 2-Fluorophenol                          | 93.0   |                 |     | ug/l  | 200         |               | 46        | 30-120     |         |            |            |
| Surrogate: Phenol-d6                               | 117    |                 |     | ug/l  | 200         |               | 58        | 35-120     |         |            |            |
| Surrogate: 2,4,6-Tribromophenol                    | 145    |                 |     | ug/l  | 200         |               | 72        | 45-120     |         |            |            |
| Surrogate: Nitrobenzene-d5                         | 68.6   |                 |     | ug/l  | 100         |               | 69        | 45-120     |         |            |            |
| Surrogate: 2-Fluorobiphenyl                        | 69.1   |                 |     | ug/l  | 100         |               | 69        | 45-120     |         |            |            |
| Surrogate: Terphenyl-d14                           | 82.7   |                 |     | ug/l  | 100         |               | 83        | 45-120     |         |            |            |
| <b>LCS Dup Analyzed: 05/03/2005 (5E01019-BSD1)</b> |        |                 |     |       |             |               |           |            |         |            |            |
| Naphthalene  | 72.4   | 10              | 4.5 | ug/l  | 100         |               | 72        | 50-120     | 8       | 20         |            |
| N-Nitrosodimethylamine                             | 60.9   | 20              | 3.7 | ug/l  | 100         |               | 61        | 40-120     | 7       | 20         |            |
| Surrogate: 2-Fluorophenol                          | 102    |                 |     | ug/l  | 200         |               | 51        | 30-120     |         |            |            |
| Surrogate: Phenol-d6                               | 127    |                 |     | ug/l  | 200         |               | 64        | 35-120     |         |            |            |
| Surrogate: 2,4,6-Tribromophenol                    | 148    |                 |     | ug/l  | 200         |               | 74        | 45-120     |         |            |            |
| Surrogate: Nitrobenzene-d5                         | 76.2   |                 |     | ug/l  | 100         |               | 76        | 45-120     |         |            |            |
| Surrogate: 2-Fluorobiphenyl                        | 73.5   |                 |     | ug/l  | 100         |               | 74        | 45-120     |         |            |            |
| Surrogate: Terphenyl-d14                           | 85.4   |                 |     | ug/l  | 100         |               | 85        | 45-120     |         |            |            |

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Project Manager



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MWH-Pasadena/Boeing  
 300 North Lake Avenue, Suite 1200  
 Pasadena, CA 91101  
 Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD2047

Sampled: 04/28/05  
 Received: 04/28/05

## METHOD BLANK/QC DATA

### INORGANICS

| Analyte   | Result | Reporting Limit | MDL   | Units | Spike Level | Source Result | %REC %REC | Limit  | RPD | RPD Limit | Data Qualifiers |
|---|--------|-----------------|-------|-------|-------------|---------------|-----------|--------|-----|-----------|-----------------|
| <b><u>Batch: 5D29065 Extracted: 04/29/05</u></b>            |        |                 |       |       |             |               |           |        |     |           |                 |
| <b>Blank Analyzed: 04/29/2005 (5D29065-BLK1)</b>            |        |                 |       |       |             |               |           |        |     |           |                 |
| Perchlorate   | ND     | 4.0             | 0.80  | ug/l  |             |               |           |        |     |           |                 |
| <b>LCS Analyzed: 04/29/2005 (5D29065-BS1)</b>               |        |                 |       |       |             |               |           |        |     |           |                 |
| Perchlorate   | 51.0   | 4.0             | 0.80  | ug/l  | 50.0        |               | 102       | 85-115 |     |           |                 |
| <b>Matrix Spike Analyzed: 04/29/2005 (5D29065-MS1)</b>      |        |                 |       |       |             |               |           |        |     |           |                 |
| Perchlorate   | 53.1   | 4.0             | 0.80  | ug/l  | 50.0        | ND            | 106       | 80-120 |     |           |                 |
| <b>Matrix Spike Dup Analyzed: 04/29/2005 (5D29065-MSD1)</b> |        |                 |       |       |             |               |           |        |     |           |                 |
| Perchlorate   | 52.9   | 4.0             | 0.80  | ug/l  | 50.0        | ND            | 106       | 80-120 | 0   | 20        |                 |
| <b><u>Batch: 5D29091 Extracted: 04/29/05</u></b>            |        |                 |       |       |             |               |           |        |     |           |                 |
| <b>Blank Analyzed: 05/04/2005 (5D29091-BLK1)</b>            |        |                 |       |       |             |               |           |        |     |           |                 |
| Biochemical Oxygen Demand                                   | ND     | 2.0             | 0.59  | mg/l  |             |               |           |        |     |           |                 |
| <b>LCS Analyzed: 05/04/2005 (5D29091-BS1)</b>               |        |                 |       |       |             |               |           |        |     |           |                 |
| Biochemical Oxygen Demand                                   | 209    | 100             | 30    | mg/l  | 198         |               | 106       | 85-115 |     |           |                 |
| <b>LCS Dup Analyzed: 05/04/2005 (5D29091-BSD1)</b>          |        |                 |       |       |             |               |           |        |     |           |                 |
| Biochemical Oxygen Demand                                   | 208    | 100             | 30    | mg/l  | 198         |               | 105       | 85-115 | 1   | 20        |                 |
| <b><u>Batch: 5D29110 Extracted: 04/29/05</u></b>            |        |                 |       |       |             |               |           |        |     |           |                 |
| <b>Blank Analyzed: 04/29/2005 (5D29110-BLK1)</b>            |        |                 |       |       |             |               |           |        |     |           |                 |
| Turbidity   | ND     | 1.0             | 0.040 | NTU   |             |               |           |        |     |           |                 |

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Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

METHOD BLANK/QC DATA

INORGANICS

| Analyte   | Result | Reporting Limit | MDL  | Units | Spike Level | Source Result | %REC %REC | RPD Limits | RPD | RPD Limit | Data Qualifiers |
|---|--------|-----------------|------|-------|-------------|---------------|-----------|------------|-----|-----------|-----------------|
| <b>Batch: 5D29110 Extracted: 04/29/05</b>                   |        |                 |      |       |             |               |           |            |     |           |                 |
| <b>Duplicate Analyzed: 04/29/2005 (5D29110-DUP1)</b>        |        |                 |      |       |             |               |           |            |     |           |                 |
| Turbidity   | 135    | 5.0             | 0.20 | NTU   |             | 130           |           |            | 4   | 20        |                 |
| <b>Batch: 5D29129 Extracted: 04/29/05</b>                   |        |                 |      |       |             |               |           |            |     |           |                 |
| <b>Blank Analyzed: 04/29/2005 (5D29129-BLK1)</b>            |        |                 |      |       |             |               |           |            |     |           |                 |
| Total Dissolved Solids                                      | ND     | 10              | 10   | mg/l  |             |               |           |            |     |           |                 |
| <b>LCS Analyzed: 04/29/2005 (5D29129-BS1)</b>               |        |                 |      |       |             |               |           |            |     |           |                 |
| Total Dissolved Solids                                      | 930    | 10              | 10   | mg/l  | 1000        |               | 93        | 90-110     |     |           |                 |
| <b>Duplicate Analyzed: 04/29/2005 (5D29129-DUP1)</b>        |        |                 |      |       |             |               |           |            |     |           |                 |
| Total Dissolved Solids                                      | 334    | 10              | 10   | mg/l  |             | 360           |           |            | 7   | 10        |                 |
| <b>Batch: 5E02067 Extracted: 05/02/05</b>                   |        |                 |      |       |             |               |           |            |     |           |                 |
| <b>Blank Analyzed: 05/02/2005 (5E02067-BLK1)</b>            |        |                 |      |       |             |               |           |            |     |           |                 |
| Ammonia-N (Distilled)                                       | ND     | 0.50            | 0.30 | mg/l  |             |               |           |            |     |           |                 |
| <b>LCS Analyzed: 05/02/2005 (5E02067-BS1)</b>               |        |                 |      |       |             |               |           |            |     |           |                 |
| Ammonia-N (Distilled)                                       | 10.4   | 0.50            | 0.30 | mg/l  | 10.0        |               | 104       | 80-115     |     |           |                 |
| <b>Matrix Spike Analyzed: 05/02/2005 (5E02067-MS1)</b>      |        |                 |      |       |             |               |           |            |     |           |                 |
| Ammonia-N (Distilled)                                       | 10.9   | 0.50            | 0.30 | mg/l  | 10.0        | 1.1           | 98        | 70-120     |     |           |                 |
| <b>Matrix Spike Dup Analyzed: 05/02/2005 (5E02067-MSD1)</b> |        |                 |      |       |             |               |           |            |     |           |                 |
| Ammonia-N (Distilled)                                       | 11.2   | 0.50            | 0.30 | mg/l  | 10.0        | 1.1           | 101       | 70-120     | 3   | 15        |                 |

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**METHOD BLANK/QC DATA**

**INORGANICS**

| Analyte  | Result | Reporting Limit | MDL  | Units | Spike Level | Source Result      | %REC %REC | Limit  | RPD | RPD Limit | Data Qualifiers |
|--|--------|-----------------|------|-------|-------------|--------------------|-----------|--------|-----|-----------|-----------------|
| <b>Batch: 5E04036 Extracted: 05/04/05</b>            |        |                 |      |       |             |                    |           |        |     |           |                 |
| <b>Blank Analyzed: 05/04/2005 (5E04036-BLK1)</b>     |        |                 |      |       |             |                    |           |        |     |           |                 |
| Oil & Grease   | ND     | 5.0             | 0.94 | mg/l  |             |                    |           |        |     |           |                 |
| <b>LCS Analyzed: 05/04/2005 (5E04036-BS1)</b>        |        |                 |      |       |             |                    |           |        |     |           |                 |
| Oil & Grease   | 18.5   | 5.0             | 0.94 | mg/l  | 20.0        |                    | 92        | 65-120 |     |           | M-NR1           |
| <b>LCS Dup Analyzed: 05/04/2005 (5E04036-BSD1)</b>   |        |                 |      |       |             |                    |           |        |     |           |                 |
| Oil & Grease   | 18.9   | 5.0             | 0.94 | mg/l  | 20.0        |                    | 94        | 65-120 | 2   | 20        |                 |
| <b>Batch: 5E04071 Extracted: 05/04/05</b>            |        |                 |      |       |             |                    |           |        |     |           |                 |
| <b>Blank Analyzed: 05/04/2005 (5E04071-BLK1)</b>     |        |                 |      |       |             |                    |           |        |     |           |                 |
| Total Suspended Solids                               | ND     | 10              | 10   | mg/l  |             |                    |           |        |     |           |                 |
| <b>LCS Analyzed: 05/04/2005 (5E04071-BS1)</b>        |        |                 |      |       |             |                    |           |        |     |           |                 |
| Total Suspended Solids                               | 1000   | 10              | 10   | mg/l  | 1000        |                    | 100       | 85-115 |     |           |                 |
| <b>Duplicate Analyzed: 05/04/2005 (5E04071-DUP1)</b> |        |                 |      |       |             |                    |           |        |     |           |                 |
| Total Suspended Solids                               | ND     | 10              | 10   | mg/l  |             | Source: IOD2054-01 |           |        |     |           |                 |
|  |        |                 |      |       |             | ND                 |           |        |     | 10        |                 |

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Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

**METHOD BLANK/QC DATA**

**1,4-DIOXANE BY GC/MS (EPA 5030B/8260B)**

| Analyte   | Result | Reporting Limit | MDL  | Units | Spike Level | Source Result                | %REC %REC | Limit  | RPD | RPD Limit | Data Qualifiers |
|---|--------|-----------------|------|-------|-------------|------------------------------|-----------|--------|-----|-----------|-----------------|
| <b>Batch: P5E0214 Extracted: 05/02/05</b>                   |        |                 |      |       |             |                              |           |        |     |           |                 |
| <b>Blank Analyzed: 05/02/2005 (P5E0214-BLK1)</b>            |        |                 |      |       |             |                              |           |        |     |           |                 |
| 1,4-Dioxane   | ND     | 1.0             | 0.49 | ug/l  |             |                              |           |        |     |           |                 |
| Surrogate: Dibromofluoromethane                             | 1.03   |                 |      | ug/l  | 1.00        |                              | 103       | 80-125 |     |           |                 |
| <b>LCS Analyzed: 05/02/2005 (P5E0214-BS1)</b>               |        |                 |      |       |             |                              |           |        |     |           |                 |
| 1,4-Dioxane   | 12.3   | 1.0             | 0.49 | ug/l  | 10.0        |                              | 123       | 70-130 |     |           |                 |
| Surrogate: Dibromofluoromethane                             | 1.00   |                 |      | ug/l  | 1.00        |                              | 100       | 80-125 |     |           |                 |
| <b>LCS Dup Analyzed: 05/02/2005 (P5E0214-BSD1)</b>          |        |                 |      |       |             |                              |           |        |     |           |                 |
| 1,4-Dioxane   | 9.82   | 1.0             | 0.49 | ug/l  | 10.0        |                              | 98        | 70-130 | 22  | 20        | R-7             |
| Surrogate: Dibromofluoromethane                             | 0.970  |                 |      | ug/l  | 1.00        |                              | 97        | 80-125 |     |           |                 |
| <b>Matrix Spike Analyzed: 05/02/2005 (P5E0214-MS1)</b>      |        |                 |      |       |             |                              |           |        |     |           |                 |
|   |        |                 |      |       |             | <b>Source: POD0904-02RE1</b> |           |        |     |           |                 |
| 1,4-Dioxane   | 176    | 1.0             | 0.49 | ug/l  | 10.0        | 140                          | 360       | 70-150 |     |           | E, M-HA         |
| Surrogate: Dibromofluoromethane                             | 0.980  |                 |      | ug/l  | 1.00        |                              | 98        | 80-125 |     |           |                 |
| <b>Matrix Spike Dup Analyzed: 05/02/2005 (P5E0214-MSD1)</b> |        |                 |      |       |             |                              |           |        |     |           |                 |
|   |        |                 |      |       |             | <b>Source: POD0904-02RE1</b> |           |        |     |           |                 |
| 1,4-Dioxane   | 151    | 1.0             | 0.49 | ug/l  | 10.0        | 140                          | 110       | 70-150 | 15  | 25        | E, M-HA         |
| Surrogate: Dibromofluoromethane                             | 0.970  |                 |      | ug/l  | 1.00        |                              | 97        | 80-125 |     |           |                 |

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Michele Harper  
Project Manager



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Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD2047

Sampled: 04/28/05  
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### DATA QUALIFIERS AND DEFINITIONS

- E** Concentration exceeds the calibration range and therefore result is semi-quantitative.
- M2** The MS and/or MSD were below the acceptance limits due to sample matrix interference. See Blank Spike (LCS).
- M-HA** Due to high levels of analyte in the sample, the MS/MSD calculation does not provide useful spike recovery information. See Blank Spike (LCS).
- M-NR1** There was no MS/MSD analyzed with this batch due to insufficient sample volume. See Blank Spike/Blank Spike Duplicate.
- R-7** LFB/LFBD RPD exceeded the method control limit. Recovery met acceptance criteria.
- ZX** Due to sample matrix effects, the surrogate recovery was outside the acceptance limits.
- ND** Analyte NOT DETECTED at or above the reporting limit or MDL, if MDL is specified.
- RPD** Relative Percent Difference

### ADDITIONAL COMMENTS

**For GRO (C4-C12):**

GRO (C4-C12) is quantitated against a gasoline standard. Quantitation begins immediately following the methanol peak.

**For Extractable Fuel Hydrocarbons (EFH, DRO, ORO) :**

Unless otherwise noted, Extractable Fuel Hydrocarbons (EFH, DRO, ORO) are quantitated against a Diesel Fuel Standard.

Del Mar Analytical, Irvine  
Michele Harper  
Project Manager



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Project ID: Alfa Outfall 012 - During Test

Report Number: IOD2047

Sampled: 04/28/05  
Received: 04/28/05

Certification Summary

Del Mar Analytical, Irvine

| Method        | Matrix | Nelac | California |
|---------------|--------|-------|------------|
| EPA 160.2     | Water  | X     | X          |
| EPA 160.5     | Water  | X     | X          |
| EPA 180.1     | Water  | X     | X          |
| EPA 314.0     | Water  | N/A   | X          |
| EPA 350.2     | Water  |       | X          |
| EPA 405.1     | Water  | X     | X          |
| EPA 413.1     | Water  | X     | X          |
| EPA 418.1     | Water  | X     | X          |
| EPA 624       | Water  | X     | X          |
| EPA 625       | Water  | X     | X          |
| EPA 8015 Mod. | Water  | X     | X          |
| EPA 8015B     | Water  | X     | X          |
| SM2540C       | Water  | X     | X          |

Nevada and NELAP provide analyte specific accreditations. Analyte specific information for Del Mar Analytical may be obtained by contacting the laboratory or visiting our website at [www.dmalabs.com](http://www.dmalabs.com).

Subcontracted Laboratories

**Del Mar Analytical - Phoenix** NELAC Cert #01109CA, California Cert #2446, Arizona Cert #AZ0426, Nevada Cert #AZ-907

9830 S. 51st Street, Suite B-120 - Phoenix, AZ 85044

Method Performed: EPA 8260B

Samples: IOD2047-01

Del Mar Analytical, Irvine  
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## SUBCONTRACT ORDER - PROJECT # IOD2047

| SENDING LABORATORY:   | RECEIVING LABORATORY:   |
|---|---|
| Del Mar Analytical, Irvine<br>17461 Derian Avenue, Suite 100<br>Irvine, CA 92614<br>Phone: (949) 261-1022<br>Fax: (949) 261-1228<br>Project Manager: Michele Harper | Del Mar Analytical - Phoenix<br>9830 S. 51st Street, Suite B-120<br>Phoenix, AZ 85044<br>Phone: (480) 785-0043<br>Fax: (480) 785-0851 |

| Analysis  | Expiration     | Due            | Comments   |
|---|----------------|----------------|--|
| <b>Sample ID: IOD2047-01 Water      Sampled: 04/28/05 14:35</b> |                |                |  |
| Dioxane-8260B-out   | 05/12/05 14:35 | 05/09/05 12:00 | Boeing-permit, sub DMAP, J flags, ID=DMA+Outfall 012 |
| Level 4 Data Package - Out                                      | 05/26/05 14:35 | 05/09/05 12:00 | Boeing   |
| <b>Containers Supplied:</b>                                     |                |                |  |
| 40 ml VOA w/HCL (IOD2047-01H)                                   |                |                |  |
| 40 ml VOA w/HCL (IOD2047-01I)                                   |                |                |  |
| 40 ml VOA w/HCL (IOD2047-01J)                                   |                |                |  |

PO00903-01

| SAMPLE INTEGRITY:      |   |                             |   |
|------------------------|---|-----------------------------|---|
| All containers intact: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No | Sample labels/COC agree:    | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No |
| Custody Seals Present: | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No | Samples Preserved Properly: | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No |
|                        |   | Samples Received On Ice::   | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No |
|                        |   | Samples Received at (temp): | 5.8°C   |

~~Released By: \_\_\_\_\_ Date: 4/29/05 Time: 17:00 Received By: \_\_\_\_\_ Date: 4/30/05 Time: \_\_\_\_\_~~

Released By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_ Received By: *[Signature]* Date: 4/30/05 Time: 9:25

74 10022047

**CHAIN OF CUSTODY FORM**

Del Mar Analytical Version 02/17/05

**Client Name/Address:**  
 MWH-Pasadena  
 300 North Lake Avenue, Suite 1200  
 Pasadena, CA 91101  
**Project Manager:** Bronwyn Kelly  
**Sampler:** Pollock

**ANALYSIS REQUIRED**

| Sample Description    | Sample Matrix | Container Type | # of Cont. | Preservative | Bottle #                     | Oil & Grease (EPA 413.1) | 8015-gas | 8015-diesel/jet fuel | 1,4-Dioxane-8260B | TRPH = Total Rec. Petroleum Hydrocarbons (EPA 418.1) | 624 (EDB, 1,2,3-TCF, MTBE, DPE, TBA) | BOD5(20 degrees C) | 625 Naphthalene +NDMA analysis | Ammonia-N, Titr. (350.2) w/ dist | Perchlorate | Turbidity, TDS, TSS | Settleable Solids | Field readings:         | Comments |
|-----------------------|---------------|----------------|------------|--------------|------------------------------|--------------------------|----------|----------------------|-------------------|--|--------------------------------------|--------------------|--------------------------------|----------------------------------|-------------|---------------------|-------------------|-------------------------|----------|
| Outfall 012           | W             | 1L Amber       | 1          | HCl          | 1A                           | X                        |          |                      |                   |  |                                      |                    |                                |                                  |             |                     |                   | Temp = 65.8<br>pH = 7.2 |          |
| Outfall 012 duplicate | W             | 1L Amber       | 1          | HCl          | 1B                           | X                        |          |                      |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | VOAs           | 1          | HCl          | 2A                           |                          | X        |                      |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012 duplicate | W             | VOAs           | 2          | HCl          | 2B, 2C                       |                          | X        |                      |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | 1L Amber       | 1          | None         | 3A                           |                          | X        |                      |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012 duplicate | W             | 1L Amber       | 1          | None         | 3B                           |                          | X        |                      |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | VOAs           | 1          | HCl          | 4A                           |                          |          | X                    |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012 duplicate | W             | VOAs           | 2          | HCl          | 4B, 4C                       |                          |          | X                    |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | 1L Amber       | 1          | HCl          | 5A                           |                          |          |                      | X                 |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012 duplicate | W             | 1L Amber       | 1          | HCl          | 5B                           |                          |          |                      |                   | X  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | VOAs           | 1          | HCl          | 6A                           |                          |          |                      |                   |  | X                                    |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012 duplicate | W             | VOAs           | 2          | HCl          | 6B, 6C                       |                          |          |                      |                   |  | X                                    |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | 1L Poly        | 1          | None         | 7A                           |                          |          |                      |                   |  | X                                    |                    |                                |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | 1L Amber       | 1          | None         | 8A                           |                          |          |                      |                   |  |                                      | X                  |                                |                                  |             |                     |                   |                         |          |
| Outfall 012 duplicate | W             | 1L Amber       | 1          | None         | 8B                           |                          |          |                      |                   |  |                                      | X                  |                                |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | 500ml Poly     | 1          | H2S04        | 9A                           |                          |          |                      |                   |  |                                      |                    | X                              |                                  |             |                     |                   |                         |          |
| Outfall 012           | W             | 1L Poly        | 1          | None         | 10A                          |                          |          |                      |                   |  |                                      |                    |                                | X                                |             |                     |                   |                         |          |
| Outfall 012           | W             | 1L Poly        | 1          | None         | 11A                          |                          |          |                      |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |
| Trip Blank            | W             | VOAs           | 1          | HCl          | 12A, 12B, 12C, 12D, 12E, 12F |                          |          | X                    |                   |  |                                      |                    |                                |                                  |             |                     |                   |                         |          |

**Relinquished By:** [Signature] Date/Time: 4-28-05 1530  
**Received By:** [Signature] Date/Time: 4/28/05 1530

**Relinquished By:** [Signature] Date/Time: 4/28/05 1815  
**Received By:** [Signature] Date/Time: 4/28/05 1815

**Relinquished By:** [Signature] Date/Time: \_\_\_\_\_  
**Received By:** [Signature] Date/Time: \_\_\_\_\_

Turn around Time: (check)  
 24 Hours \_\_\_\_\_ 5 Days \_\_\_\_\_  
 48 Hours \_\_\_\_\_ 10 Days \_\_\_\_\_  
 72 Hours \_\_\_\_\_ Normal \_\_\_\_\_  
 Perchlorate Only 72 Hours \_\_\_\_\_  
 Metals Only 72 Hours \_\_\_\_\_

Sample Integrity: (Check)  On Ice:  4C



QA/QC DATA PACKAGE: LEVEL IV



QA/QC DATA PACKAGE LEVEL IV

TABLE SUMMARY

Page Number

|                                |       |
|--------------------------------|-------|
| CASE NARRATIVE.....            | 1     |
| CHAIN OF CUSTODY.....          | 2-3   |
| ANALYTICAL REPORTS.....        | 4-5   |
| QUALITY CONTROL SUMMARIES..... | 6-9   |
| EPA 8260B RAW DATA.....        | 10-98 |



LABORATORY REPORT

Prepared For: Del Mar Analytical - Irvine  
17461 Derian Ave. Suite 100  
Irvine, CA 92614  
Attention: Michele Harper

Project: IOD2047

Sampled: 04/28/05  
Received: 04/30/05  
Issued: 05/11/05 16:11

NELAP #01109CA California ELAP#2446

*The results listed within this Laboratory Report pertain only to the samples tested in the laboratory. The analyses contained in this report were performed in accordance with the applicable certifications as noted. All soil samples are reported on a wet weight basis unless otherwise noted in the report. This Laboratory Report is confidential and is intended for the sole use of Del Mar Analytical and its client. This report shall not be reproduced, except in full, without written permission from Del Mar Analytical. The Chain of Custody, 1 page, is included and is an integral part of this report.  
This entire report was reviewed and approved for release.*

CASE NARRATIVE

LABORATORY ID  
POD0903-01

CLIENT ID  
IOD2047-01

MATRIX  
Water

- SAMPLE RECEIPT: Samples were received intact, at 6°C, on ice and with chain of custody documentation.
- HOLDING TIMES: All samples were analyzed within prescribed holding times and/or in accordance with the Del Mar Analytical Sample Acceptance Policy unless otherwise noted in the report.
- PRESERVATION: Samples requiring preservation were verified prior to sample analysis.
- QA/QC CRITERIA: All analyses met method criteria, except as noted in the report with data qualifiers.
- COMMENTS: Results that fall between the MDL and RL are 'J' flagged.
- SUBCONTRACTED: No analyses were subcontracted to an outside laboratory.

Reviewed By:

Del Mar Analytical - Phoenix  
Karen Maxwell  
Project Manager





QA/QC DATA PACKAGE: LEVEL IV

CHAIN OF CUSTODY FORMS



17461 Derian Ave. Suite 100, Irvine, CA 92614 Ph (949) 261-1022 Fax (949) 261-1228  
 1014 E. Cooley Dr., Suite A, Colton, CA 92324 Ph (909) 370-4667 Fax (909) 370-1046  
 9484 Chesapeake Drive, Suite 805, San Diego, CA 92123 Ph (619) 505-9586 Fax (619) 505-9689  
 9830 South 51st Street, Suite B-120, Phoenix, AZ 85044 Ph (480) 785-0043 Fax (480) 785-0851  
 2520 E. Sunset Rd., Suite #3, Las Vegas, NV 89120 Ph (702) 796-3620 Fax (702) 796-3621

## SUBCONTRACT ORDER - PROJECT # IOD2047

| SENDING LABORATORY:   | RECEIVING LABORATORY:   |
|---|---|
| Del Mar Analytical, Irvine<br>17461 Derian Avenue, Suite 100<br>Irvine, CA 92614<br>Phone: (949) 261-1022<br>Fax: (949) 261-1228<br>Project Manager: Michele Harper | Del Mar Analytical - Phoenix<br>9830 S. 51st Street, Suite B-120<br>Phoenix, AZ 85044<br>Phone: (480) 785-0043<br>Fax: (480) 785-0851 |

| Analysis                    | Expiration     | Due                     | Comments   |
|-----------------------------|----------------|-------------------------|--|
| Sample ID: IOD2047-01 Water |                | Sampled: 04/28/05 14:35 |  |
| Dioxane-8260B-out           | 05/12/05 14:35 | 05/09/05 12:00          | Boeing-permit, sub DMAP, J flags, ID=DMA+Outfall 012 |
| Level 4 Data Package - Out  | 05/26/05 14:35 | 05/09/05 12:00          | Boeing   |

**Containers Supplied:**

- 40 ml VOA w/HCL (IOD2047-01H)
- 40 ml VOA w/HCL (IOD2047-01I)
- 40 ml VOA w/HCL (IOD2047-01J)

P000903-01

| SAMPLE INTEGRITY:      |   |  |                             |   |                             |
|------------------------|---|--|-----------------------------|---|-----------------------------|
| All containers intact: | <input checked="" type="checkbox"/> Yes | <input type="checkbox"/> No            | Sample labels/COC agree:    | <input checked="" type="checkbox"/> Yes | <input type="checkbox"/> No |
| Custody Seals Present: | <input type="checkbox"/> Yes            | <input checked="" type="checkbox"/> No | Samples Preserved Properly: | <input checked="" type="checkbox"/> Yes | <input type="checkbox"/> No |
|                        |   |  | Samples Received On Ice:    | <input checked="" type="checkbox"/> Yes | <input type="checkbox"/> No |
|                        |   |  | Samples Received at (temp): | 5.8C                                    |                             |

|                        |                    |                  |                        |                    |      |
|------------------------|--------------------|------------------|------------------------|--------------------|------|
| <del>Released By</del> | <del>4/29/05</del> | <del>17:00</del> | <del>Received By</del> | <del>4/30/05</del> |      |
| Released By            |                    |                  | Received By            | 4/30/05            | 9:25 |



QC DATA PACKAGE: LEVEL IV

ANALYTICAL REPORTS



# Del Mar Analytical

17461 Dentan Ave., Suite 100, Irvine, CA 92614 (949) 261-1022 FAX (949) 260-3297  
 1014 E. Cooley Dr., Suite A, Colton, CA 92324 (909) 370-4667 FAX (949) 370-1046  
 9484 Chesapeake Dr., Suite 805, San Diego, CA 92123 (858) 505-8596 FAX (858) 505-9689  
 9830 South 51st St., Suite B-120, Phoenix, AZ 85044 (480) 785-0043 FAX (480) 785-0851  
 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

Del Mar Analytical - Irvine  
 17461 Derian Ave. Suite 100  
 Irvine, CA 92614  
 Attention: Michele Harper

Project ID: IOD2047

Report Number: POD0903

Sampled: 04/28/05  
 Received: 04/30/05

## 1,4-DIOXANE BY GC/MS (EPA 5030B/8260B)

| Analyte   | Method    | Batch   | MDL Limit | Reporting Limit | Sample Result | Dilution Factor | Date Extracted | Date Analyzed | Data Qualifiers |
|---|-----------|---------|-----------|-----------------|---------------|-----------------|----------------|---------------|-----------------|
| <b>Sample ID: POD0903-01 (IOD2047-01 - Water)</b> |           |         |           |                 |               |                 |                |               |                 |
| Reporting Units: ug/l                             |           |         |           |                 |               |                 |                |               |                 |
| I,4-Dioxane                                       | EPA 8260B | P5E0214 | 0.49      | 1.0             | ND            | 1               | 05/02/05       | 05/02/05      |                 |
| Surrogate: Dibromofluoromethane (80-125%)         |           |         |           |                 | 97 %          |                 |                |               |                 |

Del Mar Analytical - Phoenix  
 Karen Maxwell  
 Project Manager

*The results pertain only to the samples tested in the laboratory. This report shall not be reproduced, except in full, without written permission from Del Mar Analytical.*



QA/QC DATA PACKAGE: LEVEL IV

QUALITY CONTROL SUMMARIES

Del Mar Analytical - Irvine  
 17461 Derian Ave. Suite 100  
 Irvine, CA 92614  
 Attention: Michele Harper

Project ID: IOD2047

Report Number: POD0903

Sampled: 04/28/05  
 Received: 04/30/05

## METHOD BLANK/QC DATA

### 1,4-DIOXANE BY GC/MS (EPA 5030B/8260B)

| Analyte   | Result | Reporting Limit | MDL  | Units | Spike Level | Source Result                | %REC | %REC Limits | RPD | RPD Limit | Data Qualifiers |
|---|--------|-----------------|------|-------|-------------|------------------------------|------|-------------|-----|-----------|-----------------|
| <b>Batch: P5E0214 Extracted: 05/02/05</b>                   |        |                 |      |       |             |                              |      |             |     |           |                 |
| <b>Blank Analyzed: 05/02/2005 (P5E0214-BLK1)</b>            |        |                 |      |       |             |                              |      |             |     |           |                 |
| 1,4-Dioxane   | ND     | 1.0             | 0.49 | ug/l  |             |                              |      |             |     |           |                 |
| Surrogate: Dibromofluoromethane                             | 1.03   |                 |      | ug/l  | 1.00        |                              | 103  | 80-125      |     |           |                 |
| <b>LCS Analyzed: 05/02/2005 (P5E0214-BS1)</b>               |        |                 |      |       |             |                              |      |             |     |           |                 |
| 1,4-Dioxane   | 12.3   | 1.0             | 0.49 | ug/l  | 10.0        |                              | 123  | 70-130      |     |           |                 |
| Surrogate: Dibromofluoromethane                             | 1.00   |                 |      | ug/l  | 1.00        |                              | 100  | 80-125      |     |           |                 |
| <b>LCS Dup Analyzed: 05/02/2005 (P5E0214-BSD1)</b>          |        |                 |      |       |             |                              |      |             |     |           |                 |
| 1,4-Dioxane   | 9.82   | 1.0             | 0.49 | ug/l  | 10.0        |                              | 98   | 70-130      | 22  | 20        | R-7             |
| Surrogate: Dibromofluoromethane                             | 0.970  |                 |      | ug/l  | 1.00        |                              | 97   | 80-125      |     |           |                 |
| <b>Matrix Spike Analyzed: 05/02/2005 (P5E0214-MS1)</b>      |        |                 |      |       |             |                              |      |             |     |           |                 |
|   |        |                 |      |       |             | <b>Source: POD0904-02RE1</b> |      |             |     |           |                 |
| 1,4-Dioxane   | 176    | 1.0             | 0.49 | ug/l  | 10.0        | 140                          | 360  | 70-150      |     |           | E, M-HA         |
| Surrogate: Dibromofluoromethane                             | 0.980  |                 |      | ug/l  | 1.00        |                              | 98   | 80-125      |     |           |                 |
| <b>Matrix Spike Dup Analyzed: 05/02/2005 (P5E0214-MSD1)</b> |        |                 |      |       |             |                              |      |             |     |           |                 |
|   |        |                 |      |       |             | <b>Source: POD0904-02RE1</b> |      |             |     |           |                 |
| 1,4-Dioxane   | 151    | 1.0             | 0.49 | ug/l  | 10.0        | 140                          | 110  | 70-150      | 15  | 25        | E, M-HA         |
| Surrogate: Dibromofluoromethane                             | 0.970  |                 |      | ug/l  | 1.00        |                              | 97   | 80-125      |     |           |                 |

Del Mar Analytical - Phoenix  
 Karen Maxwell  
 Project Manager



Del Mar Analytical - Irvine  
17461 Derian Ave. Suite 100  
Irvine, CA 92614  
Attention: Michele Harper

Project ID: IOD2047

Report Number: POD0903

Sampled: 04/28/05  
Received: 04/30/05

## DATA QUALIFIERS AND DEFINITIONS

- E** Concentration exceeds the calibration range and therefore result is semi-quantitative.
- M-HA** Due to high levels of analyte in the sample, the MS/MSD calculation does not provide useful spike recovery information. See Blank Spike (LCS).
- R-7** LFB/LFBD RPD exceeded the method control limit. Recovery met acceptance criteria.
- ND** Analyte NOT DETECTED at or above the reporting limit or MDL, if MDL is specified.
- RPD** Relative Percent Difference

**Del Mar Analytical - Phoenix**  
Karen Maxwell  
Project Manager



# Del Mar Analytical

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 2520 E. Sunset Rd. #3, Las Vegas, NV 89120 (702) 798-3620 FAX (702) 798-3621

Del Mar Analytical - Irvine  
 17461 Derian Ave. Suite 100  
 Irvine, CA 92614  
 Attention: Michele Harper

Project ID: IOD2047

Report Number: POD0903

Sampled: 04/28/05  
 Received: 04/30/05

## Certification Summary

### Del Mar Analytical - Phoenix

| Method    | Matrix | Nelac | California |
|-----------|--------|-------|------------|
| EPA 8260B | Water  | X     | X          |

*Nevada and NELAP provide analyte specific accreditations. Analyte specific information for Del Mar Analytical may be obtained by contacting the laboratory or visiting our website at [www.dmalabs.com](http://www.dmalabs.com).*

**Del Mar Analytical - Phoenix**  
 Karen Maxwell  
 Project Manager

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QA/QC DATA PACKAGE: LEVEL IV

**EPA METHOD 8260B LABORATORY RAW DATA**

GCMS TUNING  
INITIAL/DAILY CALIBRATION  
RUNLOG  
CONTINUING CALBRATION  
QUANTITATION REPORTS  
CHROMATOGRAMS  
EXTRACTION LOG  
STANDARD LOG

CMS #: 1

Date Analyzed: 3/19/05

| ANALYST REVIEW | METHOD CRITERIA | PEER REVIEW |
|----------------|-----------------|-------------|
|----------------|-----------------|-------------|

- |    |  |    |
|----|--|----|
| ✓  | 1. Sequence File is printed and in the file folder?<br>Standard IDs and analyst's initials are present?  | ✓  |
| ✓  | 2. Initial Calibration met criteria?<br>a. Print calibration as Average Response Factor<br>(624: RSD ≤ 35%)<br>(8260B: ≤ 30% for CCCs, ≤ 15% for all other compounds, SPCCs met Criteria)<br>(524.2: RSD ≤ 20%)  | ✓  |
| ✓  | b. If non CCC RSD > 15%, print out the curve as Linear Regression<br>$r \geq 0.995$ or $r^2 \geq 0.99$ (do not force through zero for 8260B)   | ✓  |
| ✓  | c. If non CCC RSD > 15%, print out the curve as Quadratic<br>$r \geq 0.995$ or $r^2 \geq 0.99$ (do not force through zero for 8260B)   | ✓  |
| ✓  | d. Choose option (b or c) with the least negative intercept  | ✓  |
| ✓  | e. Requant the low (RL) standard against the curve<br>must be ± 30%, file with the calibration for reference   | ✓  |
| ✓  | f. If samples contain negative values then:<br>compare the area counts with the low standard on file<br>if <, then report as N.D. with no flag<br>if >, then report from RSD curve and flag that curve is out<br>or report at an elevated RL as compared to a curve standard | ✓  |
| ✓  | 3. Initial Midpoint / LCS / BFB Tune<br>(624: use Table 5) (524.2: ±30%) (8260B: see control chart)<br>SPCCs met criteria? CCCs met criteria (±20%)?   | ✓  |
| ✓  | 4. Checked integration of all peaks in Midpoint?   | ✓  |
| ✓  | 5. Method Blank < Report Limit, if not is data flagged?<br>(624: every 20 samples) (524.2: every 12 hours) (8260B: every 12 hours)   | ✓  |
| NA | 6. MS/MSD (every 20 samples)<br>(624: use Table 5) (524.2: N/A) (8260B: see Control Chart)   | NA |
| ✓  | 7. All samples met holding time? (Soil 72hr ext, 7/14days water)   | ✓  |
| ✓  | 8. All water samples checked to be pH < 2? (Note this on the sequence file)  | ✓  |
| ✓  | 9. LCS every 20 samples<br>(624: See Table 5) (524.2: ±30%) (8260B: See Control Chart)   | ✓  |
| ✓  | 10. Cont. Midpoint / LCS / BFB Tune done every 12 hours<br>(624: use Table 5) (524.2: ±30%) (8260B: see control chart)<br>SPCCs met criteria? CCCs met criteria (±20%)?  | ✓  |
| ✓  | 11. Surrogates within acceptance limits<br>(624 / 524.2 / 8260B: See Control Chart)  | ✓  |
| ✓  | 12. Internal Standards within acceptance limits<br>(624 / 524.2 / 8260B: response must be -50 to +100%).   | ✓  |
|    | 13. Manual re-integration(s) performed?<br>yes: _____ no: ✓  |    |
|    | 14. Corrective Action Report required?<br>yes: _____ (Attached) no: ✓  |    |
|    | 15. Reports impacted by the Corrective Action Report   |    |

Analyst: J. Galassi 3/21/05

Reviewer / Date: [Signature] 3/21/05

# DMAP GC/MS 1 DAILY LOG SUMMARY

DATE: 3/19/05

QC BATCH # (s):

CAL CURVE  
R501902J63/2/05

ANALYST: JY/MS

SEQUENCE FILE: C:\GCMS1\DATA\

CALIBRATION METHOD(S): DX021605.M/W072903.M

| POS # | FILENAME | SAMPLE ID.CLIENT      | SAMPLE VOL. | pH | EPA METHOD | MATRIX | COMMENTS                                  |
|-------|----------|-----------------------|-------------|----|------------|--------|---|
| ✓     | P0319001 | TUNE                  | 1ul         | NA | 82100      | H2O    |   |
| 1     | 2        | CCV                   | 1x10ML      |    |            |        |   |
| 2     | 3        | LCS DUP               |             |    |            |        | -DNV IS LOW                               |
| 3     | 4        | LCS DUP               |             |    |            |        | -DNV IS LOW -> Rep<br>+ xp                |
| ✓     | 5        | TUNE                  |             |    |            |        |   |
| 1     | 6        | CCV                   |             |    |            |        | -DNV, IS's still low<br>will re-calibrate |
| 2     | 7        | CCV                   |             |    |            |        |   |
| 3     | 8        | Blank                 |             |    |            |        |   |
| 4     | 9        | 1.0 ppb Cal           |             |    |            |        | DNV's Grubb's P/H<br>outlier              |
| 5     | 10       | 2.0                   |             |    |            |        |   |
| 6     | 11       | 5.0                   |             |    |            |        |   |
| 7     | 12       | 10.0                  |             |    |            |        |   |
| 8     | 13       | 20.0                  |             |    |            |        |   |
| 9     | 14       | 50.0                  |             |    |            |        |   |
| 10    | 15       | 100.0                 |             |    |            |        |   |
| 11    | 16       | clean at blank / time |             |    |            |        |   |
| 12    | 17       | MSD CV Blank          |             |    |            |        |   |
| 13    | 18       | 1.0 ppb Cal           |             |    |            |        |   |
| 14    | 19       | 55/CCV                |             |    |            |        |   |

**STANDARD ID NUMBERS**

CCV / H<sub>2</sub>O LCS / H<sub>2</sub>O SPIKE: 5030018

Internal Std: 5030259 <sup>353</sup> 3/21/05

CALIBRATION STD: 5030348/5030349

IS / Surrogate / BFB: 5030321

REVIEWER / DATE: SJ/2/05

tune / 5030090

# Injection Log

Directory: D:\HPCHEM\1\DATA\031905

| Line | Vial | FileName   | Multiplier | SampleName                          | Misc Info | Injected          |
|------|------|------------|------------|-------------------------------------|-----------|-------------------|
| 1    | 1    | P0319001.D | 1.         | <del>TUNE/BLANK</del>               | 1X 10ML   | 19 Mar 2005 06:19 |
| 2    | 2    | P0319002.D | 1.         | CCV                                 | 1X 10ML   | 19 Mar 2005 06:32 |
| 3    | 3    | P0319003.D | 1.         | LCS DUP                             | 1X 10ML   | 19 Mar 2005 07:08 |
| 4    | 4    | P0319004.D | 1.         | <del>LCS DUP DNU</del>              | 1X 10ML   | 19 Mar 2005 07:44 |
| 5    | 5    | P0319005.D | 1.         | TUNE                                | 1X 10ML   | 19 Mar 2005 08:39 |
| 6    | 6    | P0319006.D | 1.         | CCV                                 | 1X 10ML   | 19 Mar 2005 09:07 |
| 7    | 7    | P0319007.D | 1.         | <del>CCV DNU</del>                  | 1X 10ML   | 19 Mar 2005 09:39 |
| 8    | 8    | P0319008.D | 1.         | BLANK                               | 1X 10ML   | 19 Mar 2005 10:12 |
| 9    | 9    | P0319009.D | 1.         | <del>1.0 PPB CAL DNU</del>          | 1X 10ML   | 19 Mar 2005 10:54 |
| 10   | 10   | P0319010.D | 1.         | 2.0 PPB CAL                         | 1X 10ML   | 19 Mar 2005 11:26 |
| 11   | 11   | P0319011.D | 1.         | 5.0 PPB CAL                         | 1X 10ML   | 19 Mar 2005 11:59 |
| 12   | 12   | P0319012.D | 1.         | 10.0 PPB CAL                        | 1X 10ML   | 19 Mar 2005 12:32 |
| 13   | 13   | P0319013.D | 1.         | 20.0 PPB CAL                        | 1X 10ML   | 19 Mar 2005 13:05 |
| 14   | 14   | P0319014.D | 1.         | 50.0 PPB CAL                        | 1X 10ML   | 19 Mar 2005 13:38 |
| 15   | 15   | P0319015.D | 1.         | 100.0 PPB CAL                       | 1X 10ML   | 19 Mar 2005 14:11 |
| 16   | 16   | P0319016.D | 1.         | <del>CLEAN OUT BLANK/TUNE DNU</del> | 1X 10ML   | 19 Mar 2005 14:44 |
| 17   | 17   | P0319017.D | 1.         | BLANK                               | 1X 10ML   | 19 Mar 2005 15:21 |
| 18   | 18   | P0319018.D | 1.         | 1.0 PPB CAL                         | 1X 10ML   | 19 Mar 2005 15:54 |
| 19   | 19   | P0319019.D | 1.         | SS/CCV                              | 1X 10ML   | 19 Mar 2005 16:27 |

3/21/05  
JW

BFB

Data File : D:\HPCHEM\1\DATA\031905\P0319005.D

Acq On : 19 Mar 2005 8:39 am

Sample : TUNE

Misc : 1X 10ML

MS Integration Params: DIOXANE.P

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)

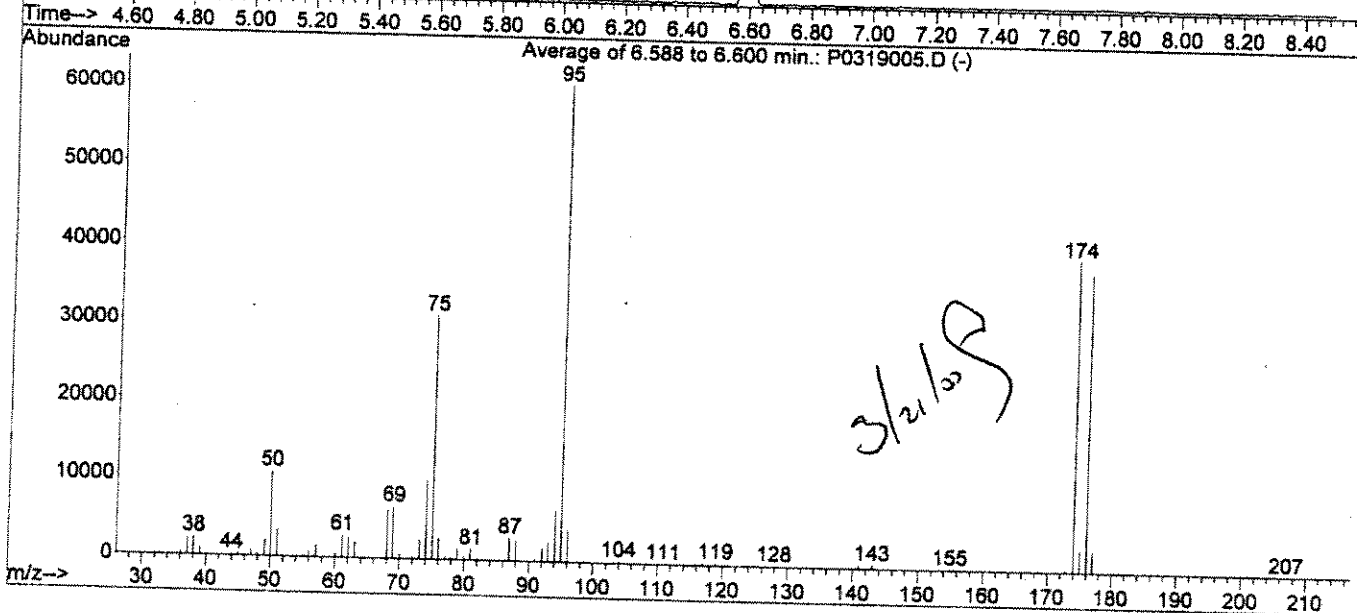
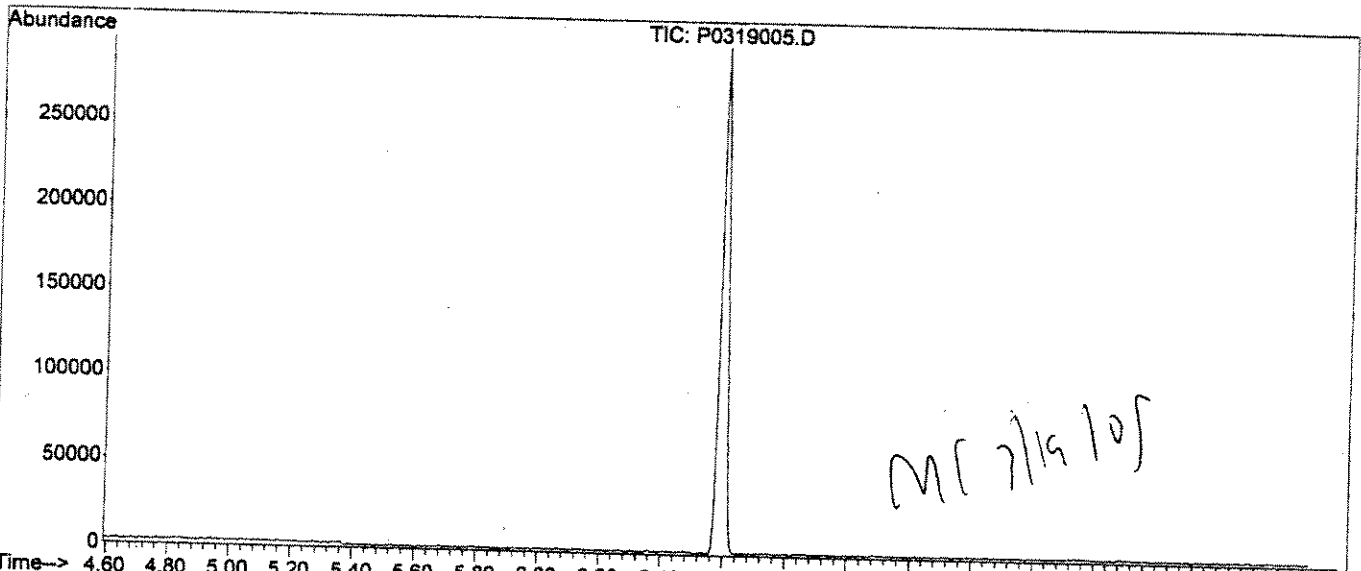
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)

Vial: 5

Operator: JG/MS/CLS

Inst : GCMS1

Multiplr: 1.00



AutoFind: Scans 411, 412, 413; Background Corrected with Scan 395

| Target Mass | Rel. to Mass | Lower Limit% | Upper Limit% | Rel. Abn% | Raw Abn | Result Pass/Fail |
|-------------|--------------|--------------|--------------|-----------|---------|------------------|
| 50          | 95           | 15           | 40           | 17.5      | 10615   | PASS             |
| 75          | 95           | 30           | 60           | 51.3      | 31037   | PASS             |
| 95          | 95           | 100          | 100          | 100.0     | 60549   | PASS             |
| 96          | 95           | 5            | 9            | 6.6       | 3996    | PASS             |
| 173         | 174          | 0.00         | 2            | 0.6       | 226     | PASS             |
| 174         | 95           | 50           | 100          | 65.5      | 39648   | PASS             |
| 175         | 174          | 5            | 9            | 6.9       | 2752    | PASS             |
| 176         | 174          | 95           | 101          | 96.0      | 38059   | PASS             |
| 177         | 176          | 5            | 9            | 6.9       | 2638    | PASS             |

| Grubbs Test for curve |  | 1.0ppb  | 2.0ppb  | 5.0ppb   | 10ppb    | 20ppb    | 50ppb    | 100ppb   | MEAN     | STDEV     |
|-----------------------|--|---------|---------|----------|----------|----------|----------|----------|----------|-----------|
| Response factors      |  | 3.099   | 2.478   | 2.101    | 1.905    | 1.995    | 1.822    | 1.905    | 2.186429 | 0.4566975 |
| Grubbs value          |  | 1.99698 | 0.63805 | 0.186944 | 0.615851 | 0.418904 | 0.797481 | 0.615851 |          |           |
| 5pts Grubbs values <  |  |         |         |          |          |          |          |          |          |           |
| 6pts Grubbs values <  |  |         |         |          |          |          |          |          |          |           |
| 7pts Grubbs values <  |  |         |         |          |          |          |          |          |          |           |
| 8pts Grubbs values <  |  |         |         |          |          |          |          |          |          |           |
| 9pts Grubbs values <  |  |         |         |          |          |          |          |          |          |           |
| 10pts Grubbs values < |  |         |         |          |          |          |          |          |          |           |

outlier

MS 3/19/05

*Handwritten signature*

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\PO319009.D  
 Acq On : 19 Mar 2005 10:54 am  
 Sample : 1.0 PPB CAL  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 13:42 2005

Vial: 9  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX021605.RES

Quant Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Wed Feb 16 15:53:54 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*MI 3/19/05*

*See Grubb's Test JG 3/21/05*

| Internal Standards            | R.T.  | QIon  | Response | Conc     | Units | Dev(Min)  |
|-------------------------------|-------|-------|----------|----------|-------|-----------|
| 1) Pentafluorobenzene (IS)    | 10.56 | 99    | 42761    | 1.00     | ug/L  | 0.00      |
| 3) 1,4-DIOXANE-d8             | 12.35 | 64    | 4961     | 25.00    | ug/L  | 0.00      |
| 5) 1,2,3-Trichloropropane-d5  | 0.00  | 79    | 0        | 0.00     | ug/L  | -15.08    |
| <i>NT</i>                     |       |       |          |          |       |           |
| System Monitoring Compounds   |       |       |          |          |       |           |
| 2) Dibromofluoromethane (SU1) | 10.07 | 113   | 3531     | 0.11     | ug/L  | 0.00      |
| Spiked Amount                 | 1.000 | Range | 80 - 120 | Recovery | =     | 11.00%#   |
| Target Compounds              |       |       |          |          |       |           |
| 4) 1,4-DIOXANE                | 12.43 | 88    | 615      | 1.50     | ug/L  | Qvalue 97 |

*DNV*

*Q*

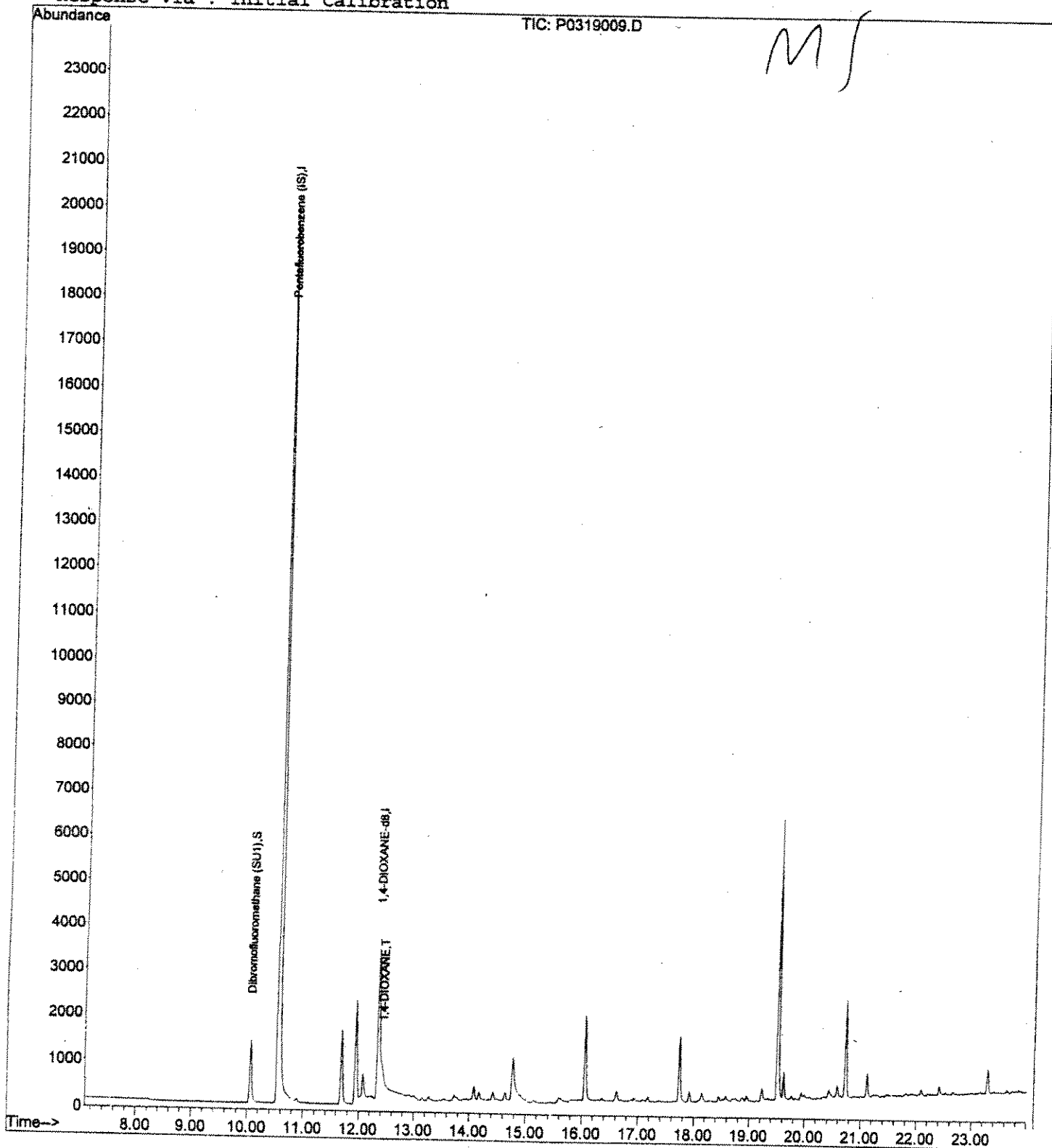
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\0319009.D  
Acq On : 19 Mar 2005 10:54 am  
Sample : 1.0 PPB CAL  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 19 13:42 2005

Vial: 9  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

Quant Results File: DX021605.RES

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
Last Update : Wed Feb 16 15:53:54 2005  
Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\P0319010.D  
 Acq On : 19 Mar 2005 11:26 am  
 Sample : 2.0 PPB CAL  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 13:43 2005

Vial: 10  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX021605.RES

Quant Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Wed Feb 16 15:53:54 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*MS 3/14/05*

| Internal Standards                 | R.T.  | QIon  | Response | Conc     | Units | Dev(Min)     |
|------------------------------------|-------|-------|----------|----------|-------|--------------|
| 1) Pentafluorobenzene (IS)         | 10.57 | 99    | 45768    | 1.00     | ug/L  | 0.00         |
| 3) 1,4-DIOXANE-d8                  | 12.35 | 64    | 5185     | 25.00    | ug/L  | 0.00         |
| 5) 1,2,3-Trichloropropane-d5       | 0.00  | 79    | ONT      | 0.00     | ug/L  | -15.08       |
| <b>System Monitoring Compounds</b> |       |       |          |          |       |              |
| 2) Dibromofluoromethane (SU1)      | 10.07 | 113   | 7585     | 0.21     | ug/L  | 0.00         |
| Spiked Amount                      | 1.000 | Range | 80 - 120 | Recovery | =     | 21.00%#      |
| <b>Target Compounds</b>            |       |       |          |          |       |              |
| 4) 1,4-DIOXANE                     | 12.43 | 88    | 1028     | 2.69     | ug/L  | Qvalue<br>94 |

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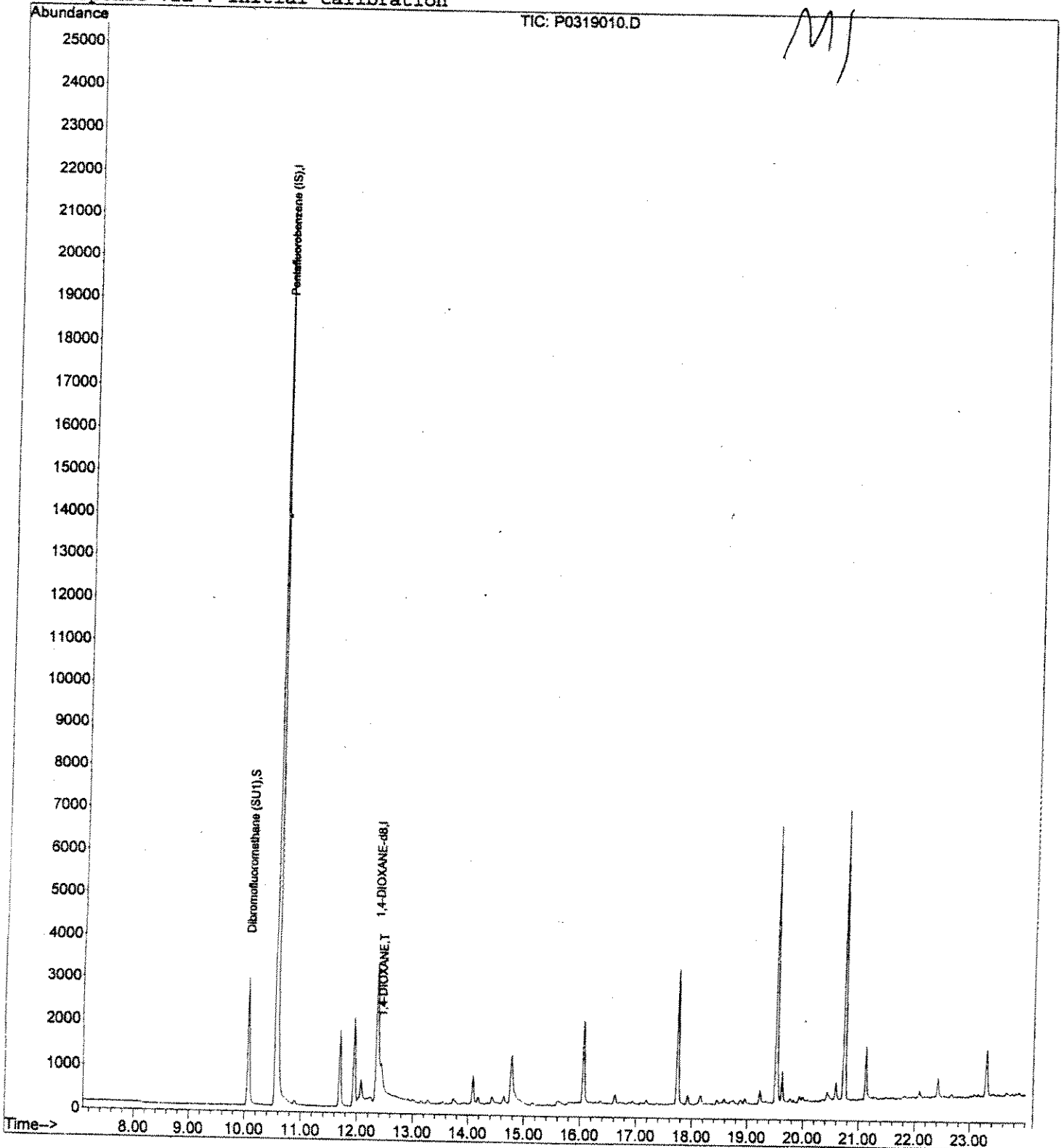
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\0319010.D  
Acq On : 19 Mar 2005 11:26 am  
Sample : 2.0 PPB CAL  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 19 13:43 2005

Vial: 10  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

Quant Results File: DX021605.RES

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
Last Update : Wed Feb 16 15:53:54 2005  
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\P0319011.D  
 Acq On : 19 Mar 2005 11:59 am  
 Sample : 5.0 PPB CAL  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 13:43 2005

Vial: 11  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX021605.RES

Quant Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Wed Feb 16 15:53:54 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*M (3/19/05)*

| Internal Standards            | R.T.  | QIon           | Response | Conc  | Units   | Dev(Min)  |
|-------------------------------|-------|----------------|----------|-------|---------|-----------|
| 1) Pentafluorobenzene (IS)    | 10.56 | 99             | 47558    | 1.00  | ug/L    | 0.00      |
| 3) 1,4-DIOXANE-d8             | 12.35 | 64             | 5263     | 25.00 | ug/L    | 0.00      |
| 5) 1,2,3-Trichloropropane-d5  | 0.00  | 79             | 0NT      | 0.00  | ug/L    | -15.08    |
| System Monitoring Compounds   |       |                |          |       |         |           |
| 2) Dibromofluoromethane (SU1) | 10.06 | 113            | 19072    | 0.52  | ug/L    | 0.00      |
| Spiked Amount                 | 1.000 | Range 80 - 120 | Recovery | =     | 52.00%# |           |
| Target Compounds              |       |                |          |       |         |           |
| 4) 1,4-DIOXANE                | 12.43 | 88             | 2211     | 6.25  | ug/L    | Qvalue 99 |

*3/21/05*

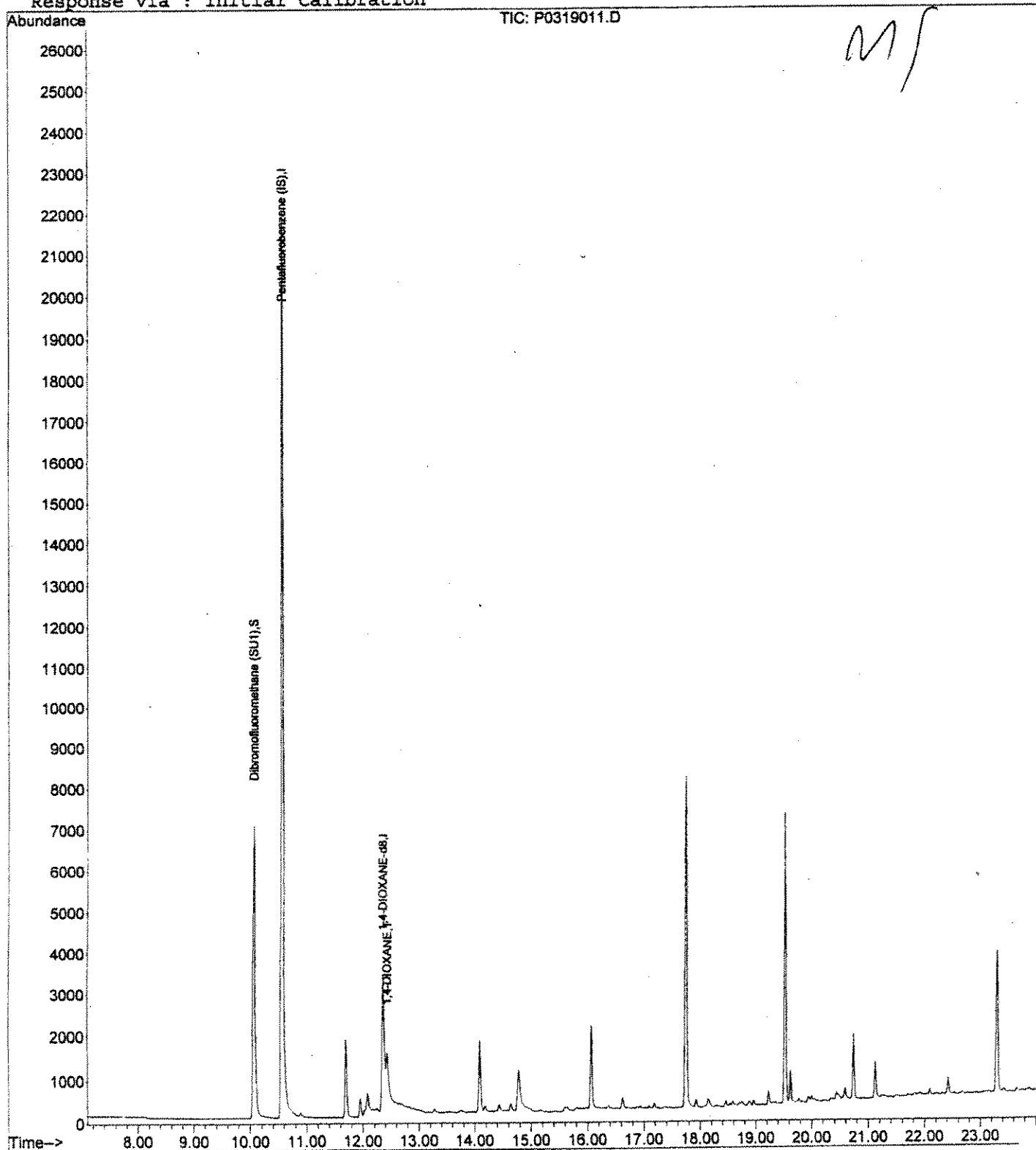
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\P0319011.D  
Acq On : 19 Mar 2005 11:59 am  
Sample : 5.0 PPB CAL  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 19 13:43 2005

Vial: 11  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

Quant Results File: DX021605.RES

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
Last Update : Wed Feb 16 15:53:54 2005  
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\PO319012.D  
 Acq On : 19 Mar 2005 12:32 pm  
 Sample : 10.0 PPB CAL  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 13:37 2005

Vial: 12  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX021605.RES

Quant Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Wed Feb 16 15:53:54 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*M 3/19/05*

| Internal Standards           | R.T.  | QIon | Response   | Conc  | Units | Dev(Min) |
|------------------------------|-------|------|------------|-------|-------|----------|
| 1) Pentafluorobenzene (IS)   | 10.57 | 99   | 47071      | 1.00  | ug/L  | 0.00     |
| 3) 1,4-DIOXANE-d8            | 12.35 | 64   | 5034       | 25.00 | ug/L  | 0.00     |
| 5) 1,2,3-Trichloropropane-d5 | 0.00  | 79   | <i>0.1</i> | 0.00  | ug/L  | -15.08   |

System Monitoring Compounds  
 2) Dibromofluoromethane (SU1) 10.07 113 34373 0.95 ug/L 0.00  
 Spiked Amount 1.000 Range 80 - 120 Recovery = 95.00%

Target Compounds  
 4) 1,4-DIOXANE 12.43 88 3835 11.74 ug/L Qvalue 99

*3/21/05*

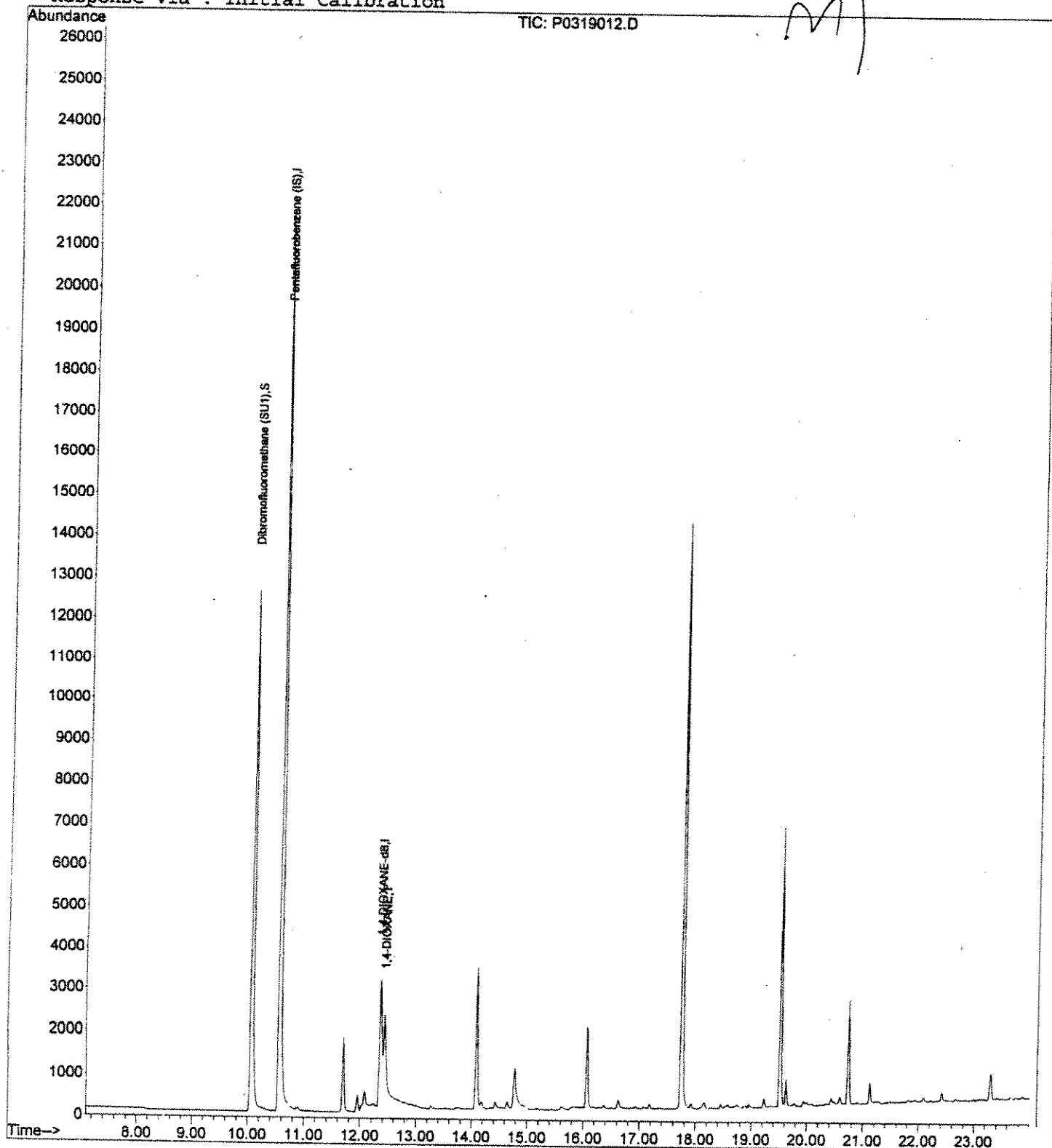
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\PO319012.D  
Acq On : 19 Mar 2005 12:32 pm  
Sample : 10.0 PPB CAL  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 19 13:37 2005

Vial: 12  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

Quant Results File: DX021605.RES

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
Last Update : Wed Feb 16 15:53:54 2005  
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\0319013.D  
 Acq On : 19 Mar 2005 1:05 pm  
 Sample : 20.0 PPB CAL  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 13:37 2005

Vial: 13  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX021605.RES

Quant Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Wed Feb 16 15:53:54 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*M (3/19/05)*

| Internal Standards                 | R.T.  | QIon  | Response | Conc     | Units | Dev(Min)  |
|------------------------------------|-------|-------|----------|----------|-------|-----------|
| 1) Pentafluorobenzene (IS)         | 10.56 | 99    | 47635    | 1.00     | ug/L  | 0.00      |
| 3) 1,4-DIOXANE-d8                  | 12.35 | 64    | 4790     | 25.00    | ug/L  | 0.00      |
| 5) 1,2,3-Trichloropropane-d5       | 0.00  | 79    | 0.00     | 0.00     | ug/L  | -15.08    |
| <b>System Monitoring Compounds</b> |       |       |          |          |       |           |
| 2) Dibromofluoromethane (SU1)      | 10.07 | 113   | 68573    | 1.86     | ug/L  | 0.00      |
| Spiked Amount                      | 1.000 | Range | 80 - 120 | Recovery | =     | 186.00%#  |
| <b>Target Compounds</b>            |       |       |          |          |       |           |
| 4) 1,4-DIOXANE                     | 12.43 | 88    | 7646     | 25.14    | ug/L  | Qvalue 97 |

*3/21/05*

Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\0319013.D

Vial: 13

Acq On : 19 Mar 2005 1:05 pm

Operator: JG/MS/CLS

Sample : 20.0 PPB CAL

Inst : GCMS1

Misc : 1X 10ML

Multiplr: 1.00

MS Integration Params: DIOXANE.P

Quant Time: Mar 19 13:37 2005

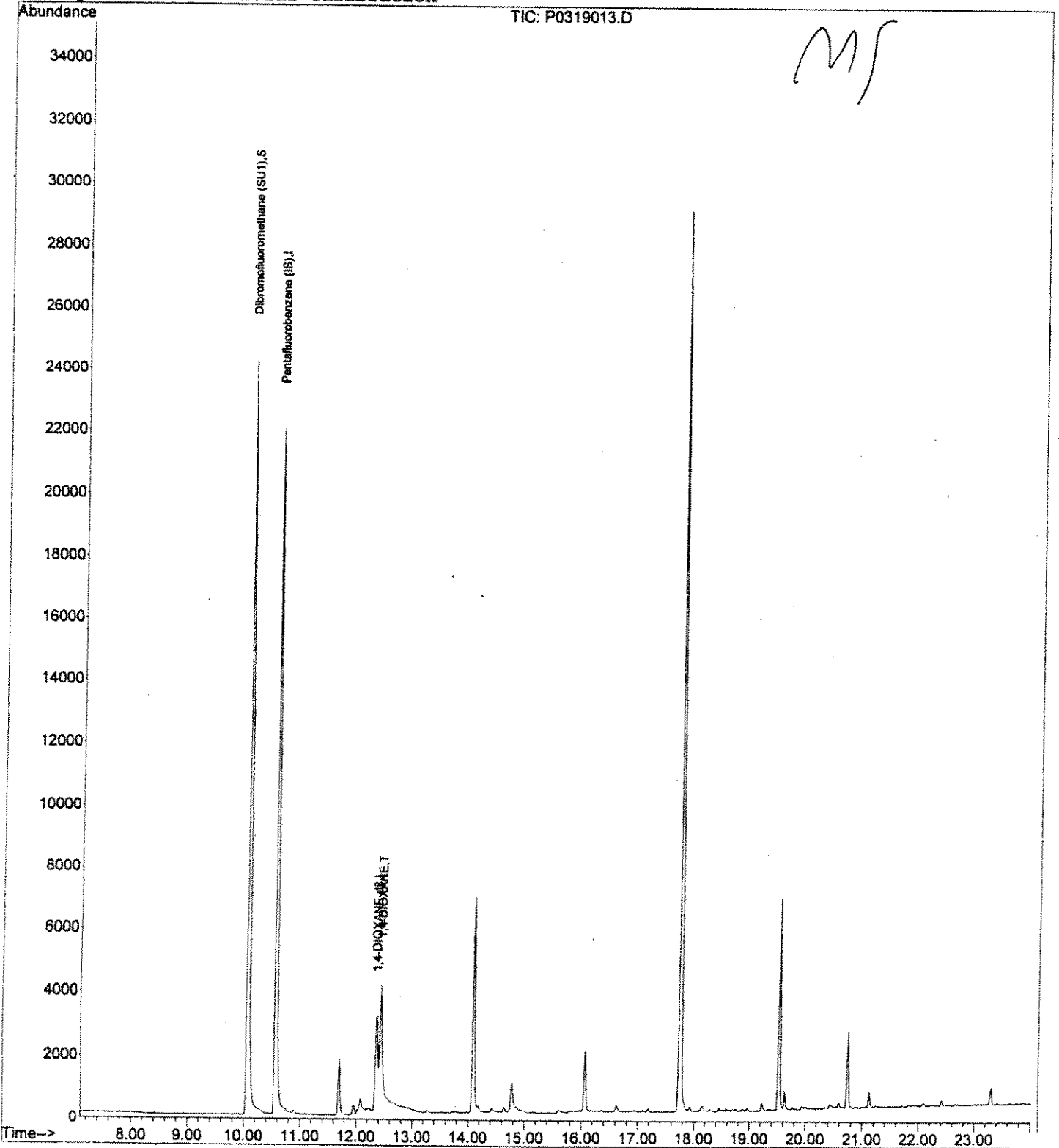
Quant Results File: DX021605.RES

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)

Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)

Last Update : Wed Feb 16 15:53:54 2005

Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\0319014.D  
 Acq On : 19 Mar 2005 1:38 pm  
 Sample : 50.0 PPB CAL  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 14:18 2005

Vial: 14  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX021605.RES

Quant Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Wed Feb 16 15:53:54 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*MS 3/19/05*

| Internal Standards           | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|------------------------------|-------|------|----------|-------|-------|----------|
| 1) Pentafluorobenzene (IS)   | 10.56 | 99   | 47704    | 1.00  | ug/L  | 0.00     |
| 3) 1,4-DIOXANE-d8            | 12.35 | 64   | 5034     | 25.00 | ug/L  | 0.00     |
| 5) 1,2,3-Trichloropropane-d5 | 0.00  | 79   | 0NT      | 0.00  | ug/L  | -15.08   |

| System Monitoring Compounds   | R.T.  | QIon  | Response | Conc     | Units | Dev(Min) |
|-------------------------------|-------|-------|----------|----------|-------|----------|
| 2) Dibromofluoromethane (SU1) | 10.07 | 113   | 164450   | 4.46     | ug/L  | 0.00     |
| Spiked Amount                 | 1.000 | Range | 80 - 120 | Recovery | =     | 446.00%# |

| Target Compounds | R.T.  | QIon | Response | Conc  | Units | Qvalue |
|------------------|-------|------|----------|-------|-------|--------|
| 4) 1,4-DIOXANE   | 12.43 | 88   | 18344    | 58.04 | ug/L  | 99     |

*gubing*

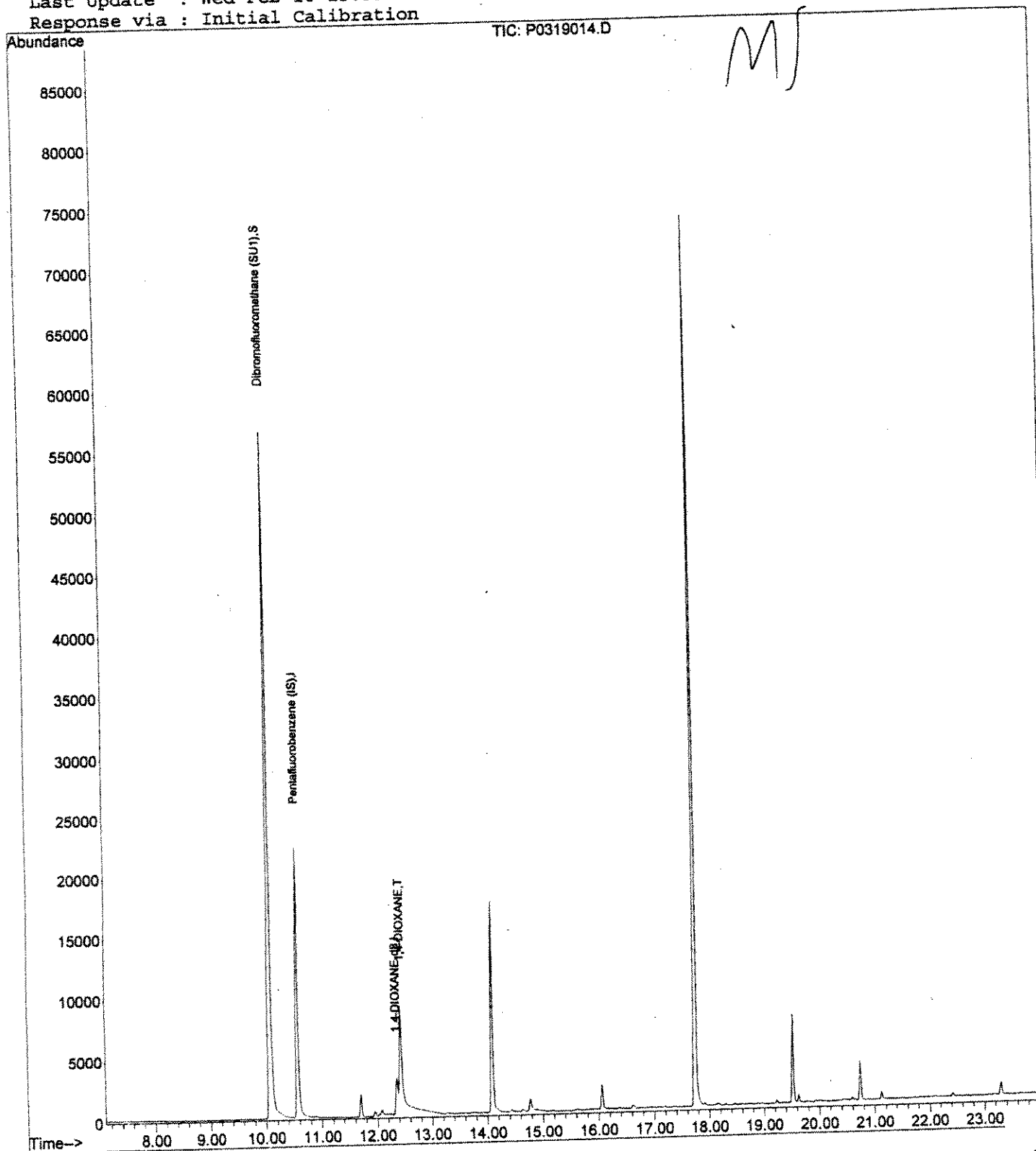
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\0319014.D  
Acq On : 19 Mar 2005 1:38 pm  
Sample : 50.0 PPB CAL  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 19 14:18 2005

Vial: 14  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

Quant Results File: DX021605.RES

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
Last Update : Wed Feb 16 15:53:54 2005  
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\0319015.D  
 Acq On : 19 Mar 2005 2:11 pm  
 Sample : 100.0 PPB CAL  
 Misc : LX 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 14:54 2005

Vial: 15  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX021605.RES

Quant Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Wed Feb 16 15:53:54 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*M 3/19/05*

| Internal Standards           | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|------------------------------|-------|------|----------|-------|-------|----------|
| 1) Pentafluorobenzene (IS)   | 10.56 | 99   | 48150    | 1.00  | ug/L  | 0.00     |
| 3) 1,4-DIOXANE-d8            | 12.35 | 64   | 5834     | 25.00 | ug/L  | 0.00     |
| 5) 1,2,3-Trichloropropane-d5 | 0.00  | 79   | 0        | 0.00  | ug/L  | -15.08   |

System Monitoring Compounds  
 2) Dibromofluoromethane (SU1) 10.07 113 307967 8.28 ug/L 0.00  
 Spiked Amount 1.000 Range 80 - 120 Recovery = 828.00%#

| Target Compounds | R.T.  | QIon | Response | Conc   | Units | Qvalue |
|------------------|-------|------|----------|--------|-------|--------|
| 4) 1,4-DIOXANE   | 12.43 | 88   | 44445    | 121.87 | ug/L  | 98     |

*3/21/05*

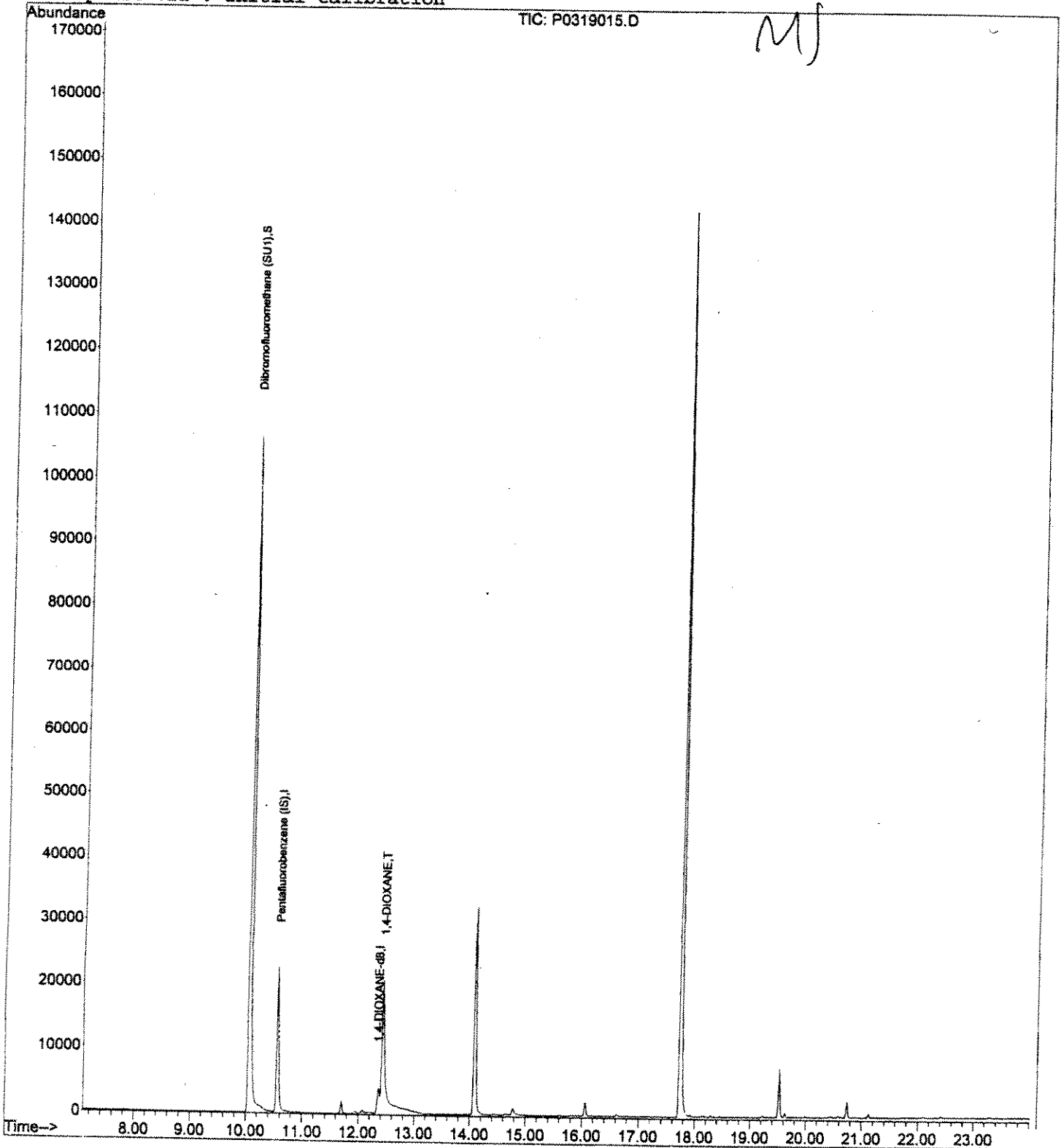
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\0319015.D  
Acq On : 19 Mar 2005 2:11 pm  
Sample : 100.0 PPB CAL  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 19 14:54 2005

Vial: 15  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

Quant Results File: DX021605.RES

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
Last Update : Wed Feb 16 15:53:54 2005  
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\F0319018.D  
 Acq On : 19 Mar 2005 3:54 pm  
 Sample : 1.0 PPB CAL  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 21 7:48 2005

Vial: 18  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX021605.RES

Quant Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Wed Feb 16 15:53:54 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*3/21/05  
JG*

| Internal Standards           | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|------------------------------|-------|------|----------|-------|-------|----------|
| 1) Pentafluorobenzene (IS)   | 10.56 | 99   | 42387    | 1.00  | ug/L  | 0.00     |
| 3) 1,4-DIOXANE-d8            | 12.35 | 64   | 6173     | 25.00 | ug/L  | 0.00     |
| 5) 1,2,3-Trichloropropane-d5 | 0.00  | 79   | 0        | 0.00  | ug/L  | -15.08   |

System Monitoring Compounds  
 2) Dibromofluoromethane (SU1) 10.07 113 3733 0.11 ug/L 0.00  
 Spiked Amount 1.000 Range 80 - 120 Recovery = 11.00%#

| Target Compounds          | R.T.  | QIon | Response | Conc | Units | Qvalue |
|---------------------------|-------|------|----------|------|-------|--------|
| 4) 1,4-DIOXANE            | 12.43 | 88   | 668      | 1.24 | ug/L  | 97     |
| 6) 1,2,3-Trichloropropane | 0.00  | 75   | 0        | N.D. |       |        |

*3/21/05  
JG*

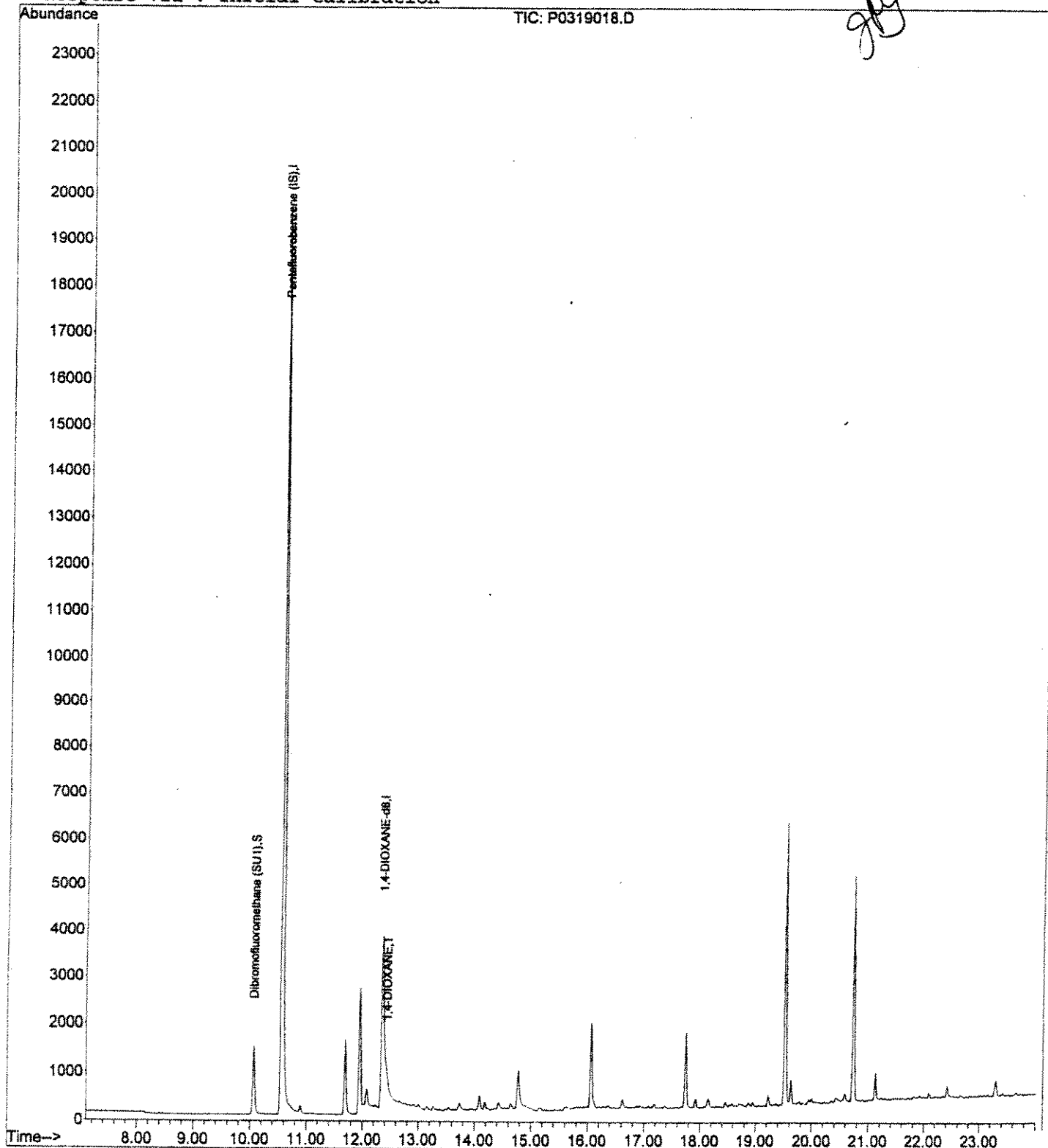
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\0319018.D  
Acq On : 19 Mar 2005 3:54 pm  
Sample : 1.0 PPB CAL  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 21 7:48 2005

Vial: 18  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

Quant Results File: DX021605.RES

Method : D:\HPCHEM\1\METHODS\DX021605.M (RTE Integrator)  
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
Last Update : Wed Feb 16 15:53:54 2005  
Response via : Initial Calibration



Calibration Status Report GCMS1

Method : D:\HPCHEM\1\METHODS\DX031905.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Mon Mar 21 12:54:07 2005  
 Response via : Initial Calibration

*3/21/05  
 Jly*

| # | ID  | Conc | ISTD<br>Conc | Path\File                          |
|---|-----|------|--------------|------------------------------------|
| 1 | 1   | 0    | 1            | D:\HPCHEM\1\DATA\031905\P0319018.D |
| 2 | 2   | 0    | 1            | D:\HPCHEM\1\DATA\031905\P0319010.D |
| 3 | 5   | 1    | 1            | D:\HPCHEM\1\DATA\031905\P0319011.D |
| 4 | 10  | 1    | 1            | D:\HPCHEM\1\DATA\031905\P0319012.D |
| 5 | 20  | 2    | 1            | D:\HPCHEM\1\DATA\031905\P0319013.D |
| 6 | 50  | 5    | 1            | D:\HPCHEM\1\DATA\031905\P0319014.D |
| 7 | 100 | 10   | 1            | D:\HPCHEM\1\DATA\031905\P0319015.D |

| # | ID  | Update Time       | Quant Time         | Acquisition Time     |
|---|-----|-------------------|--------------------|----------------------|
| 1 | 1   | Mar 21 07:49 2005 | Mar 21 07:48 19105 | 19 Mar 2005 3:54 pm  |
| 2 | 2   | Mar 19 14:55 2005 | Mar 19 13:43 19105 | 19 Mar 2005 11:26 am |
| 3 | 5   | Mar 19 14:55 2005 | Mar 19 13:43 19105 | 19 Mar 2005 11:59 am |
| 4 | 10  | Mar 19 14:55 2005 | Mar 19 13:37 19105 | 19 Mar 2005 12:32 pm |
| 5 | 20  | Mar 19 14:55 2005 | Mar 19 13:37 19105 | 19 Mar 2005 1:05 pm  |
| 6 | 50  | Mar 19 14:55 2005 | Mar 19 14:18 19105 | 19 Mar 2005 1:38 pm  |
| 7 | 100 | Mar 19 14:55 2005 | Mar 19 14:54 19105 | 19 Mar 2005 2:11 pm  |

DX031905.M

Mon Mar 21 12:55:30 2005

GCMS1

*3/22/05  
 Jly*

Compound List Report GCMS1

Method : D:\HPCHEM\1\METHODS\DX031905.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Mon Mar 21 12:54:07 2005  
 Response via : Initial Calibration  
 Total Cpnds : 6

3/21/05  
 gky

| PK# | Compound Name              | QIon | Exp_RT | Rel_RT | Cal | #Qual | A/H | ID |
|-----|----------------------------|------|--------|--------|-----|-------|-----|----|
| 1 I | Pentafluorobenzene (IS)    | 99   | 10.57  | 1.000  | A   | 1     | A   | B  |
| 2 S | Dibromofluoromethane (SU1) | 113  | 10.07  | 0.953  | A   | 0     | A   | B  |
| 3 I | 1,4-DIOXANE-d8             | 64   | 12.35  | 1.000  | A   | 1     | A   | B  |
| 4 T | 1,4-DIOXANE                | 88   | 12.43  | 1.007  | L   | 2     | A   | B  |
| 5 I | 1,2,3-Trichloropropane-d5  | 79   | 15.08  | 1.000  | A   | 2     | A   | B  |
| 6 T | 1,2,3-Trichloropropane     | 75   | 15.08  | 1.000  | A   | 2     | A   | B  |

Cal A = Average L = Linear LO = Linear w/origin Q = Quad QC = Quad w/origin  
 #Qual = number of qualifiers  
 A/H = Area or Height  
 ID R = R.T. B = R.T. & Q Q = Qvalue L = Largest A = All

DX031905.M Mon Mar 21 12:55:24 2005 GCMS1

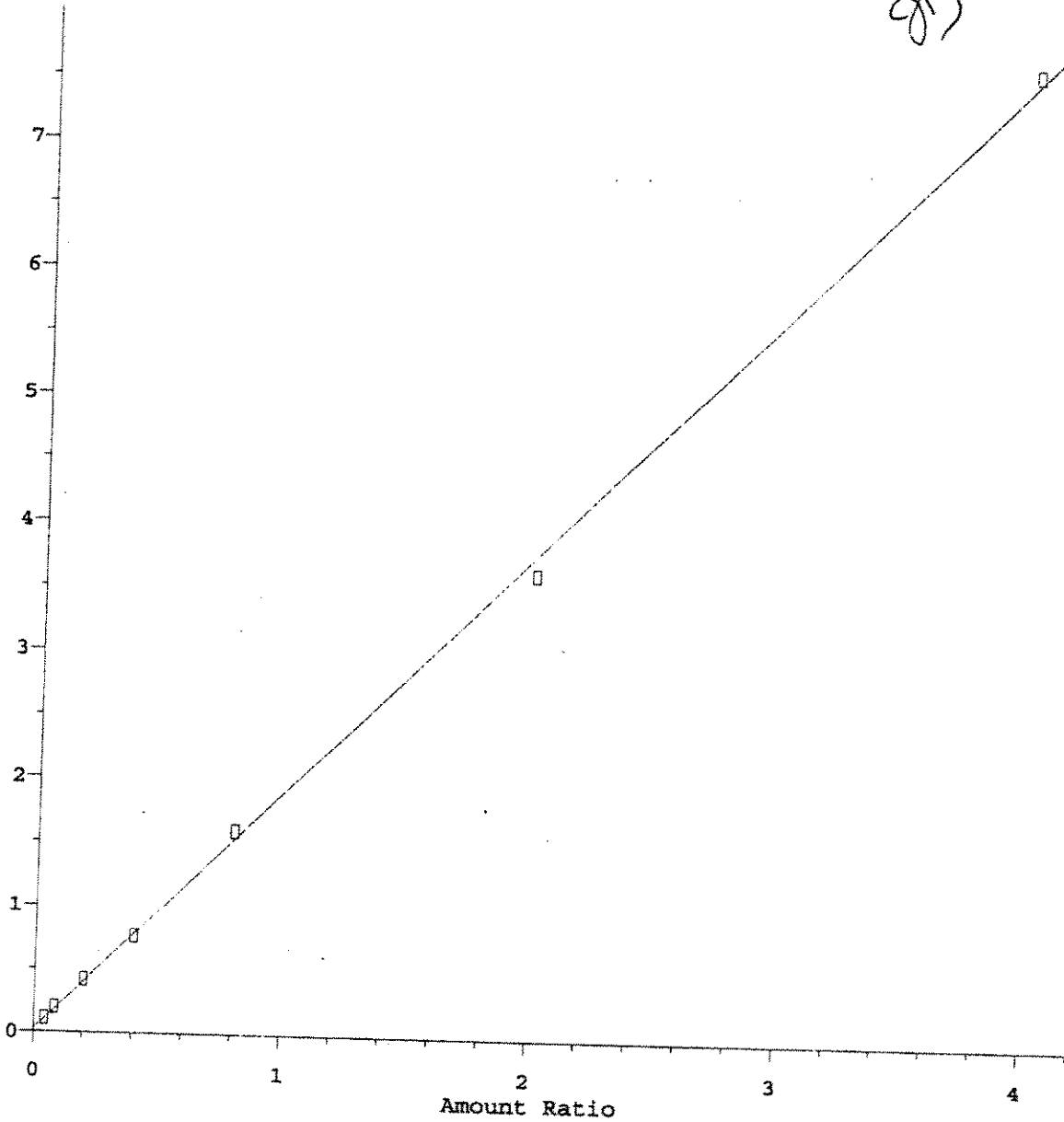
3/22/05



3/21/05  
gfy

Response Ratio

1,4-DIOXANE



Resp Ratio =  $1.88e+000 * Amt + 2.76e-002$   
Coef of Det ( $r^2$ ) = 0.999 Curve Fit: Linear

Method Name: D:\HPCHEM\1\METHODS\DX031905.M  
Calibration Table Last Updated: Mon Mar 21 12:54:07 2005

3/22/05

Response Factor Report GCMS1

Method : D:\HPCHEM\1\METHODS\DX031905.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Mon Mar 21 12:54:07 2005  
 Response via : Initial Calibration

3/21/05  
*[Signature]*

Calibration File

1 =P0319018.D 2 =P0319010.D 5 =P0319011.D 10 =P0319012.D  
 20 =P0319013.D 50 =P0319014.D 100 =P0319015.D

| Compound                        | 1 | 2 | 5 | 10 | 20 | 50 | 100 | Avg | %RSD |
|---------------------------------|---|---|---|----|----|----|-----|-----|------|
| 1) I Pentafluorobenzene (IS)    |   |   |   |    |    |    |     |     |      |
| 2) S Dibromofluoromethane (SU1) |   |   |   |    |    |    |     |     |      |
| 3) I 1,4-DIOXANE-d8             |   |   |   |    |    |    |     |     |      |
| 4) T 1,4-DIOXANE                |   |   |   |    |    |    |     |     |      |
| 5) I 1,2,3-Trichloropropane-d5  |   |   |   |    |    |    |     |     |      |
| 6) T 1,2,3-Trichloropropane     |   |   |   |    |    |    |     |     |      |
| (#) = Out of Range              |   |   |   |    |    |    |     |     |      |

DX031905.M Tue Mar 22 12:15:58 2005 GCMS1

3/21/05  
*[Signature]*

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\0319018.D  
 Acq On : 19 Mar 2005 3:54 pm  
 Sample : 1.0 PPB CAL  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 21 12:54 2005

Vial: 18  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

Quant Results File: DX031905.RES

Quant Method : D:\HPCHEM\1\METHODS\DX031905.M (RTE Integrator)  
 Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
 Last Update : Mon Mar 21 12:54:07 2005  
 Response via : Initial Calibration  
 DataAcq Meth : DX021605

*Re-Calc.*

*3/21/05  
JG*

| Internal Standards                 | R.T.  | QIon           | Response   | Conc    | Units | Dev(Min) |
|------------------------------------|-------|----------------|------------|---------|-------|----------|
| 1) Pentafluorobenzene (IS)         | 10.56 | 99             | 42387      | 1.00    | ug/L  | 0.00     |
| 3) 1,4-DIOXANE-d8                  | 12.35 | 64             | 6173       | 25.00   | ug/L  | 0.00     |
| 5) 1,2,3-Trichloropropane-d5       | 0.00  | 79             | 0          | 0.00    | ug/L  | -15.08   |
| <b>System Monitoring Compounds</b> |       |                |            |         |       |          |
| 2) Dibromofluoromethane (SU1)      | 10.07 | 113            | 3733       | 0.12    | ug/L  | 0.00     |
| Spiked Amount                      | 1.000 | Range 80 - 120 | Recovery = | 12.00%# |       |          |
| <b>Target Compounds</b>            |       |                |            |         |       |          |
| 4) 1,4-DIOXANE                     | 12.43 | 88             | 668        | 1.07    | ug/L  | 96       |
| 6) 1,2,3-Trichloropropane          | 0.00  | 75             | 0          | N.D.    |       |          |

*5-1.5*

*3/21/05 JG*

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\P0319018.D  
Acq On : 19 Mar 2005 3:54 pm  
Sample : 1.0 PPB CAL  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 21 12:54 2005

Vial: 18  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

Quant Results File: DX031905.RES

Method : D:\HPCHEM\1\METHODS\DX031905.M (RTE Integrator)  
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)  
Last Update : Mon Mar 21 12:54:07 2005  
Response via : Initial Calibration

