

## 1. INTRODUCTION

Task Order Title: NPDES Monitoring  
Contract Task Order #: 313150010  
SDG#: IOD0229  
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 6, 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	IOD0229-01	water	624
Trip Blank	Trip Blank	IOD0229-02	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 04/01/05 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 04/10/05 was associated with the sample analyses in this SDG. The %D for diisopropyl ether (DIPE) exceeded 20% in the continuing calibration dated 04/10/05; 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 (5D10007-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 (5D10007-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.



# Del Mar Analytical

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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: IOD0229

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

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

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers	
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water)					Sampled: 04/04/05					Rev Qual   Qual Code
Reporting Units: ug/l										
1,2-Dibromoethane (EDB)	EPA 624	5D10007	0.32	2.0	ND	1	04/10/05	04/10/05	u	
Methyl-tert-butyl Ether (MTBE)	EPA 624	5D10007	0.32	5.0	ND	1	04/10/05	04/10/05	u	
1,2,3-Trichloropropane	EPA 624	5D10007	0.85	10	ND	1	04/10/05	04/10/05	u	
Di-isopropyl Ether (DIPE)	EPA 624	5D10007	0.25	5.0	ND	1	04/10/05	04/10/05	u	
tert-Butanol (TBA)	EPA 624	5D10007	3.1	25	ND	1	04/10/05	04/10/05	u	
Surrogate: Dibromofluoromethane (80-120%)					103 %					
Surrogate: Toluene-d8 (80-120%)					102 %					
Surrogate: 4-Bromofluorobenzene (80-120%)					104 %					
Sample ID: IOD0229-02 (DRAFT: Trip Blank - Water)					Sampled: 04/04/05					Rev Qual   Qual Code
Reporting Units: ug/l										
1,2-Dibromoethane (EDB)	EPA 624	5D10007	0.32	2.0	ND	1	04/10/05	04/10/05	u	
Methyl-tert-butyl Ether (MTBE)	EPA 624	5D10007	0.32	5.0	ND	1	04/10/05	04/10/05	u	
1,2,3-Trichloropropane	EPA 624	5D10007	0.85	10	ND	1	04/10/05	04/10/05	u	
Di-isopropyl Ether (DIPE)	EPA 624	5D10007	0.25	5.0	ND	1	04/10/05	04/10/05	u	
tert-Butanol (TBA)	EPA 624	5D10007	3.1	25	ND	1	04/10/05	04/10/05	u	
Surrogate: Dibromofluoromethane (80-120%)					109 %					
Surrogate: Toluene-d8 (80-120%)					111 %					
Surrogate: 4-Bromofluorobenzene (80-120%)					107 %					

**AMEC VALIDATED**

**LEVEL IV**

DRAFT REPORT  
 DRAFT REPORT  
 DATA SUBJECT TO CHANGE

u  
 05/06/05

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.

**CONTRACT COMPLIANCE SCREENING FORM FOR HARDCOPY DATA**

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

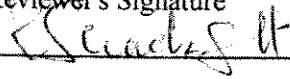
Package ID T711VO98  
 Task Order 313150010  
 SDG No. IOD0229

No. of Analyses 1

Laboratory Del Mar Analytical

Reviewer K. Shadowlight

Analysis/Method 1,4-Dioxane by 8260

Date May 11, 2005  
 Reviewer's Signature  


<b>ACTION ITEMS*</b>	
<b>1. Case Narrative</b>	
<b>Deficiencies</b>	
<b>2. Out of Scope</b>	
<b>Analyses</b>	
<b>3. Analyses Not Conducted</b>	
<b>4. Missing Hardcopy</b>	
<b>Deliverables</b>	
<b>5. Incorrect Hardcopy</b>	
<b>Deliverables</b>	
<b>6. Deviations from Analysis</b>	<u>Qualification was assigned for a %D continuing calibration outlier</u>
GC/MS Tune/Inst. Perform	
Calibrations	
Blanks	
Surrogates	
Matrix Spike/Dup LCS	
Field QC	
Internal Standard Performance	
Compound Identification and	
Quantitation	
System Performance	

**COMMENTS<sup>b</sup>**

<sup>a</sup> Subcontracted analytical laboratory is not meeting contract and/or method requirements.  
<sup>b</sup> Differences in protocol have been adopted by the laboratory but no action against the laboratory is required.

### 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: IOD0229

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 #: IOD0229  
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 8, 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. Del Mar, CA	Lab No. Del Mar, AZ	Matrix	Method
Outfall 012	Outfall 012	IOD0229-01	POD0176-01	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. 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. The EPA ID was added to the sample result summary report by the reviewer. 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 (P5D1317-BS1) from the same analysis. As the analysis cannot be reported as both a CCV and a blank spike, the reviewer reported P5D1317-BS1 as the continuing calibration. The RRF for 1,4-dioxane was  $\geq 0.05$  and the %D was  $\leq 20\%$ . 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 qualifications were required.

## 2.4 BLANKS

One water method blank (P5D1317-BLK1) was associated with this SDG. Target compound 1,4-dioxane was not detected 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 (P5D1317-BS1/BS1D) with this SDG; however, P5D1317-BS1 was reported as the CCV (see section 2.3); therefore, P5D1317-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.





### 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: GENERAL MINERALS

SAMPLE DELIVERY GROUP: IOD0229

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 #: IOD0229  
Project Manager: B. McIlvaine  
Matrix: Water  
Analysis: General Minerals  
QC Level: Level IV  
No. of Samples: 1  
Reviewer: L. Jarusewic  
Date of Review: May 10, 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 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	IOD0229-01	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 presented 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 48-hour holding time for turbidity was exceeded; therefore, turbidity detected was qualified as estimated, "J." The 28-day analytical holding time for ammonia, oil and grease, and total recoverable hydrocarbons, the 7-day holding time for total dissolved solids and total suspended solids, and the 48-hour holding time for total settleable solids and BOD were met. No further qualifications were required.

### 2.2 CALIBRATION

For the applicable analyses, the initial calibration correlation coefficients were  $\geq 0.995$ . The 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 suspended solids, total settleable solids, or oil and grease. No qualifications were required.

### 2.3 BLANKS

Turbidity was detected in a bracketing CCB at 0.040 NTU; however, the turbidity CCB result was insufficient to qualify Outfall 012. The remaining 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 and laboratory control sample duplicate (total recoverable hydrocarbons, BOD, and oil and grease only) recoveries were within the laboratory-established control limits. The LCS is not applicable to total 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 assessed based on LCS results. No qualifications were required.

## 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 sample in this data package. 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. 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: IOD0229

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

**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: IOD0229-01 (DRAFT: Outfall 012 - Water)					Sampled: 04/04/05					REV QUAL QUAL CODE
Reporting Units: mg/l										
Total Recoverable Hydrocarbons	EPA 418.1	5D07088	0.31	1.0	7.7	1	04/07/05	04/07/05		

**AMEC VALIDATED**

**LEVEL III**

DRAFT REPORT  
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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: IOD0229

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

## DRAFT: INORGANICS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water) - cont.					Sampled: 04/04/05				REV QUAL
Reporting Units: mg/l									
Ammonia-N (Distilled)	EPA 350.2	5D08083	0.30	0.50	ND	1	04/08/05	04/08/05	U
Biochemical Oxygen Demand	EPA 405.1	5D06055	0.59	2.0	3.6	1	04/06/05	04/11/05	
Oil & Grease	EPA 413.1	5D11073	0.94	5.0	5.5	1	04/11/05	04/11/05	
Total Dissolved Solids	SM2540C	5D07121	10	10	220	1	04/07/05	04/07/05	
Total Suspended Solids	EPA 160.2	5D07079	10	10	20	1	04/07/05	04/07/05	
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water)					Sampled: 04/04/05				
Reporting Units: ml/hr									
Total Settleable Solids	EPA 160.5	5D06057	0.10	0.10	ND	1	04/06/05	04/06/05	U
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water)					Sampled: 04/04/05				
Reporting Units: NTU									
Turbidity	EPA 180.1	5D07071	0.040	1.0	32	1	04/07/05	04/07/05	J H H
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water)					Sampled: 04/04/05				
Reporting Units: ug/l									
Perchlorate	EPA 314.0	5D06050	0.80	4.0	ND	1	04/06/05	04/06/05	*

# AMEC VALIDATED

# LEVEL IV

\*Analysis Not Validated

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.	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: PERCHLORATE

SAMPLE DELIVERY GROUP: IOD0229

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 #: IOD0229  
Project Manager: B. McIlvaine  
Matrix: Water  
Analysis: Perchlorate  
QC Level: Level IV  
No. of Samples: 1  
Reviewer: L. Jarusewic  
Date of Review: May 10, 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	IOD0229-01	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 result reported on the summary form and in the raw data for the blank analysis associated with the sample was a nondetect at the reporting limit. Perchlorate was detected in a bracketing CCB at  $0.85 \mu\text{g/L}$ ; however, as perchlorate was not detected in the sample, 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: IOD0229

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

## DRAFT: INORGANICS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers	
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water) - cont.					Sampled: 04/04/05					PEP QUAL CODE
Reporting Units: mg/l										
Ammonia-N (Distilled)	EPA 350.2	5D08083	0.30	0.50	ND	1	04/08/05	04/08/05	*	
Biochemical Oxygen Demand	EPA 405.1	5D06055	0.59	2.0	3.6	1	04/06/05	04/11/05		
Oil & Grease	EPA 413.1	5D11073	0.94	5.0	5.5	1	04/11/05	04/11/05		
Total Dissolved Solids	SM2540C	5D07121	10	10	220	1	04/07/05	04/07/05		
Total Suspended Solids	EPA 160.2	5D07079	10	10	20	1	04/07/05	04/07/05		
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water)					Sampled: 04/04/05					
Reporting Units: ml/l/hr										
Total Settleable Solids	EPA 160.5	5D06057	0.10	0.10	ND	1	04/06/05	04/06/05		
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water)					Sampled: 04/04/05					
Reporting Units: NTU										
Turbidity	EPA 180.1	5D07071	0.040	1.0	32	1	04/07/05	04/07/05	H	
Sample ID: IOD0229-01 (DRAFT: Outfall 012 - Water)					Sampled: 04/04/05					
Reporting Units: ug/l										
Perchlorate	EPA 314.0	5D06050	0.80	4.0	ND	1	04/06/05	04/06/05	u	

**AMEC VALIDATED  
 LEVEL IV**

\*Analysis Not Validated

DRAFT REPORT  
 DRAFT REPORT  
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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: 04/13/05  
Received: 04/13/05  
Issued: 05/20/05 11:30

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
IOD0948-01	Outfall 012	Water
IOD0948-02	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: IOD0948

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

**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: IOD0948-01 (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	

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: IOD0948

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

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

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IOD0948-01 (Outfall 012 - Water) - cont.</b>									
Reporting Units: mg/l									
EFH (C13 - C22)	EPA 8015B	5D15050	0.082	0.50	1.7	0.99	04/15/05	04/16/05	
Surrogate: n-Octacosane (40-125%)					67 %				

Del Mar Analytical, Irvine  
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Project Manager



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Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD0948

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

### 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: IOD0948-01 (Outfall 012 - Water) - cont.</b>									
Reporting Units: mg/l									
GRO (C4 - C12)	EPA 8015 Mod.	5D21046	1.0	2.0	2.5	20	04/21/05	04/21/05	
Surrogate: 4-BFB (FID) (65-140%)					100 %				
<b>Sample ID: IOD0948-02 (Trip Blank - Water)</b>									
Reporting Units: mg/l									
GRO (C4 - C12)	EPA 8015 Mod.	5D21046	0.050	0.10	ND	1	04/21/05	04/21/05	
Surrogate: 4-BFB (FID) (65-140%)					79 %				

Del Mar Analytical, Irvine  
Michele Harper  
Project Manager

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Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD0948

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

**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: IOD0948-01 (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	
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%)					113 %				
Surrogate: Toluene-d8 (80-120%)					106 %				
Surrogate: 4-Bromofluorobenzene (80-120%)					106 %				
<b>Sample ID: IOD0948-02 (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	
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 %				

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: IOD0948

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

**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: IOD0948-01 (Outfall 012 - Water)</b>									
Reporting Units: ug/l									
Naphthalene	EPA 625	5D14041	4.5	10	37	0.966	04/14/05	04/19/05	
N-Nitrosodimethylamine	EPA 625	5D14041	3.7	20	ND	0.966	04/14/05	04/19/05	
Surrogate: 2-Fluorophenol (30-120%)					57 %				
Surrogate: Phenol-d6 (35-120%)					60 %				
Surrogate: 2,4,6-Tribromophenol (45-120%)					69 %				
Surrogate: Nitrobenzene-d5 (45-120%)					73 %				
Surrogate: 2-Fluorobiphenyl (45-120%)					69 %				
Surrogate: Terphenyl-d14 (45-120%)					98 %				

**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: IOD0948

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

INORGANICS

Analyte	Method	Batch	MDL Limit	Reporting Limit	Sample Result	Dilution Factor	Date Extracted	Date Analyzed	Data Qualifiers
<b>Sample ID: IOD0948-01 (Outfall 012 - Water) - cont.</b>									
Reporting Units: mg/l									
Ammonia-N (Distilled)	EPA 350.2	5D14077	0.30	0.50	ND	1	04/14/05	04/14/05	
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	J
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	
<b>Sample ID: IOD0948-01 (Outfall 012 - Water)</b>									
Reporting Units: ml/hr									
Total Settleable Solids	EPA 160.5	5D15069	0.10	0.10	ND	1	04/15/05	04/15/05	
<b>Sample ID: IOD0948-01 (Outfall 012 - Water)</b>									
Reporting Units: NTU									
Turbidity	EPA 180.1	5D15076	0.040	1.0	25	1	04/15/05	04/15/05	
<b>Sample ID: IOD0948-01 (Outfall 012 - Water)</b>									
Reporting Units: ug/l									
Perchlorate	EPA 314.0	5D14043	0.80	4.0	ND	1	04/14/05	04/14/05	

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Attention: Bronwyn Kelly

Project ID: Alfa Outfall 012 - During Test

Report Number: IOD0948

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

### 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
Sample ID: IOD0948-01 (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	
Surrogate: Dibromofluoromethane (80-125%)					117 %				

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

**SHORT HOLD TIME DETAIL REPORT**

Sample ID: Outfall 012 (IOD0948-01) - Water	Hold Time (in days)	Date/Time Sampled	Date/Time Received	Date/Time Extracted	Date/Time Analyzed
EPA 160.5	2	04/13/2005 13:00	04/13/2005 19:00	04/15/2005 07:20	04/15/2005 08:20
EPA 180.1	2	04/13/2005 13:00	04/13/2005 19:00	04/15/2005 08:00	04/15/2005 10:30
EPA 405.1	2	04/13/2005 13:00	04/13/2005 19:00	04/14/2005 18:15	04/19/2005 10:30

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Received: 04/13/05

**METHOD BLANK/QC DATA**

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

Analyte	Result	Reporting Limit	MDL	Units	Spike Level	Source Result	%REC %REC	Limit	RPD	RPD Limit	Data Qualifiers
<b>Batch: 5D21036 Extracted: 04/21/05</b>											
<b>Blank Analyzed: 04/21/2005 (5D21036-BLK1)</b>											
Total Recoverable Hydrocarbons	ND	1.0	0.31	mg/l							
<b>LCS Analyzed: 04/21/2005 (5D21036-BS1)</b>											
Total Recoverable Hydrocarbons	4.48	1.0	0.31	mg/l	5.00		90	65-120			M-NR1
<b>LCS Dup Analyzed: 04/21/2005 (5D21036-BSD1)</b>											
Total Recoverable Hydrocarbons	4.60	1.0	0.31	mg/l	5.00		92	65-120	3	20	



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 Received: 04/13/05

**METHOD BLANK/QC DATA**

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

Analyte	Result	Reporting Limit	MDL	Units	Spike Level	Source Result	%REC %REC	Limit	RPD	RPD Limit	Data Qualifiers
<b>Batch: 5D15050 Extracted: 04/15/05</b>											
<b>Blank Analyzed: 04/15/2005 (5D15050-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	0.0878			mg/l	0.200		44	40-125			
<b>LCS Analyzed: 04/16/2005 (5D15050-BS1)</b>											
EFH (C13 - C40)	0.513	0.50	0.082	mg/l	0.775		66	40-120			<b>M-NRI</b>
Surrogate: n-Octacosane	0.105			mg/l	0.200		52	40-125			
<b>LCS Dup Analyzed: 04/15/2005 (5D15050-BSD1)</b>											
EFH (C13 - C40)	0.511	0.50	0.082	mg/l	0.775		66	40-120	0	25	
Surrogate: n-Octacosane	0.0853			mg/l	0.200		43	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 RPD	Limit Limits	RPD Limit	Data Qualifiers
<b>Batch: 5D21046 Extracted: 04/21/05</b>											
<b>Blank Analyzed: 04/21/2005 (5D21046-BLK1)</b>											
GRO (C4 - C12)	ND	0.10	0.050	mg/l							
Surrogate: 4-BFB (FID)	0.00961			mg/l	0.0100		96		65-140		
<b>LCS Analyzed: 04/21/2005 (5D21046-BS1)</b>											
GRO (C4 - C12)	0.797	0.10	0.050	mg/l	0.800		100		70-140		
Surrogate: 4-BFB (FID)	0.0308			mg/l	0.0300		103		65-140		
<b>Matrix Spike Analyzed: 04/21/2005 (5D21046-MS1) Source: IOD0931-07</b>											
GRO (C4 - C12)	0.291	0.10	0.050	mg/l	0.220	ND	132		60-140		
Surrogate: 4-BFB (FID)	0.0117			mg/l	0.0100		117		65-140		
<b>Matrix Spike Dup Analyzed: 04/21/2005 (5D21046-MSD1) Source: IOD0931-07</b>											
GRO (C4 - C12)	0.213	0.10	0.050	mg/l	0.220	ND	97		60-140	31	20
Surrogate: 4-BFB (FID)	0.00938			mg/l	0.0100		94		65-140		R



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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	RPD	Data Qualifiers
<b>Batch: 5D24007 Extracted: 04/24/05</b>									
<b>Blank Analyzed: 04/24/2005 (5D24007-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	27.1			ug/l	25.0		108	80-120	
Surrogate: Toluene-d8	26.4			ug/l	25.0		106	80-120	
Surrogate: 4-Bromofluorobenzene	25.3			ug/l	25.0		101	80-120	
<b>LCS Analyzed: 04/24/2005 (5D24007-BS1)</b>									
1,2-Dibromoethane (EDB)	26.4	2.0	0.32	ug/l	25.0		106	70-125	
Methyl-tert-butyl Ether (MTBE)	29.0	5.0	0.32	ug/l	25.0		116	55-140	
1,2,3-Trichloropropane	27.9	10	0.85	ug/l	25.0		112	55-130	
Di-isopropyl Ether (DIPE)	29.6	5.0	0.25	ug/l	25.0		118	60-135	
tert-Butanol (TBA)	131	25	3.1	ug/l	125		105	65-135	
Surrogate: Dibromofluoromethane	27.3			ug/l	25.0		109	80-120	
Surrogate: Toluene-d8	26.2			ug/l	25.0		105	80-120	
Surrogate: 4-Bromofluorobenzene	26.0			ug/l	25.0		104	80-120	
<b>Matrix Spike Analyzed: 04/24/2005 (5D24007-MS1)</b>					<b>Source: IOD0931-05</b>				
1,2-Dibromoethane (EDB)	21.0	2.0	0.32	ug/l	25.0	ND	84	65-130	
Methyl-tert-butyl Ether (MTBE)	24.3	5.0	0.32	ug/l	25.0	ND	97	50-150	
1,2,3-Trichloropropane	21.7	10	0.85	ug/l	25.0	ND	87	50-135	
Di-isopropyl Ether (DIPE)	27.2	5.0	0.25	ug/l	25.0	ND	109	60-140	
tert-Butanol (TBA)	128	25	3.1	ug/l	125	ND	102	60-145	
Surrogate: Dibromofluoromethane	27.2			ug/l	25.0		109	80-120	
Surrogate: Toluene-d8	26.3			ug/l	25.0		105	80-120	
Surrogate: 4-Bromofluorobenzene	26.0			ug/l	25.0		104	80-120	



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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 Limits	RPD	RPD Limit	Data Qualifiers
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**Batch: 5D24007 Extracted: 04/24/05**

**Matrix Spike Dup Analyzed: 04/24/2005 (5D24007-MSD1)**

**Source: IOD0931-05**

I,2-Dibromoethane (EDB)	25.2	2.0	0.32	ug/l	25.0	ND	101	65-130	18	25	
Methyl-tert-butyl Ether (MTBE)	28.0	5.0	0.32	ug/l	25.0	ND	112	50-150	14	25	
1,2,3-Trichloropropane	26.8	10	0.85	ug/l	25.0	ND	107	50-135	21	30	
Di-isopropyl Ether (DIPE)	29.0	5.0	0.25	ug/l	25.0	ND	116	60-140	6	25	
tert-Butanol (TBA)	132	25	3.1	ug/l	125	ND	106	60-145	3	25	
Surrogate: Dibromofluoromethane	27.6			ug/l	25.0		110	80-120			
Surrogate: Toluene-d8	26.3			ug/l	25.0		105	80-120			
Surrogate: 4-Bromofluorobenzene	26.1			ug/l	25.0		104	80-120			

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**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 Limit	Data Qualifiers
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**Batch: 5D14041 Extracted: 04/14/05**

**Blank Analyzed: 04/18/2005 (5D14041-BLK1)**

Naphthalene	ND	10	4.5	ug/l						
N-Nitrosodimethylamine	ND	20	3.7	ug/l						
Surrogate: 2-Fluorophenol	104			ug/l	200		52	30-120		
Surrogate: Phenol-d6	119			ug/l	200		60	35-120		
Surrogate: 2,4,6-Tribromophenol	151			ug/l	200		76	45-120		
Surrogate: Nitrobenzene-d5	69.2			ug/l	100		69	45-120		
Surrogate: 2-Fluorobiphenyl	71.4			ug/l	100		71	45-120		
Surrogate: Terphenyl-d14	76.3			ug/l	100		76	45-120		

**LCS Analyzed: 04/18/2005-04/19/2005 (5D14041-BS1)**

Naphthalene	67.6	10	4.5	ug/l	100		68	50-120		
N-Nitrosodimethylamine	53.9	20	3.7	ug/l	100		54	40-120		
Surrogate: 2-Fluorophenol	112			ug/l	200		56	30-120		
Surrogate: Phenol-d6	114			ug/l	200		57	35-120		
Surrogate: 2,4,6-Tribromophenol	147			ug/l	200		74	45-120		
Surrogate: Nitrobenzene-d5	70.1			ug/l	100		70	45-120		
Surrogate: 2-Fluorobiphenyl	67.9			ug/l	100		68	45-120		
Surrogate: Terphenyl-d14	71.5			ug/l	100		72	45-120		

**M-NRI**

**LCS Dup Analyzed: 04/18/2005-04/19/2005 (5D14041-BSD1)**

Naphthalene	71.8	10	4.5	ug/l	100		72	50-120	6	20
N-Nitrosodimethylamine	57.4	20	3.7	ug/l	100		57	40-120	6	20
Surrogate: 2-Fluorophenol	117			ug/l	200		58	30-120		
Surrogate: Phenol-d6	123			ug/l	200		62	35-120		
Surrogate: 2,4,6-Tribromophenol	157			ug/l	200		78	45-120		
Surrogate: Nitrobenzene-d5	74.3			ug/l	100		74	45-120		
Surrogate: 2-Fluorobiphenyl	75.1			ug/l	100		75	45-120		
Surrogate: Terphenyl-d14	72.2			ug/l	100		72	45-120		



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**METHOD BLANK/QC DATA**

**INORGANICS**

Analyte	Result	Reporting Limit	MDL	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Data Qualifiers
<b>Batch: 5D14043 Extracted: 04/14/05</b>											
<b>Blank Analyzed: 04/14/2005 (5D14043-BLK1)</b>											
Perchlorate	ND	4.0	0.80	ug/l							
<b>LCS Analyzed: 04/14/2005 (5D14043-BS1)</b>											
Perchlorate	48.3	4.0	0.80	ug/l	50.0		97	85-115			
<b>Matrix Spike Analyzed: 04/14/2005 (5D14043-MS1)</b>											
						<b>Source: IOD0930-02</b>					
Perchlorate	61.0	4.0	0.80	ug/l	50.0	1.6	119	80-120			
<b>Matrix Spike Dup Analyzed: 04/14/2005 (5D14043-MSD1)</b>											
						<b>Source: IOD0930-02</b>					
Perchlorate	50.6	4.0	0.80	ug/l	50.0	1.6	98	80-120	19	20	
<b>Batch: 5D14060 Extracted: 04/14/05</b>											
<b>Blank Analyzed: 04/19/2005 (5D14060-BLK1)</b>											
Biochemical Oxygen Demand	ND	2.0	0.59	mg/l							
<b>LCS Analyzed: 04/19/2005 (5D14060-BS1)</b>											
Biochemical Oxygen Demand	209	100	30	mg/l	198		106	85-115			
<b>LCS Dup Analyzed: 04/19/2005 (5D14060-BSD1)</b>											
Biochemical Oxygen Demand	207	100	30	mg/l	198		105	85-115	1	20	
<b>Batch: 5D14077 Extracted: 04/14/05</b>											
<b>Blank Analyzed: 04/14/2005 (5D14077-BLK1)</b>											
Ammonia-N (Distilled)	ND	0.50	0.30	mg/l							

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**METHOD BLANK/QC DATA**

**INORGANICS**

Analyte	Result	Reporting Limit	MDL	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Data Qualifiers
<b>Batch: 5D14077 Extracted: 04/14/05</b>											
<b>LCS Analyzed: 04/14/2005 (5D14077-BS1)</b>											
Ammonia-N (Distilled)	10.6	0.50	0.30	mg/l	10.0		106	80-115			
<b>Matrix Spike Analyzed: 04/14/2005 (5D14077-MS1) Source: IOD0922-02</b>											
Ammonia-N (Distilled)	10.6	0.50	0.30	mg/l	10.0	0.56	100	70-120			
<b>Matrix Spike Dup Analyzed: 04/14/2005 (5D14077-MSD1) Source: IOD0922-02</b>											
Ammonia-N (Distilled)	10.4	0.50	0.30	mg/l	10.0	0.56	98	70-120	2	15	
<b>Batch: 5D14081 Extracted: 04/14/05</b>											
<b>Blank Analyzed: 04/14/2005 (5D14081-BLK1)</b>											
Oil & Grease	ND	5.0	0.94	mg/l							
<b>LCS Analyzed: 04/14/2005 (5D14081-BS1) M-NRI</b>											
Oil & Grease	19.8	5.0	0.94	mg/l	20.0		99	65-120			
<b>LCS Dup Analyzed: 04/14/2005 (5D14081-BSD1)</b>											
Oil & Grease	17.5	5.0	0.94	mg/l	20.0		88	65-120	12	20	
<b>Batch: 5D15076 Extracted: 04/15/05</b>											
<b>Blank Analyzed: 04/15/2005 (5D15076-BLK1)</b>											
Turbidity	ND	1.0	0.040	NTU							
<b>Duplicate Analyzed: 04/15/2005 (5D15076-DUP1) Source: IOD1013-01</b>											
Turbidity	0.460	1.0	0.040	NTU		0.48			4	20	J

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**METHOD BLANK/QC DATA**

**INORGANICS**

Analyte	Result	Reporting Limit	MDL	Units	Spike Level	Source Result	%REC %REC	RPD RPD	Limit Limits	RPD Limit	Data Qualifiers
<b>Batch: 5D18089 Extracted: 04/18/05</b>											
<b>Blank Analyzed: 04/18/2005 (5D18089-BLK1)</b>											
Total Suspended Solids	ND	10	10	mg/l							
<b>LCS Analyzed: 04/18/2005 (5D18089-BS1)</b>											
Total Suspended Solids	951	10	10	mg/l	1000		95		85-115		
<b>Duplicate Analyzed: 04/18/2005 (5D18089-DUP1)</b>											
						<b>Source: IOD1125-01</b>					
Total Suspended Solids	12.0	10	10	mg/l		21		55		10	R-4
<b>Batch: 5D18095 Extracted: 04/18/05</b>											
<b>Blank Analyzed: 04/18/2005 (5D18095-BLK1)</b>											
Total Dissolved Solids	ND	10	10	mg/l							
<b>LCS Analyzed: 04/18/2005 (5D18095-BS1)</b>											
Total Dissolved Solids	988	10	10	mg/l	1000		99		90-110		
<b>Duplicate Analyzed: 04/18/2005 (5D18095-DUP1)</b>											
						<b>Source: IOD0830-01</b>					
Total Dissolved Solids	345	10	10	mg/l		350		1		10	

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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: P5D1803 Extracted: 04/18/05</b>											
<b>Blank Analyzed: 04/18/2005 (P5D1803-BLK1)</b>											
1,4-Dioxane	ND	1.0	0.49	ug/l							
Surrogate: Dibromofluoromethane	1.06			ug/l	1.00		106	80-125			
<b>LCS Analyzed: 04/18/2005 (P5D1803-BS1)</b>											
1,4-Dioxane	8.69	1.0	0.49	ug/l	10.0		87	70-130			
Surrogate: Dibromofluoromethane	1.06			ug/l	1.00		106	80-125			
<b>LCS Dup Analyzed: 04/18/2005 (P5D1803-BSD1)</b>											
1,4-Dioxane	8.50	1.0	0.49	ug/l	10.0		85	70-130	2	20	
Surrogate: Dibromofluoromethane	1.07			ug/l	1.00		107	80-125			
<b>Matrix Spike Analyzed: 04/18/2005 (P5D1803-MS1) Source: POD0370-01</b>											
1,4-Dioxane	8.69	1.0	0.49	ug/l	10.0	ND	87	70-150			
Surrogate: Dibromofluoromethane	1.10			ug/l	1.00		110	80-125			
<b>Matrix Spike Dup Analyzed: 04/18/2005 (P5D1803-MSD1) Source: POD0370-01</b>											
1,4-Dioxane	8.17	1.0	0.49	ug/l	10.0	ND	82	70-150	6	25	
Surrogate: Dibromofluoromethane	1.13			ug/l	1.00		113	80-125			

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: IOD0948

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

### DATA QUALIFIERS AND DEFINITIONS

- J** Estimated value. Analyte detected at a level less than the Reporting Limit (RL) and greater than or equal to the Method Detection Limit (MDL). The user of this data should be aware that this data is of unknown quality.
- M-NRI** There was no MS/MSD analyzed with this batch due to insufficient sample volume. See Blank Spike/Blank Spike Duplicate.
- R** The RPD exceeded the method control limit due to sample matrix effects. The individual analyte QA/QC recoveries, however, were within acceptance limits.
- R-4** Due to the low levels of analyte in the sample, the duplicate RPD calculation does not provide useful information.
- 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





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

**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

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

Method Performed: EPA 8260B

Samples: IOD0948-01

**Del Mar Analytical, Irvine**  
Michele Harper  
Project Manager



17461 Derian Ave. Suite 100, Irvine, CA 92614 Ph (949) 261-1022 Fax (949) 261-1228  
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 9484 Chesapeake Drive, Suite 805, San Diego, CA 92123 Ph (619) 505-9596 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 # IOD0948

**SENDING LABORATORY:**

Del Mar Analytical, Irvine  
 17461 Derian Avenue, Suite 100  
 Irvine, CA 92614  
 Phone: (949) 261-1022  
 Fax: (949) 261-1228  
 Project Manager: Michele Harper

**RECEIVING LABORATORY:**

Del Mar Analytical - Phoenix  
 9830 S. 51st Street, Suite B-120  
 Phoenix, AZ 85044  
 Phone : (480) 785-0043  
 Fax: (480) 785-0851

Analysis	Expiration	Due	Comments
Sample ID: IOD0948-01 Water	Sampled: 04/13/05 13:00		
Dioxane-8260B-out	04/27/05 13:00	04/22/05 12:00	Boeing-permit, sub DMAP, J flags, ID=DMA+Outfall 012
Level 4 Data Package - Out	05/11/05 13:00	04/22/05 12:00	Boeing

**Containers Supplied:**

- 40 ml VOA w/HCL (IOD0948-01H)
- 40 ml VOA w/HCL (IOD0948-01I)
- 40 ml VOA w/HCL (IOD0948-01J)

PODO411-01

**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	Samples Received On Ice: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
Custody Seals Present: <input type="checkbox"/> Yes <input type="checkbox"/> No	Samples Preserved Properly: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No	Samples Received at (temp): <u>26°C</u>

~~Released By: Date: 4-14-05 Time: 1700 Received By: Date: 4/15/05 Time: 9:45~~

Released By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_ Received By: Date: 4/15/05 Time: 9:45

(281) I000948

**CHAIN OF CUSTODY FORM**

Del Mar Analytical Version 5 8/12/04

Client Name/Address:				Project:				ANALYSIS REQUIRED												Field readings:	
<b>MWH-Pasadena</b> 300 North Lake Avenue, Suite 1200 Pasadena, CA 91101 Project Manager: Bronwyn Kelly Sampler: <i>Pollock</i>				Boeing-SSFL NPDES <b>During Test -- Outfall 012</b> Alfa Test Stand Phone Number: (626) 568-6691 Fax Number: (626) 568-6515				ANALYSIS REQUIRED												Field readings:	
								Oil & Grease (EPA 413.1)	8015-gas	8015-diesel/et fuel	1,4-Dioxane-8260B	TRPH=Total Rec. (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)	Perchlorate	Turbidity, TDS, TSS	Settleable Solids	Temp = 71.8 pH = 7.4	
Sample Description	Sample Matrix	Container Type	# of Cont.	Sampling Date/Time	Preservative	Bottle #	Oil & Grease (EPA 413.1)	8015-gas	8015-diesel/et fuel	1,4-Dioxane-8260B	TRPH=Total Rec. (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)	Perchlorate	Turbidity, TDS, TSS	Settleable Solids			
Outfall 012	W	1L Amber	1	4-13-05 ~13:00	HCl	1A	X														
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											X				
Trip Blank	W	VOAs	6		HCl	12A, 12B, 12C, 12D, 12E, 12F						X									
Relinquished By				Date/Time:	Received By				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 _____											
Relinquished By				Date/Time:	Received By				Date/Time:	Sample Integrity: (Check) Intact _____ On Ice: _____											
Relinquished By				Date/Time:	Received By				Date/Time:	3C											

*Pollock*

4-13-05 1535

4/13/05 1535

4/13/05 1900

4/13/05 1900



QA/QC DATA PACKAGE: LEVEL IV



QA/QC DATA PACKAGE LEVEL IV

TABLE SUMMARY

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LABORATORY REPORT

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

Project: IOD0948

Sampled: 04/13/05  
Received: 04/15/05  
Issued: 04/21/05 14:57

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

POD0411-01

CLIENT ID

IOD0948-01

MATRIX

Water

- SAMPLE RECEIPT: Samples were received intact, at 3°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, Cotton, CA 92324 Ph (909) 370-4667 Fax (909) 370-1046  
 9484 Chesapeake Drive, Suite 805, San Diego, CA 92123 Ph (619) 505-9596 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 # IOD0948

**SENDING LABORATORY:**

Del Mar Analytical, Irvine  
 17461 Derian Avenue, Suite 100  
 Irvine, CA 92614  
 Phone: (949) 261-1022  
 Fax: (949) 261-1228  
 Project Manager: Michele Harper

**RECEIVING LABORATORY:**

Del Mar Analytical - Phoenix  
 9830 S. 51st Street, Suite B-120  
 Phoenix, AZ 85044  
 Phone : (480) 785-0043  
 Fax: (480) 785-0851

Analysis	Expiration	Due	Comments
<b>Sample ID: IOD0948-01 Water</b>	<b>Sampled: 04/13/05 13:00</b>		
Dioxane-8260B-out	04/27/05 13:00	04/22/05 12:00	Boeing-permit, sub DMAP, J flags, ID=DMA+Outfall 012
Level 4 Data Package - Out	05/11/05 13:00	04/22/05 12:00	Boeing

**Containers Supplied:**

- 40 ml VOA w/HCL (IOD0948-01H)
- 40 ml VOA w/HCL (IOD0948-01I)
- 40 ml VOA w/HCL (IOD0948-01J)

PODO411-01

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	Samples Received On Ice::	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No
Custody Seals Present:	<input type="checkbox"/> Yes <input type="checkbox"/> No	Samples Preserved Properly:	<input checked="" type="checkbox"/> Yes <input type="checkbox"/> No	Samples Received at (temp):	<u>26°C</u>

~~Released By: Date: 4-14-05 Time: 1700 Received By: Date: 4/15/05 Time: 9:45~~

Released By: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_ Received By: Date: 4/15/05 Time: 9:45





QC DATA PACKAGE: LEVEL IV

ANALYTICAL REPORTS



17461 Derian Ave., Suite 100, Irvine, CA 92614 (949) 261-1022 FAX (949) 260-3297  
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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: IOD0948

Report Number: POD0411

Sampled: 04/13/05  
Received: 04/15/05

### 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: POD0411-01 (IOD0948-01 - Water)</b>									
Reporting Units: ug/l									
1,4-Dioxane	EPA 8260B	P5D1803	0.49	1.0	ND	1	04/18/05	04/18/05	
Surrogate: Dibromofluoromethane (80-125%)					117 %				

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

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 1014 E. Cooley Dr., Suite A, Colton, CA 92324 (909) 370-4667 FAX (949) 370-1046  
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 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: IOD0948

Report Number: POD0411

Sampled: 04/13/05  
 Received: 04/15/05

## 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: P5D1803 Extracted: 04/18/05</b>											
<b>Blank Analyzed: 04/18/2005 (P5D1803-BLK1)</b>											
1,4-Dioxane	ND	1.0	0.49	ug/l							
Surrogate: Dibromofluoromethane	1.06			ug/l	1.00		106	80-125			
<b>LCS Analyzed: 04/18/2005 (P5D1803-BS1)</b>											
1,4-Dioxane	8.69	1.0	0.49	ug/l	10.0		87	70-130			
Surrogate: Dibromofluoromethane	1.06			ug/l	1.00		106	80-125			
<b>LCS Dup Analyzed: 04/18/2005 (P5D1803-BSD1)</b>											
1,4-Dioxane	8.50	1.0	0.49	ug/l	10.0		85	70-130	2	20	
Surrogate: Dibromofluoromethane	1.07			ug/l	1.00		107	80-125			
<b>Matrix Spike Analyzed: 04/18/2005 (P5D1803-MS1)</b>											
						<b>Source: POD0370-01</b>					
1,4-Dioxane	8.69	1.0	0.49	ug/l	10.0	ND	87	70-150			
Surrogate: Dibromofluoromethane	1.10			ug/l	1.00		110	80-125			
<b>Matrix Spike Dup Analyzed: 04/18/2005 (P5D1803-MSD1)</b>											
						<b>Source: POD0370-01</b>					
1,4-Dioxane	8.17	1.0	0.49	ug/l	10.0	ND	82	70-150	6	25	
Surrogate: Dibromofluoromethane	1.13			ug/l	1.00		113	80-125			

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.



# 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: IOD0948

Report Number: POD0411

Sampled: 04/13/05  
Received: 04/15/05

## DATA QUALIFIERS AND DEFINITIONS

- 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

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# 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: IOD0948

Report Number: POD0411

Sampled: 04/13/05  
 Received: 04/15/05

## 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



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



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?  
Standard IDs and analyst's initials are present?
2. Initial Calibration met criteria? 
  - a. Print calibration as Average Response Factor  
(624: RSD  $\leq$  35%)  
(8260B:  $\leq$  30% for CCCs,  $\leq$  15% for all other compounds, SPCCs met Criteria)  
(524.2: RSD  $\leq$  20%)
  - b. If non CCC RSD > 15%, print out the curve as Linear Regression  
 $r \geq 0.995$  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  
 $r \geq 0.995$  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  
must be  $\pm 30\%$ , file with the calibration for reference
  - f. If samples contain negative values then:  
compare the area counts with the low standard on file  
if <, then report as N.D. with no flag  
if >, then report from RSD curve and flag that curve is out  
or report at an elevated RL as compared to a curve standard
3. Initial Midpoint / LCS / BFB Tune  
(624: use Table 5) (524.2:  $\pm 30\%$ ) (8260B: see control chart)  
SPCCs met criteria?  CCCs met criteria ( $\pm 20\%$ )?
4. Checked integration of all peaks in Midpoint?
5. Method Blank < Report Limit, if not is data flagged?  
(624: every 20 samples) (524.2: every 12 hours) (8260B: every 12 hours)
6. MS/MSD (every 20 samples)  
(624: use Table 5) (524.2: N/A) (8260B: see Control Chart)
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  
(624: See Table 5) (524.2:  $\pm 30\%$ ) (8260B: See Control Chart)
10. Cont. Midpoint / LCS / BFB Tune done every 12 hours  
(624: use Table 5) (524.2:  $\pm 30\%$ ) (8260B: see control chart)  
SPCCs met criteria?  CCCs met criteria ( $\pm 20\%$ )?
11. Surrogates within acceptance limits  
(624 / 524.2 / 8260B: See Control Chart)
12. Internal Standards within acceptance limits  
(624 / 524.2 / 8260B: response must be -50 to +100%).
13. Manual re-integration(s) performed?  
yes: \_\_\_\_\_ no:
14. Corrective Action Report required?  
yes: \_\_\_\_\_ (Attached) no:
15. Reports impacted by the Corrective Action Report

Analyst: Joe Galassi 3/21/05

Reviewer / Date: [Signature] 3/22/05

# DMAP GC/MS 1 DAILY LOG SUMMARY

CAL CURVE

DATE: 3/19/05

QC BATCH # (s): P501902J63/2/05

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	8260	H2O	
1	2	CCV	1x10ML				
2	3	LCS DUP					-DNU IS LOW
3	4	LCS DUP					-DNU IS LOW -> replace + xp
✓	5	TUNE					
1	6	CCV					-DNU, IS's still low
2	7	CCV					↓ will re-calibrate
3	8	Blank					
4	9	1.0 ppb Cal					DNU's Grubby PPH out of spec
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 up blank / Tune					
12	17	MSCCV Blank					
13	18	1.0 ppb Cal					
14	19	SS/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: JY/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</del> DNU	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.	CCV DNU	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</del> DNU	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</del> DNU	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  
gwy

BFB

Data File : D:\HPCHEM\1\DATA\031905\0319005.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)

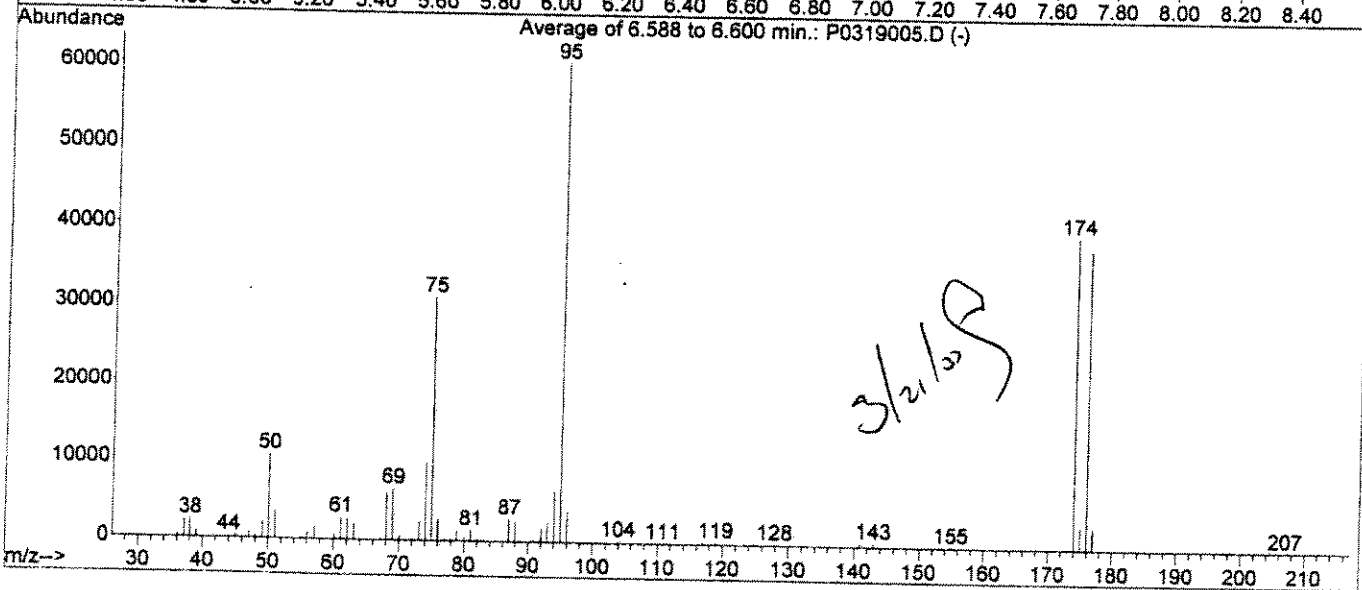
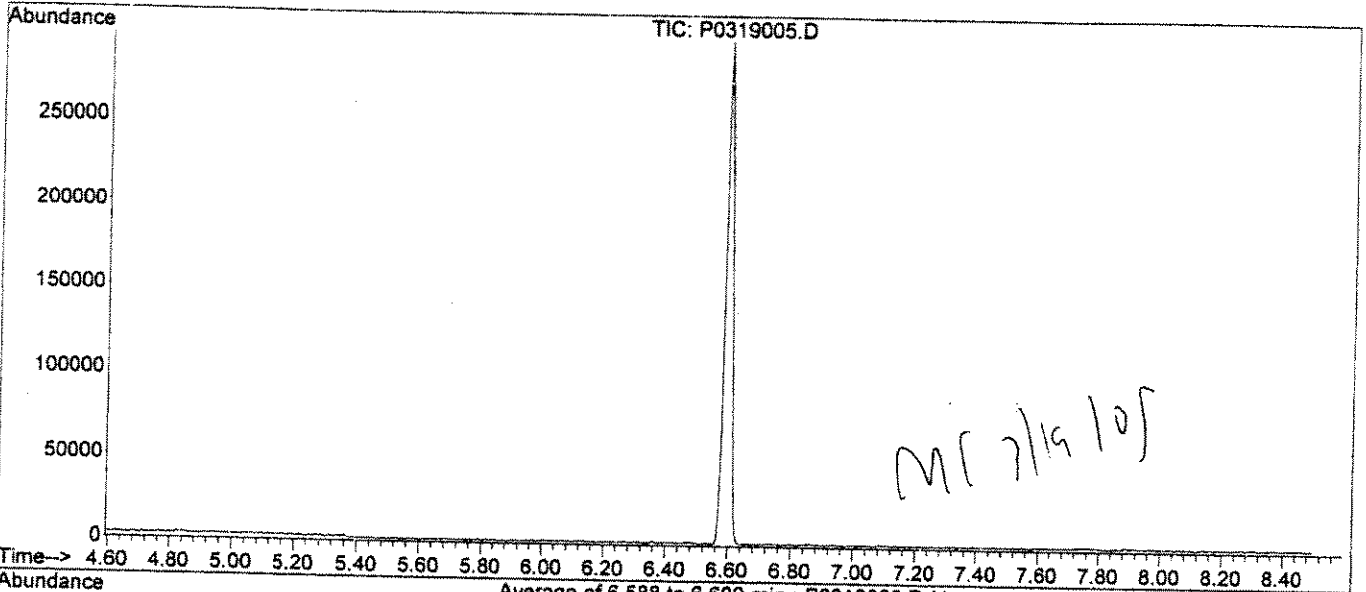
Title : 8260 1,4-Dioxane Ini. Cal. (05/02/02)

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

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\0319008.D  
 Acq On : 19 Mar 2005 10:12 am  
 Sample : BLANK  
 Misc : 1X 10ML  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 10:34 2005

Vial: 8  
 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

*MT 3/19/05*

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene (IS)	10.56	99	46878	1.00	ug/L	0.00
3) 1,4-DIOXANE-d8	12.35	64	6171	25.00	ug/L	0.00
5) 1,2,3-Trichloropropane-d5	0.00	79	0 <sub>NT</sub>	0.00	ug/L	-15.08
System Monitoring Compounds						
2) Dibromofluoromethane (SU1)	10.07	113	37890	1.05	ug/L	0.00
Spiked Amount	1.000	Range 80 - 120	Recovery	=	105.00%	✓
Target Compounds						
4) 1,4-DIOXANE	12.45	88	278	0.23	ug/L	Qvalue <i>ND</i> 92

*3/21/05*

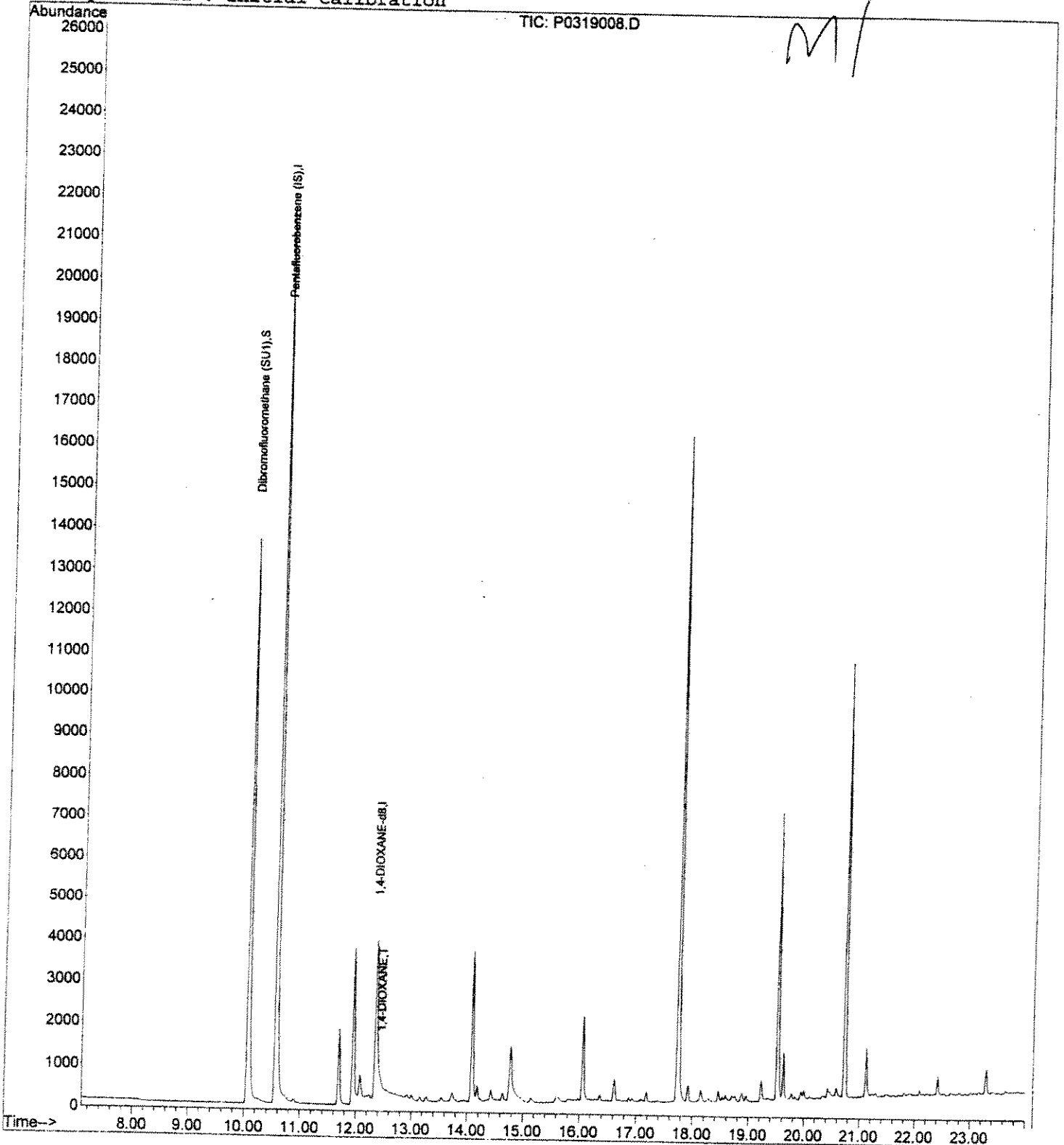
Quantitation Report

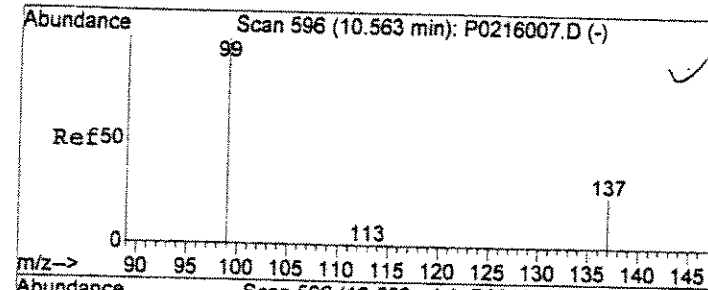
Data File : D:\HPCHEM\1\DATA\031905\P0319008.D  
Acq On : 19 Mar 2005 10:12 am  
Sample : BLANK  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 19 10:34 2005

Vial: 8  
Operator: JG/MS/CLS  
Inst : GCMS1  
Multiplr: 1.00

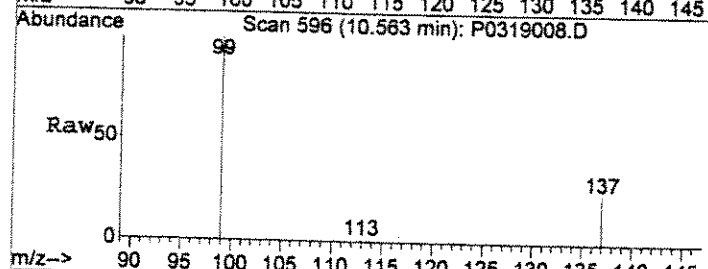
Quant Results File: DX021605.RES

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

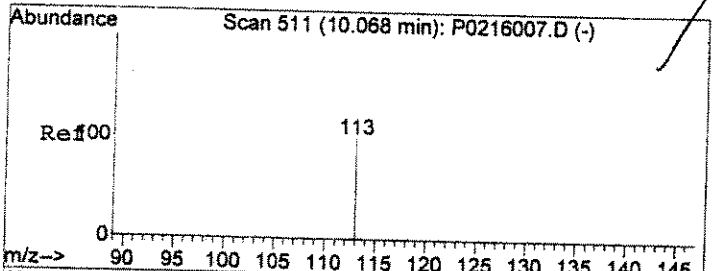
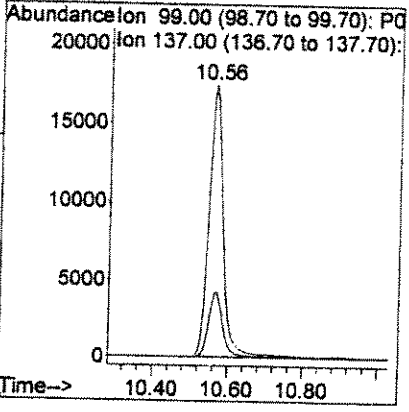
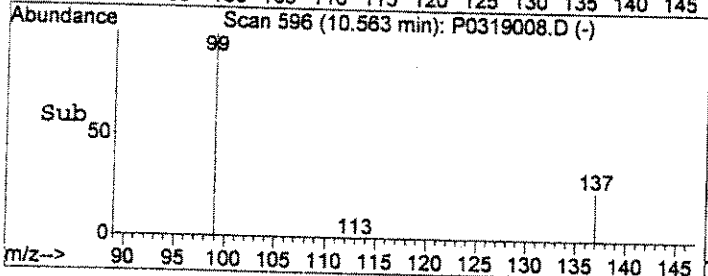




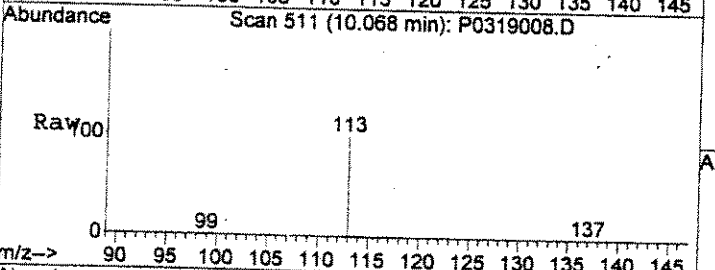
#1  
 Pentafluorobenzene (IS)  
 Concen: 1.00 ug/L  
 RT: 10.56 min Scan# 596  
 Delta R.T. 0.00 min  
 Lab File: P0319008.D  
 Acq: 19 Mar 2005 10:12 am



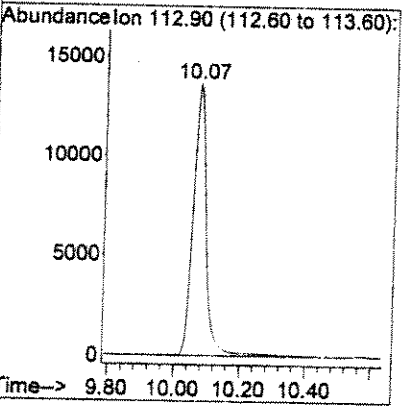
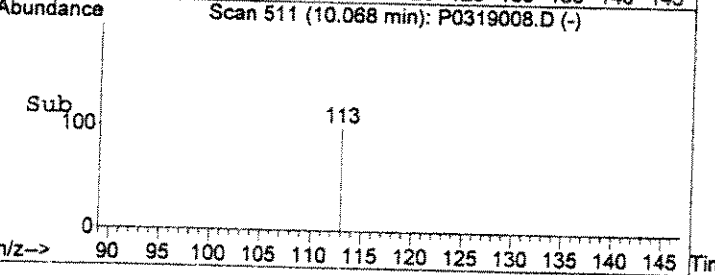
Tgt Ion: 99 Resp: 46878  
 Ion Ratio Lower Upper  
 99 100  
 137 23.9 3.7 43.7



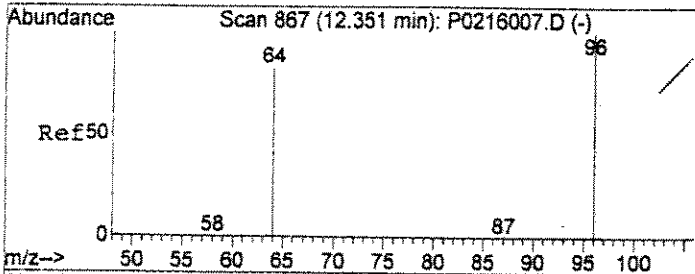
#2  
 Dibromofluoromethane (SU1)  
 Concen: 1.00 ug/L  
 RT: 10.07 min Scan# 511  
 Delta R.T. 0.00 min  
 Lab File: P0319008.D  
 Acq: 19 Mar 2005 10:12 am



Tgt Ion: 113 Resp: 37890

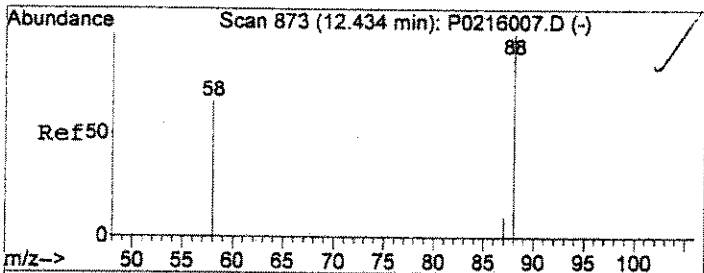
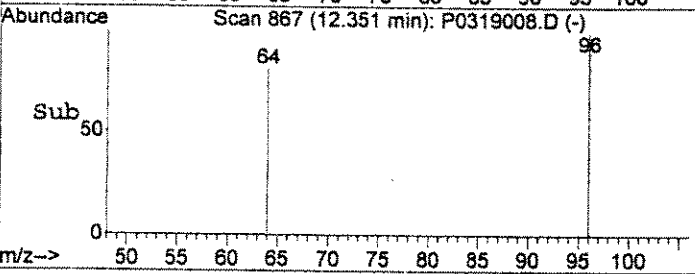
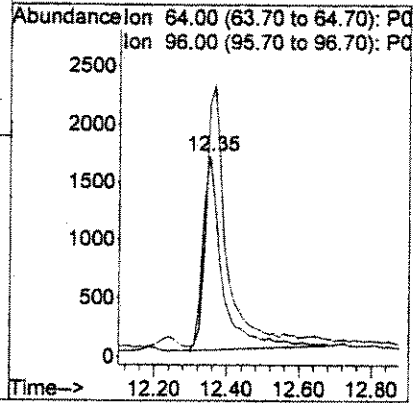
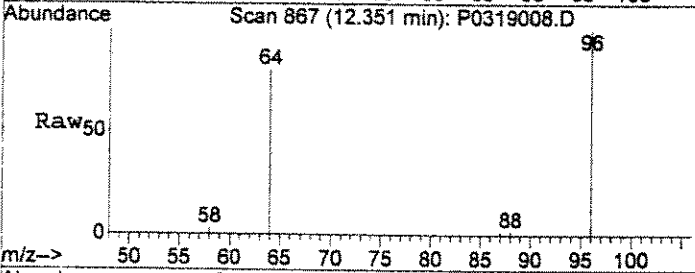


8



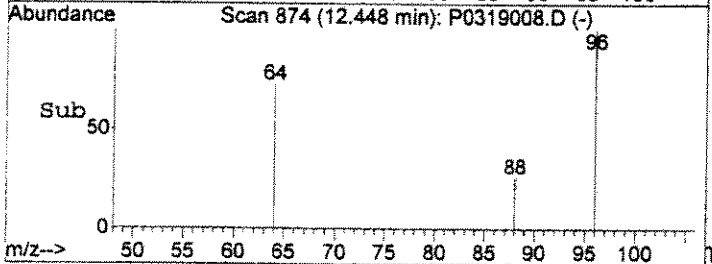
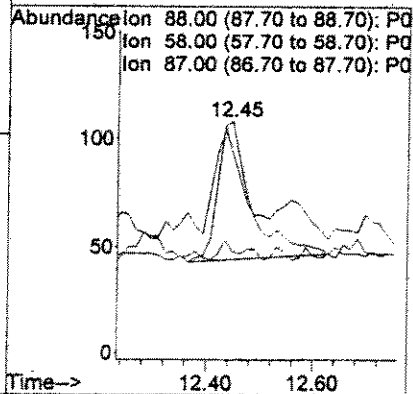
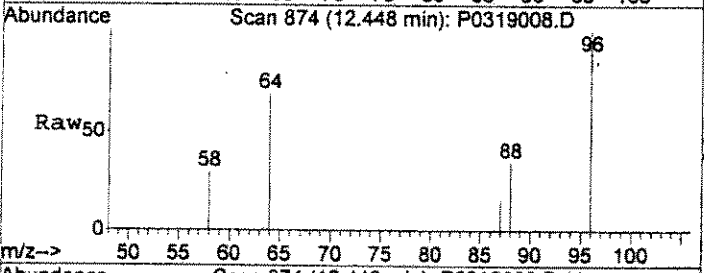
#3  
 1,4-DIOXANE-d8  
 Concen: 25.00 ug/L  
 RT: 12.35 min Scan# 867  
 Delta R.T. -0.00 min  
 Lab File: P0319008.D  
 Acq: 19 Mar 2005 10:12 am

Tgt Ion: 64 Resp: 6171  
 Ion Ratio Lower Upper  
 64 100  
 96 123.7 70.1 170.1



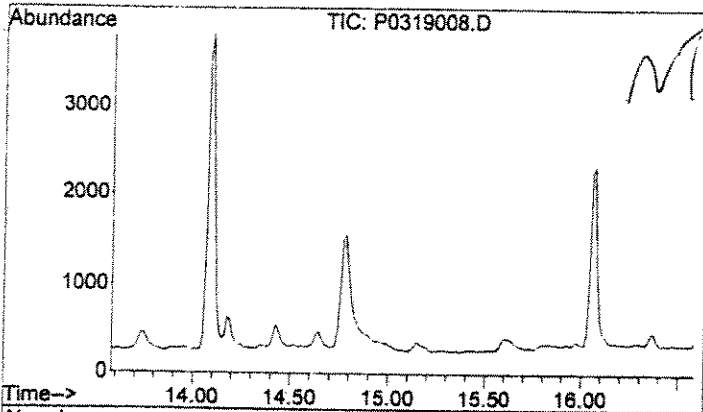
#4  
 1,4-DIOXANE  
 Concen: 0.23 ug/L  
 RT: 12.45 min Scan# 874  
 Delta R.T. 0.01 min  
 Lab File: P0319008.D  
 Acq: 19 Mar 2005 10:12 am

Tgt Ion: 88 Resp: 278  
 Ion Ratio Lower Upper  
 88 100  
 58 61.5 16.3 116.3  
 87 3.1 0.0 59.9



*S*

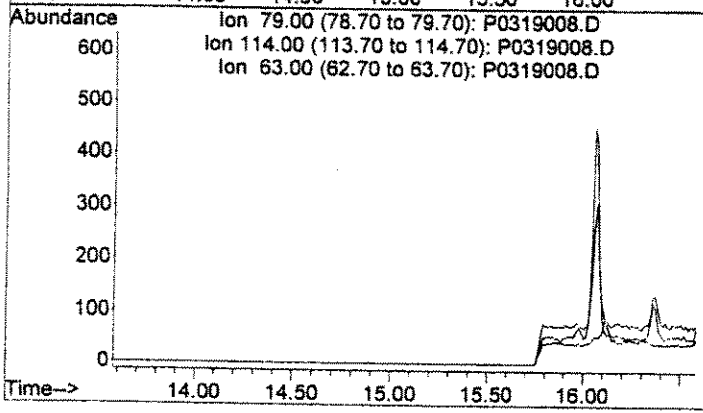




#5 *MT*  
 1,2,3-Trichloropropane-d5  
 Concen: 0.00 ug/L  
 Expected RT: 15.08 min

Lab File: P0319008.D  
 Acq: 19 Mar 2005 10:12 am

Sig	Exp Ratio
79	100
114	0.0
63	98.0



Grubbs Test for curve		MEAN	STDEV
Response factors		2.186429	0.456975
Grubbs value			
1.0ppb	2.0ppb	5.0ppb	10ppb
3.099	2.478	2.101	1.905
1.99698	0.63805	0.186944	0.615851
2.0ppb	5.0ppb	20ppb	50ppb
1.822	1.995	0.418904	0.797481
100ppb			1.905
			0.615851
5pts Grubbs values <	1.672		
6pts Grubbs values <	1.822		
7pts Grubbs values <	1.938		
8pts Grubbs values <	2.032		
9pts Grubbs values <	2.11		
10pts Grubbs values <	2.176		

outlier

MS 3/19/05

83/1/05

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

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

MS Integration Params: DIOXANE.P  
Quant Time: Mar 19 13:42 2005

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 NT	0.00	ug/L	-15.08
<b>System Monitoring Compounds</b>						
2) Dibromofluoromethane (SU1)	10.07	113	3531	0.11	ug/L	0.00
Spiked Amount	1.000	Range 80 - 120	Recovery	=	11.00%#	
<b>Target Compounds</b>						
4) 1,4-DIOXANE	12.43	88	615	1.50	ug/L	Qvalue 97

*DNV*

*Q*

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

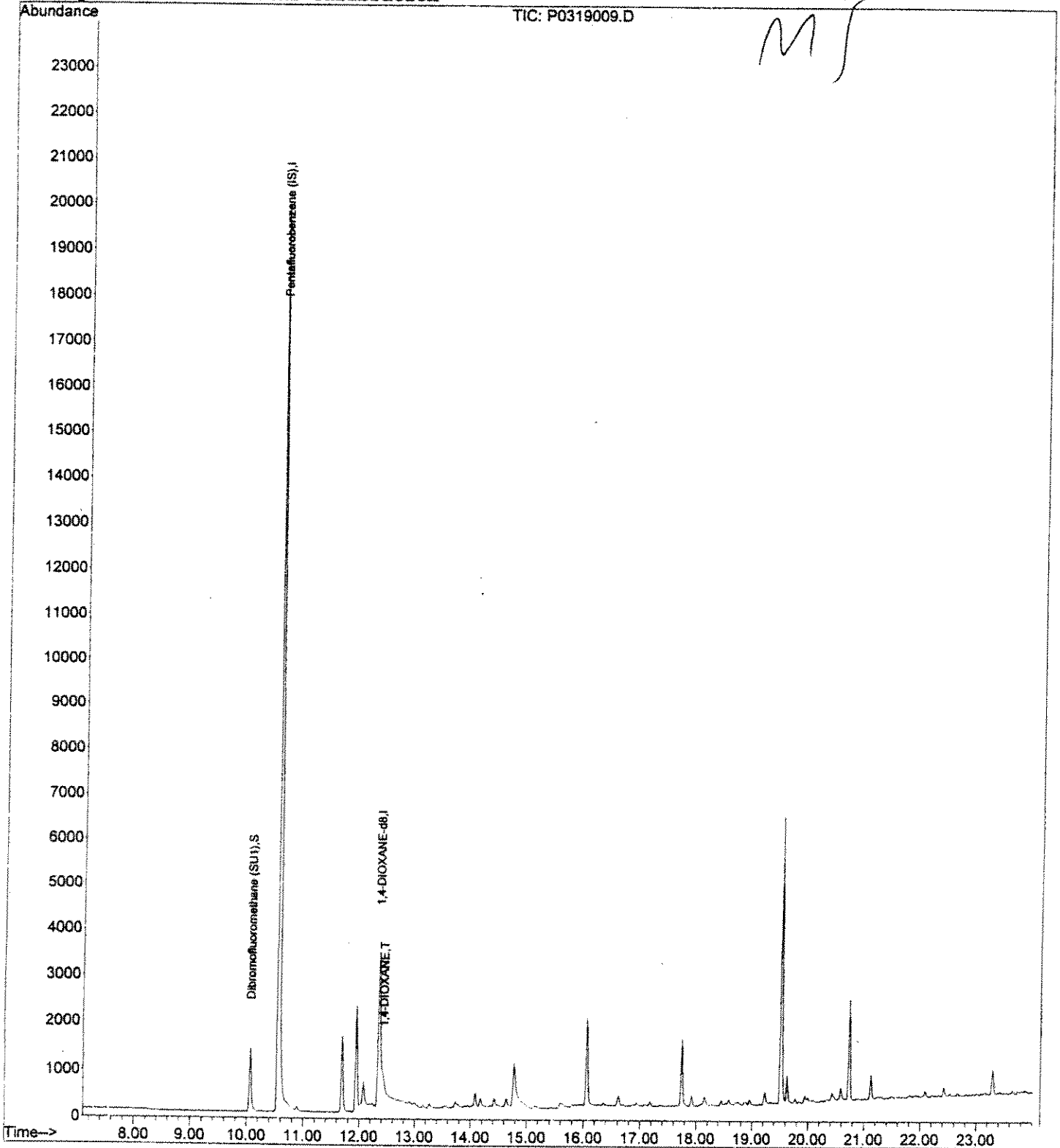
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\P0319009.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	0NT	0.00	ug/L	-15.08

System Monitoring Compounds

2) Dibromofluoromethane (SU1)	10.07	113	7585	0.21	ug/L	0.00
Spiked Amount	1.000	Range	80 - 120	Recovery	=	21.00%#

Target Compounds

4) 1,4-DIOXANE	12.43	88	1028	2.69	ug/L	Qvalue 94
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*8/21/05*

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

P0319010.D DX021605.M

Sat Mar 19 13:43:20 2005

GCMS1

Page

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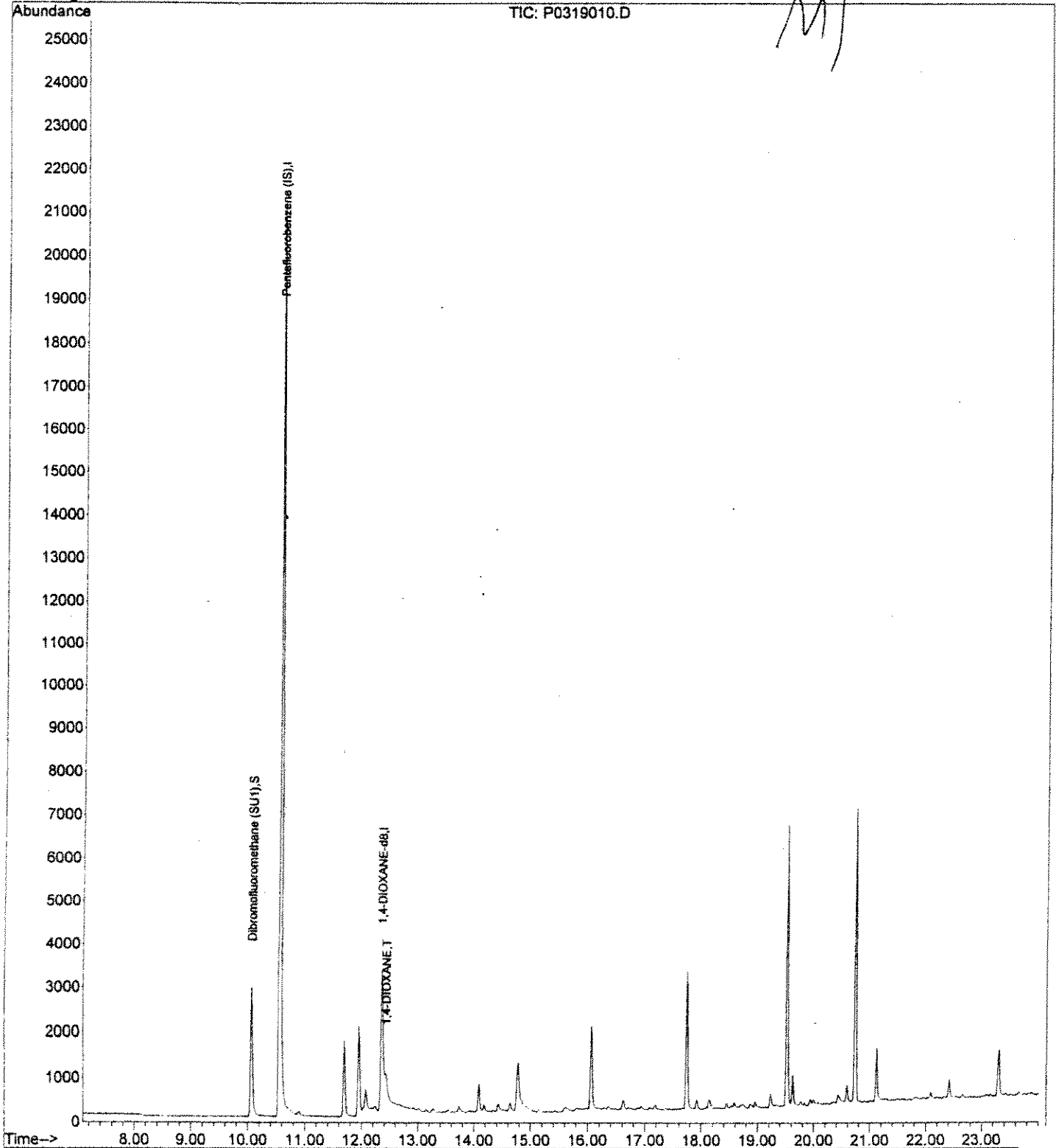
Quantitation Report

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	0 <sub>NT</sub>	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 (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\031905\PO319011.D  
 Acq On : 19 Mar 2005 11:59 am  
 Sample : 5.0 PPB CAL  
 Misc : 1X 10ML

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

MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 13:43 2005

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*



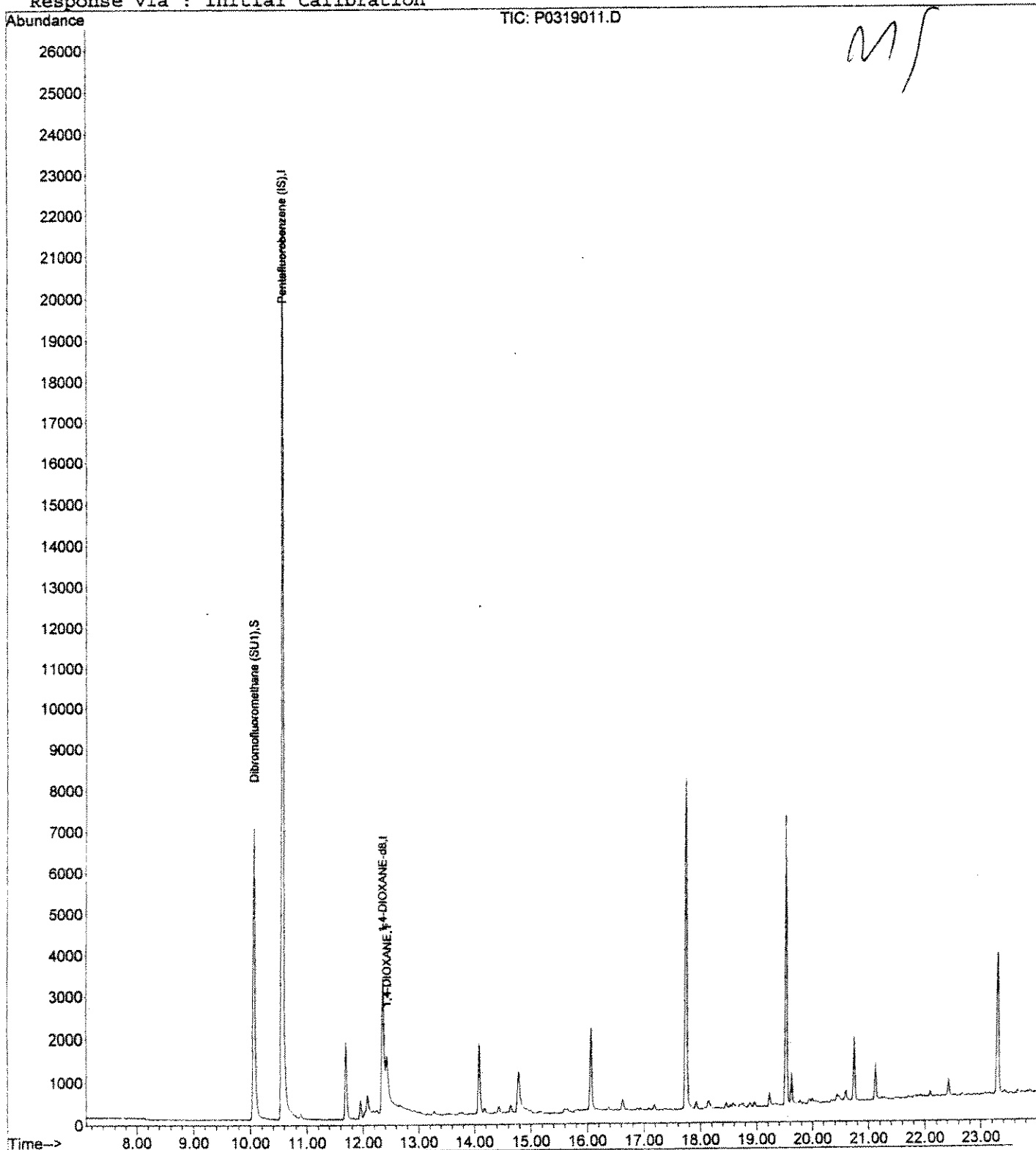
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\0319011.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\P0319012.D  
 Acq On : 19 Mar 2005 12:32 pm  
 Sample : 10.0 PPB CAL  
 Misc : 1X 10ML

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

MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 13:37 2005

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 13/19/05*

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	0.21	0.00	ug/L	-15.08
<b>System Monitoring Compounds</b>						
2) Dibromofluoromethane (SU1)	10.07	113	34373	0.95	ug/L	0.00
Spiked Amount	1.000	Range 80 - 120	Recovery	=	95.00%	
<b>Target Compounds</b>						
4) 1,4-DIOXANE	12.43	88	3835	11.74	ug/L	Qvalue 99

*3/21/05*

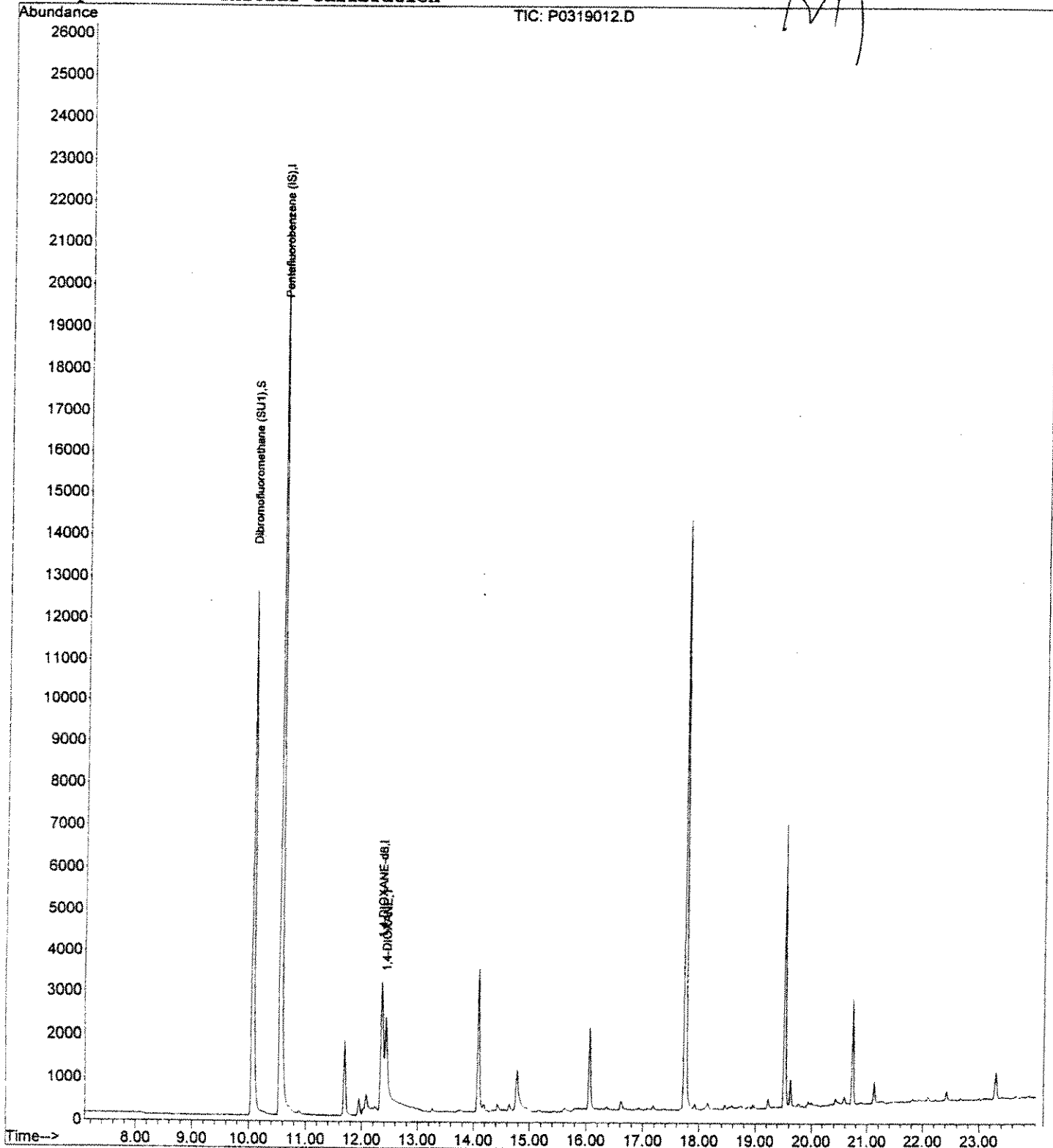
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\P0319012.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\P0319013.D  
 Acq On : 19 Mar 2005 1:05 pm  
 Sample : 20.0 PPB CAL  
 Misc : 1X 10ML

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

MS Integration Params: DIOXANE.P  
 Quant Time: Mar 19 13:37 2005

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

*MC 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 <sub>√T</sub>	0.00	ug/L	-15.08

System Monitoring Compounds

2) Dibromofluoromethane (SU1) 10.07 113 68573 1.86 ug/L 0.00  
 Spiked Amount 1.000 Range 80 - 120 Recovery = 186.00%#

Target Compounds

4) 1,4-DIOXANE 12.43 88 7646 25.14 ug/L Qvalue 97

*3/21/05*

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

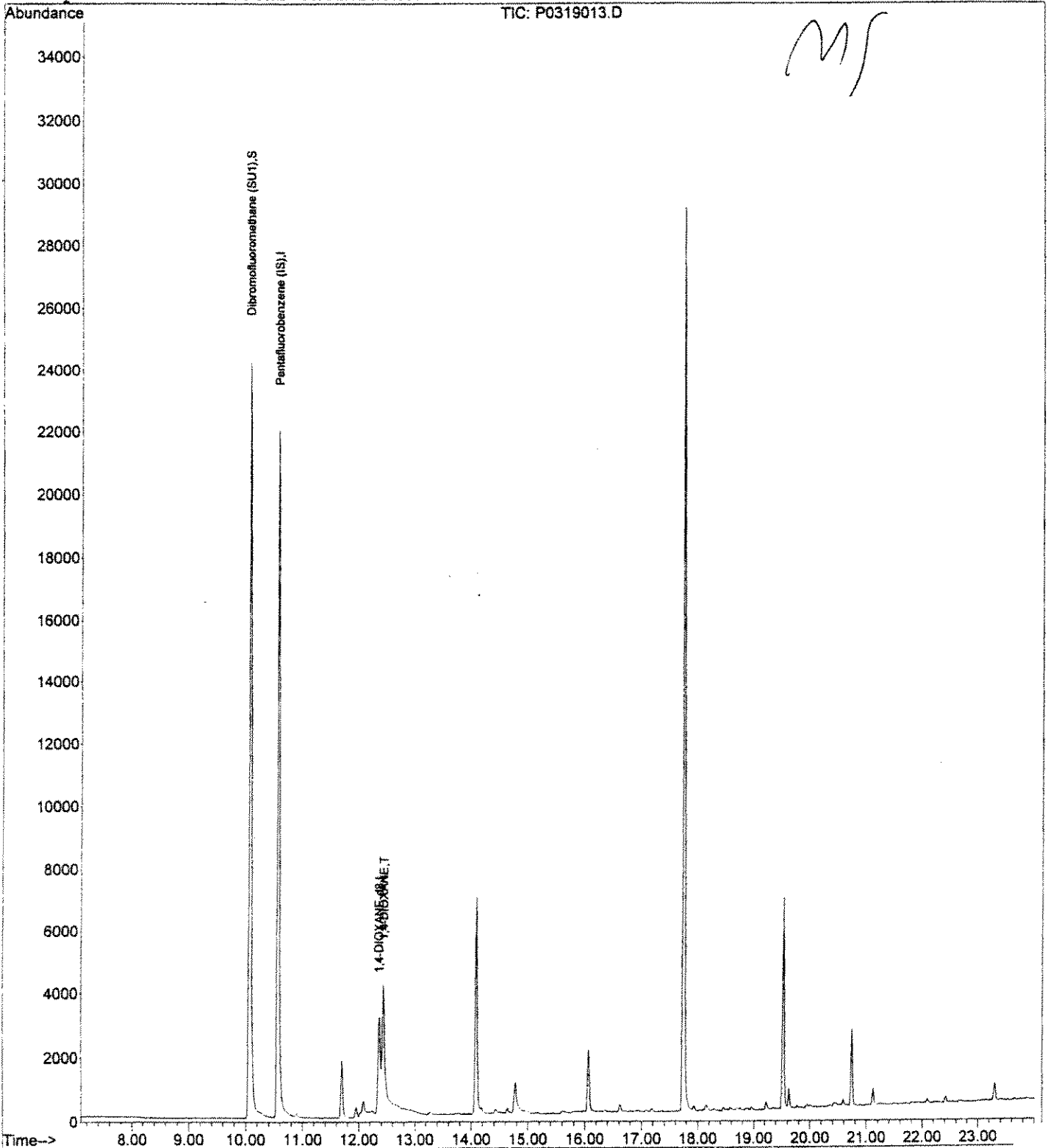
Quantitation Report

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

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\P0319014.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

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

4) 1,4-DIOXANE 12.43 88 18344 58.04 ug/L Qvalue 99

*JG/MS*

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

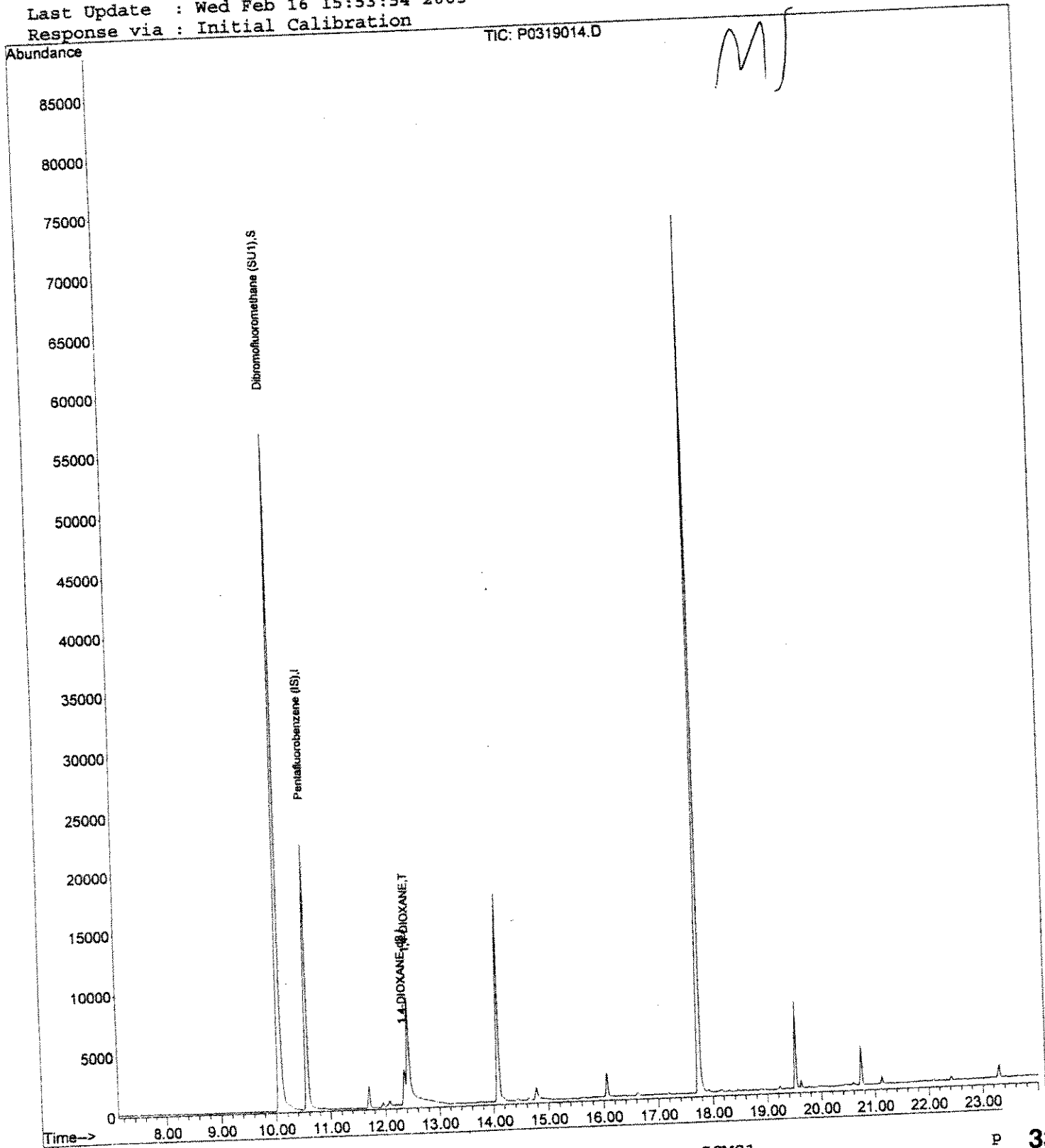
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\P0319014.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\PO319015.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

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 17/19/05*

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 <sup>√</sup>	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  
 4) 1,4-DIOXANE 12.43 88 44445 121.87 ug/L Qvalue 98

*3/17/05*



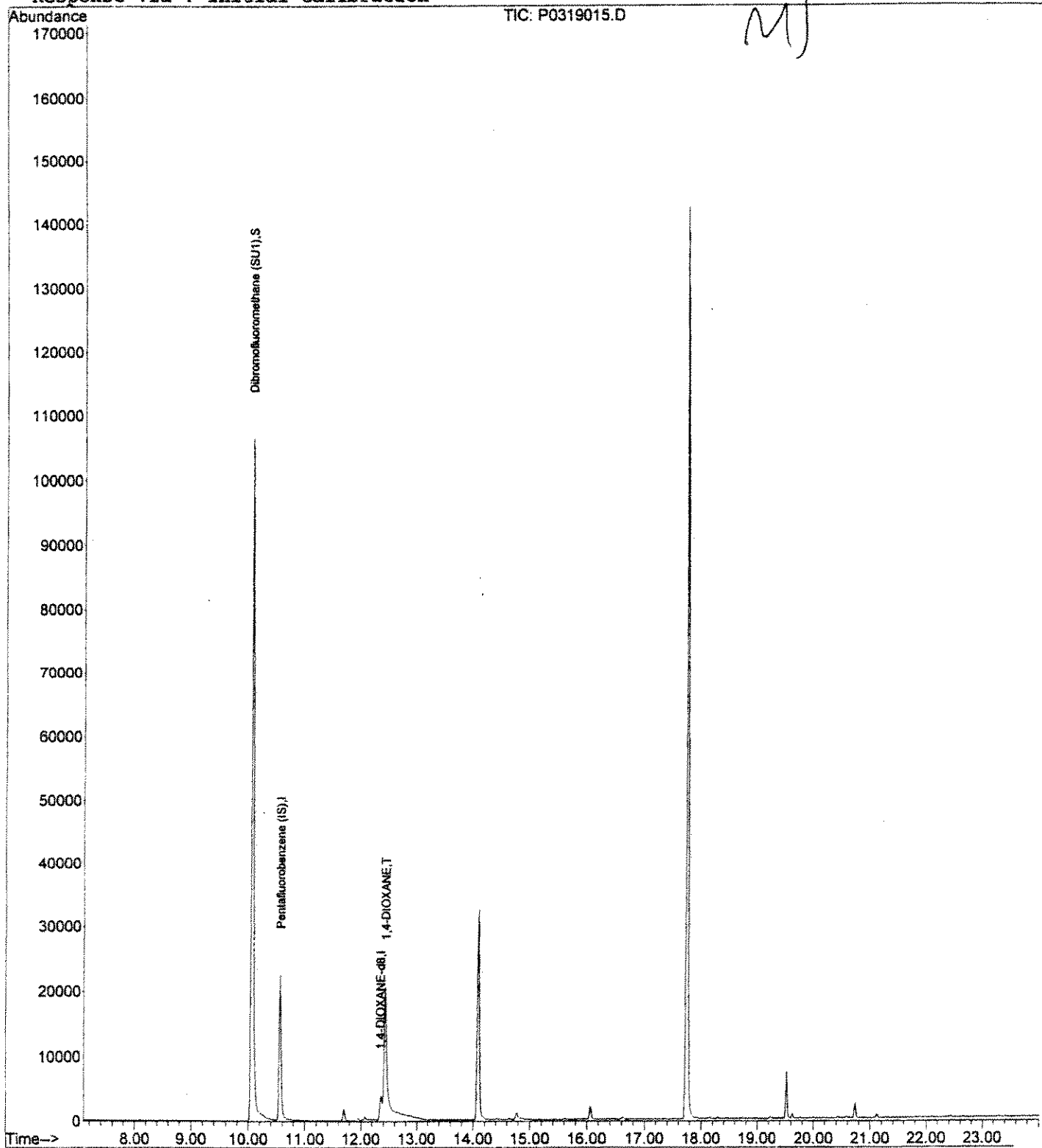
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\P0319015.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\P0319016.D  
 Acq On : 19 Mar 2005 2:44 pm  
 Sample : CLEAN OUT BLANK/TUNE  
 Misc : 1X 10ML

Vial: 16  
 Operator: JG/MS/CLS  
 Inst : GCMS1  
 Multiplr: 1.00

MS Integration Params: DIOXANE.P

Quant Time: Mar 21 7:48 2005

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 : W072903

3/21/05  
 Jly

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene (IS)	10.56	99	168438	1.00	ug/L	0.00
3) 1,4-DIOXANE-d8	12.36	64	64	25.00	ug/L	0.00
5) 1,2,3-Trichloropropane-d5	15.15	79	57	500.00	ug/L	0.07

System Monitoring Compounds

2) Dibromofluoromethane (SU1)	10.06	113	129670	1.00	ug/L	0.00
Spiked Amount	1.000	Range	80 - 120	Recovery	=	100.00%

Target Compounds

Qvalue

*DMU*

*Q*

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

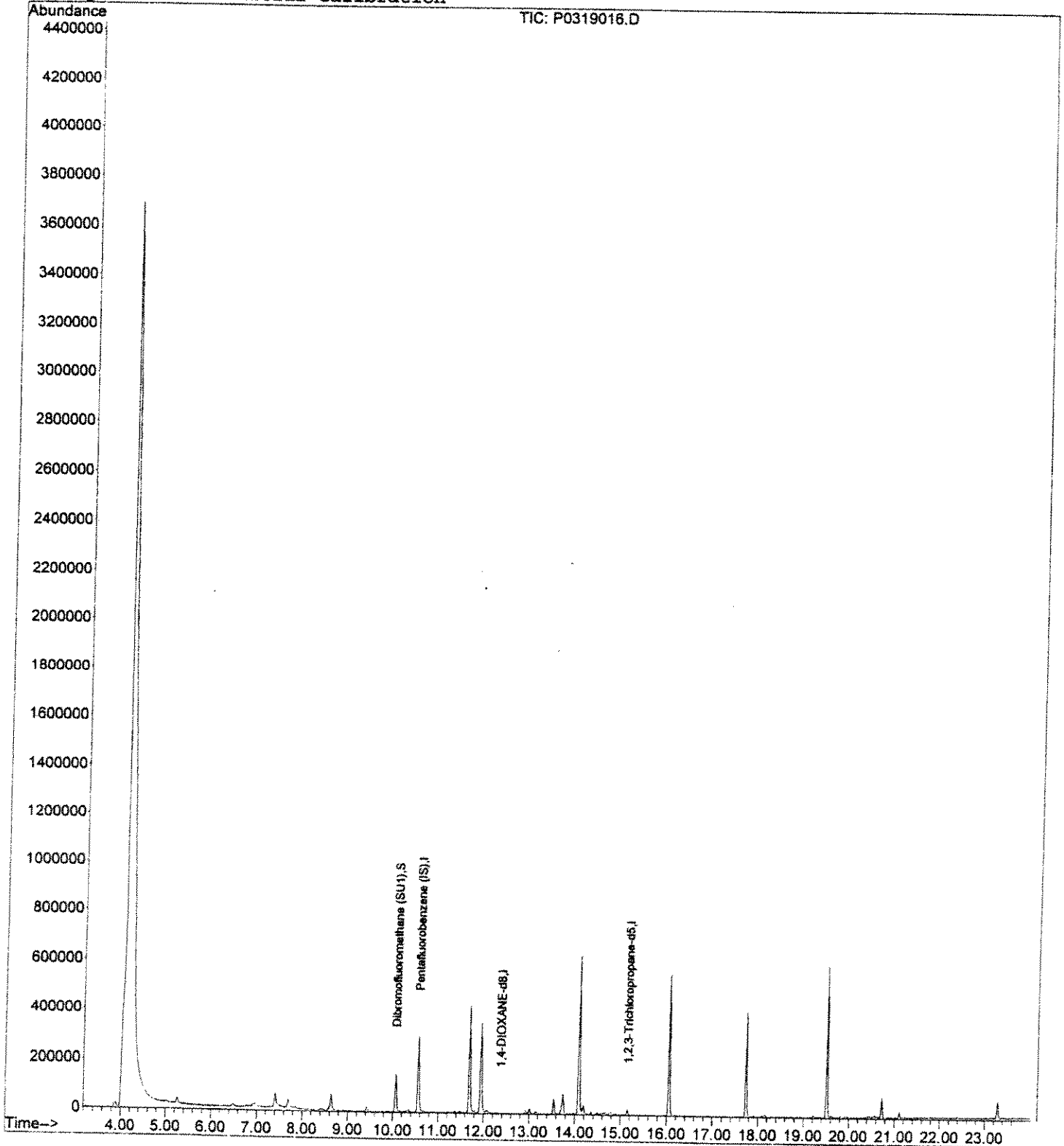
Quantitation Report

Data File : D:\HPCHEM\1\DATA\031905\0319016.D  
Acq On : 19 Mar 2005 2:44 pm  
Sample : CLEAN OUT BLANK/TUNE  
Misc : 1X 10ML  
MS Integration Params: DIOXANE.P  
Quant Time: Mar 21 7:48 2005

Vial: 16  
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\0319017.D Vial: 17  
 Acq On : 19 Mar 2005 3:21 pm Operator: JG/MS/CLS  
 Sample : BLANK Inst : GCMS1  
 Misc : 1X 10ML Multiplr: 1.00  
 MS Integration Params: DIOXANE.P  
 Quant Time: Mar 21 7:48 2005 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	41664	1.00	ug/L	0.00
3) 1,4-DIOXANE-d8	12.35	64	6641	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 34219 1.06 ug/L 0.00  
 Spiked Amount 1.000 Range 80 - 120 Recovery = 106.00%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) 1,4-DIOXANE	12.43	88	233	N.D.		
6) 1,2,3-Trichloropropane	0.00	75	0	N.D.		

*skul*