

APPENDIX A COPY OF BWSC108 TRANSMITTAL FORM



Massachusetts Department of Environmental Protection Bureau of Waste Site Cleanup

BWSC108

Release Tracking Number

4 _

_ 17786

COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H)

Pulsualitio 3 to Civil 40.0404 (Subpart D) and 40.0000 (Subpart D)
A. SITE LOCATION: 1. Site Name:Barge B120 Spill
2. Street Address:
3. City/Town: Buzzards Bay 4. ZIP Code: N/A
5. UTM Coordinates: a. UTM N: b. UTM E:
6. Check here if a Tier Classification Submittal has been provided to DEP for this disposal site. a. Tier IA b. Tier IB c. Tier IC d. Tier II 7. If applicable, provide the Permit Number:
B. THIS FORM IS BEING USED TO: (check all that apply)
1. Submit a Phase I Completion Statement, pursuant to 310 CMR 40.0484.
2. Submit a Revised Phase I Completion Statement, pursuant to 310 CMR 40.0484.
3. Submit a Phase II Scope of Work, pursuant to 310 CMR 40.0834.
4. Submit an interim Phase II Report . This report does not satisfy the response action deadline requirements in 310 CMR 40.0500.
5. Submit a final Phase II Report and Completion Statement, pursuant to 310 CMR 40,0836.
6. Submit a Revised Phase II Report and Completion Statement, pursuant to 310 CMR 40.0836.
7. Submit a Phase III Remedial Action Plan and Completion Statement, pursuant to 310 CMR 40.0862.
8. Submit a Revised Phase III Remedial Action Plan and Completion Statement, pursuant to 310 CMR 40.0862.
9. Submit a Phase IV Remedy Implementation Plan , pursuant to 310 CMR 40.0874.
10. Submit a Modified Phase IV Remedy Implementation Plan, pursuant to 310 CMR 40.0874.
11. Submit an As-Built Construction Report , pursuant to 310 CMR 40.0875.
12. Submit a Phase IV Status Report , pursuant to 310 CMR 40.0877.
13. Submit a Phase IV Completion Statement, pursuant to 310 CMR 40.0878 and 40.0879.
Specify the outcome of Phase IV activities: (check one)
a. Phase V Operation, Maintenance or Monitoring of the Comprehensive Remedial Action is necessary to achieve a Response Action Outcome.
b. The requirements of a Class A Response Action Outcome have been met. No additional Operation, Maintenance or Monitoring is necessary to ensure the integrity of the Response Action Outcome. A completed Response Action Outcome Statement and Report (BWSC104) will be submitted to DEP.
c. The requirements of a Class C Response Action Outcome have been met. No additional Operation, Maintenance or Monitoring is necessary to ensure the integrity of the Response Action Outcome. A completed Response Action Outcome Statement and Report (BWSC104) will be submitted to DEP.
d. The requirements of a Class C Response Action Outcome have been met. Further Operation, Maintenance or Monitoring of the remedial action is necessary to ensure that conditions are maintained and that further progress is made toward a Permanent Solution. A completed Response Action Outcome Statement and Report (BWSC104) will be submitted to DEP.
(All sections of this transmittal form must be filled out unless otherwise noted above)



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COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H) B. THIS FORM IS BEING USED TO (cont.): (check all that apply) 14. Submit a Revised Phase IV Completion Statement, pursuant to 310 CMR 40.0878 and 40.0879. 15. Submit a Phase V Status Report, pursuant to 310 CMR 40.0892. 16. Submit a Remedial Monitoring Report. (This report can only be submitted through eDEP.) ii. Interim Report a. Type of Report: (check one) i. Initial Report iii. Final Report b. Frequency of Submittal: (check all that apply) i. A Remedial Monitoring Report(s) submitted monthly to address an Imminent Hazard. ii. A Remedial Monitoring Report(s) submitted monthly to address a Condition of Substantial Release Migration. iii. A Remedial Monitoring Report(s) submitted concurrent with a Status Report. iii. Class C RAO i. Phase V ii. Remedy Operation Status c. Status of Site: (check one) d. Number of Remedial Systems and/or Monitoring Programs: A separate BWSC108A, CRA Remedial Monitoring Report, must be filled out for each Remedial System and/or Monitoring Program addressed by this transmittal form. 17. Submit a Remedy Operation Status, pursuant to 310 CMR 40.0893. 18. Submit a Status Report to maintain a Remedy Operation Status, pursuant to 310 CMR 40,0893(2). 19. Submit a Modification of a Remedy Operation Status, pursuant to 310 CMR 40.0893(5), 20. Submit a Termination of a Remedy Operation Status, pursuant to 310 CMR 40.0893(6). 21. Submit a Phase V Completion Statement, pursuant to 310 CMR 40,0894. Specify the outcome of Phase V activities: (check one) a. The requirements of a Class A Response Action Outcome have been met. No additional Operation, Maintenance or Monitoring is necessary to ensure the integrity of the Response Action Outcome. A completed Response Action Outcome Statement (BWSC104) will be submitted to DEP. b. The requirements of a Class C Response Action Outcome have been met. No additional Operation, Maintenance or Monitoring is necessary to ensure the integrity of the Response Action Outcome. A completed Response Action Outcome Statement and Report (BWSC104) will be submitted to DEP. c. The requirements of a Class C Response Action Outcome have been met. Further Operation, Maintenance or Monitoring of the remedial action is necessary to ensure that conditions are maintained and/or that further progress is made toward a Permanent Solution. A completed Response Action Outcome Statement and Report (BWSC104) will be submitted to DEP. 22. Submit a Revised Phase V Completion Statement, pursuant to 310 CMR 40.0894. 23. Submit a Post-Class C Response Action Outcome Status Report, pursuant to 310 CMR 40,0898.

(All sections of this transmittal form must be filled out unless otherwise noted above)



Massachusetts Department of Environmental Protection Bureau of Waste Site Cleanup

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COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

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Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H)

C. LSP SIGNATURE AND STAMP:

I attest under the pains and penalties of perjury that I have personally examined and am familiar with this transmittal form, including any and all documents accompanying this submittal. In my professional opinion and judgment based upon application of (i) the standard of care in 309 CMR 4.02(1), (ii) the applicable provisions of 309 CMR 4.02(2) and (3), and 309 CMR 4.03(2), and (iii) the provisions of 309 CMR 4.03(3), to the best of my knowledge, information and belief.

- > if Section B indicates that a **Phase I, Phase II, Phase III, Phase IV or Phase V Completion Statement** is being submitted, the response action(s) that is (are) the subject of this submittal (i) has (have) been developed and implemented in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal;
- > if Section B indicates that a **Phase II Scope of Work** or a **Phase IV Remedy Implementation Plan** is being submitted, the response action(s) that is (are) the subject of this submittal (i) has (have) been developed in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal:
- > if Section B indicates that an As-Built Construction Report, a Remedy Operation Status, a Phase IV, Phase V or Post-Class C RAO Status Report, a Status Report to Maintain a Remedy Operation Status and/or a Remedial Monitoring Report is being submitted, the response action(s) that is (are) the subject of this submittal (i) is (are) being implemented in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal.

I am aware that significant penalties may result, including, but not limited to, possible fines and imprisonment, if I submit information which I know to be false, inaccurate or materially incomplete.

1. LSP#: 5463		
2. First Name: Richard	3. Last Name: Wozmak	
4. Telephone: (603) 421-2777 5. Ext.:	6. FAX: (603) 421-9880	
7. Signature:		j
8. Date: <u>07/23/2007</u> (mm/dd/yyyy)	9. LSP Stamp: WOZMAK No. 5463 REGISTER STEEPHOTE	

Revised: 2/15/2005 Page 3 of 5

Massachusetts Department of Environmental Protection Bureau of Waste Site Cleanup

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COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

Release Tracking Number 17786

Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H)			
D. PERSON UNDERTAKING RESPONSE ACTIONS:			
1. Check all that apply: a. change in contact name b. change of address c. change in the person undertaking response actions			
2. Name of Organization: Bouchard Transportation Co., Inc.			
3. Contact First Name: W. Lawrence 4. Last Name: Lopez			
5. Street: 58 South Service Road, Suite 150 6. Title: Risk Manager			
7. City/Town: Melville 8. State: NY 9. ZIP Code: 11747			
10. Telephone: <u>(516)</u> 681-4900 <u>11. Ext.:</u> 12. FAX:			
E. RELATIONSHIP TO SITE OF PERSON UNDERTAKING RESPONSE ACTIONS:			
1. RP or PRP a. Owner b. Operator c. Generator d. Transporter			
e. Other RP or PRP Specify:			
2. Fiduciary, Secured Lender or Municipality with Exempt Status (as defined by M.G.L. c. 21E, s. 2)			
3. Agency or Public Utility on a Right of Way (as defined by M.G.L. c. 21E, s. 5(j))			
4. Any Other Person Undertaking Response Actions Specify Relationship:			
F. REQUIRED ATTACHMENT AND SUBMITTALS:			
Check here if the Response Action(s) on which this opinion is based, if any, are (were) subject to any order(s), permit(s) and/or approval(s) issued by DEP or EPA. If the box is checked, you MUST attach a statement identifying the applicable provisions thereof.			
2. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of the submittal of any Phase Reports to DEP.			
3. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of the availability of a Phase III Remedial Action Plan.			
4. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of the availability of a Phase IV Remedy Implementation Plan.			
5. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of any field work involving the implementation of a Phase IV Remedial Action.			
6. If submitting a Modification of a Remedy Operation Status, check here to certify that a statement detailing the compliance history, as per 310 CMR 40.0893(5), for the person making this submittal is attached.			
7. If submitting a Modification of a Remedy Operation Status, check here to certify that written consent of the person who submitted the Remedy Operation Status submittal, as per 310 CMR 40.0893(5), is attached.			
8. Check here if any non-updatable information provided on this form is incorrect, e.g. Site Name. Send corrections to the DEP Regional Office.			
L DEP Regional Office.			
9. Check here to certify that the LSP Opinion containing the material facts, data, and other information is attached.			



Massachusetts Department of Environmental Protection Bureau of Waste Site Cleanup

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COMPREHENSIVE RESPONSE ACTION TRANSMITTAL FORM & PHASE I COMPLETION STATEMENT

Pursuant to 310 CMR 40.0484 (Subpart D) and 40.0800 (Subpart H)

G. CERTIFICATION OF PERSON UNDERTAKING RESPONSE ACTIONS:			
1. I. Richard Wozmak attest under the pains and	nonalting of pariury (i) that they appropriate		
1. I, Richard Wozmak , attest under the pains and penalties of perjury (i) that I have personally examined and am familiar with the information contained in this submittal, including any and all documents accompanying this			
transmittal form, (ii) that, based on my inquiry of those individuals immediately res	sponsible for obtaining the information, the		
material information contained in this submittal is, to the best of my knowledge and	d belief, true, accurate and complete, and (iii)		
that I am fully authorized to make this attestation on behalf of the entity legally respentity on whose behalf this submittal is made am/is aware that there are significa	onsible for this submittal. I/the person or		
possible times and imprisonment, for willfully submitting false, inaccurate, or incomposition of the same significant possible times and imprisonment, for willfully submitting false, inaccurate, or incomposition of the same significant possible times and imprisonment.	nt penalties, including, but not limited to, molete information		
2. By: Signature	3. Title: Licensed Oile i Tolessional		
Signature			
4. For: Agent for Bouchard Transportation Co., Inc. (Name of person or entity recorded in Section D)	5. Date: 07/23/2007		
(Name of person or entity recorded in Section D)	* (mm/dd/yyyy)		
6. Check here if the address of the person providing certification is different fr	om address recorded in Section D.		
7. Street:			
8. City/Town: 9. State:	10. ZIP Code:		
11. Telephone: 12. Ext.: 13. FA	v.		
12. Ext. 13. PA	^		
YOU ARE SUBJECT TO AN ANNUAL COMPLIANCE ASSURANCE			
BILLABLE YEAR FOR THIS DISPOSAL SITE. YOU MUST LEGIBL			
SECTIONS OF THIS FORM OR DEP MAY RETURN THE DOCUME SUBMIT AN INCOMPLETE FORM, YOU MAY BE PENALIZED FOR M			
SUBMIT AN INCOMPLETE FORM, TOO MAT BE PENALIZED FOR W	ISSING A REQUIRED DEADLINE.		
Date Stamp (DEP USE ONLY:)			
This claim, (SET COL CIVET.)			

Supplement to BWSC108, Section F

Barge B120 Release Buzzards Bay, Massachusetts 4-17786

Section F – Required Attachments and Submittals

1. Check here if the Response Action(s) on which this opinion is based, if any, are (were) subject to any order(s), permit(s) and/or approval(s) issued by DEP or EPA. If the box is checked, you MUST attach a statement identifying the applicable provisions thereof.

Massachusetts Department of Environmental Protection Orders, Permits, and/or Approvals:

- September 8, 2003 Request for IRA with Interim Deadlines;
- July 27, 2004 Decision to Grant Permit;
- January 18, 2006 Phase II Scope of Work Conditional Approval/Interim Deadline;
- June 27, 2006 Phase II SOW Addendum Approval; and
- October 19, 2006 Phase II Comprehensive Site Assessment Report and Phase III Remedial Action Plan Approval.



APPENDIX B LABORATORY ANALYTICAL REPORT

TDI - BROOKS INTERNATIONAL, INC. B&B Laboratories, Inc. College Station, TX

GeoInsight, Inc.
Buzzards Bay Spill Project

Determination of: Total Petroleum Hydrocarbons and Polycyclic Aromatic Hydrocarbons in an Oiled Rock Sample

(QC Batch ENV 1635)

May 8, 2007

Technical Report 07-1857

Geolnsight, Inc. Buzzards Bay Spill Project Table of Contents B&B Laboratories 08-May-2007

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Technical Report 07-1857 Geolnsight, Inc. Buzzards Bay Spill Project Oiled Rock (Product) Sample

May 8, 2007

Introduction

B&B Laboratories received one (1) ice chest that contained one (1) bag that contained an oiled rock sample that was sent on April 11, 2007 and arrived on April 12, 2007 at B&B Laboratories in College Station, Texas sealed and in good condition. The internal temperature of the cooler was 14.8°C. The oiled rock sample was collected in support of the Buzzards Bay Spill Project (GeoInsight Project 3871-002). The oiled rock sample was stored in an access-controlled refrigerator (4.0°C) until processing. The oiled rock sample was analyzed for Total Petroleum Hydrocarbons (TPH) by GC/FID and Polycyclic Aromatic Hydrocarbons (PAHs) by GC/MS-SIM.

The results for TPH, PAH and hopanes are included in this report.

Analytical Methods

The analytical methods employed for TPH and PAH are listed in Table 1.

Table 1. Standard Operating Procedures for each analytical test.

Matrix	Extraction	ТРН	PAH
Product	SW-846 3580A	B&B 1013	B&B 1006

Data Reporting

The reporting units for each analyte are listed in Table 2. The method detection limits (MDL) for each analyte are listed in Table 3. Analytes that are detected below the method detection limit are qualified as "J". Analytes that are detected in the procedural blanks greater than 3X MDL are qualified with a "B". Analytical interference's that are detected in the sample are qualified with an "I". Analytes not detected in the samples are qualified with a "U". RPD for analytes in duplicate samples that are <2X MDL are qualified with a "X". Spiked levels of analytes in matrix spikes that are <50% of the native levels are considered invalid spikes and are qualified with a "Y". Any QC result reported to be outside the corresponding QC criteria is discussed in the QA/QC variance section of this report.

Table 2. Analytical Reporting Units.

Matrix	TPH/ALI	PAH
Product	ug/mg	ng/mg

Table 3. Reporting Limits.

TPH	RLs
Sample size	1ml final extract volume
Unit of measure	μg/mg
Total Petroleum Hydrocarbons Total Resolved Hydrocarbons Unresolved Complex Mixture Extractable Organic Matter	1.4 1.4 1.4 100

Table 3 (Continued). Reporting Limits.

PAH	RL
Unit of measure	ng/ug
Naphthalene	10
C1-Naphthalenes	10
C2-Naphthalenes	10
C3-Naphthalenes	10
C4-Naphthalenes	10
Benzothiophene	10
C1-Benzothiophenes	10
C2-Benzothiophenes	10
C3-Benzothiophenes	10
Biphenyl	10
Acenaphthylene	10
Acenaphthene	10
Dibenzofuran	10
Fluorene	10
C1-Fluorenes	10
C2-Fluorenes	10
C3-Fluorenes	10
Carbazole	10
Anthracene	10
Phenanthrene	10
C1-Phenanthrenes/Anthracenes	10
C2-Phenanthrenes/Anthracenes	10
C3-Phenanthrenes/Anthracenes	10
C4-Phenanthrenes/Anthracenes	10
Dibenzothiophene	10
C1-Dibenzothiophenes	10
C2-Dibenzothiophenes	10
C3-Dibenzothiophenes	10
Fluoranthene	10
Pyrene	10
C1-Fluoranthenes/Pyrenes	10
C2-Fluoranthenes/Pyrenes	10
C3-Fluoranthenes/Pyrenes	10
Benz(a)anthracene	10

PAH (Continued)	RL
Unit of measure	ng/ug
Chrysene	10
C1-Chrysenes	10
C2-Chrysenes	10
C3-Chrysenes	10
C4-Chrysenes	10
Benzo(b)fluoranthene	10
Benzo(k)fluoranthene	10
Benzo(e)pyrene	10
Benzo(a)pyrene	10
Perylene	10
Indeno(1,2,3-c,d)pyrene	10
Dibenzo(a,h)anthracene	10
Benzo(g,h,i)perylene	10
Individual Alkyl Isomers and Hopanes	
2-Methylnaphthalene	10
1-Methylnaphthalene	10
2,6-Dimethylnaphthalene	10
1,6,7-Trimethylnaphthalene	10
1-Methylphenanthrene	10
C29-Hopane	10
18a-Oleanane	10
C30-Hopane	10

Quality Assurance/Quality Control

Oiled Rock (mg of extract)

TPH

The quality assurance/quality control procedure for this program included the analyses of a laboratory control sample (LCS) that was analyzed with each data set. The LCS is a diesel sample that is analyzed with each TPH/ALI run and for which controls are established based on performance. The QC criterion for the LCS is between 85 - 115% of the laboratory determined mean.

Surrogate solutions equivalent to 5-10X the MDL are prepared for various hydrocarbon analyses. The appropriate surrogate solution is added to every sample including quality control samples. The data are corrected based on surrogate recovery up to 100%. The QC criteria for surrogate recoveries are between 40-120%.

PAH

The quality assurance/quality control procedure for this program included a standard reference oil (NIST 1582) and a laboratory control solution that were analyzed with this data set. A SRM is a material for which a mean and confidence interval are certified for specific analytes. SRMs are selected based on matrix similarities as well as type and level of certified analytes. All SRMs are traceable to NIST. SRMs are used to verify analytical accuracy. All QC samples are subject to the identical preparation and analysis steps as samples. The QC criterion for the reference oil SRM and the laboratory control material is \pm 15% the laboratory derived mean.

Surrogate solutions equivalent to 5-10X the MDL are prepared for various hydrocarbon analyses. The appropriate surrogate solution is added to every sample including quality control samples. The data are corrected based on surrogate recovery up to 100%. The QC criteria for surrogate recoveries are between 40-120%, except d12-perylene.

Quality Assurance/Quality Control Variances

Oiled Rock (mg of extract)

Total Petroleum Hydrocarbons (TPH)

Surrogate Recoveries

Observation

No variances were observed.

Laboratory Control Standard

Observation

No variances were observed.

PAH

Surrogate Recoveries

Observation

• No variances were observed.

Standard Reference Materials/Laboratory Control Materials

Observation

No variances were observed.

We appreciate the opportunity to serve your analytical needs and please do not hesitate to contact us should you have any questions.

Thomas J. McDonald Project Manager

Project Quality Officer

Donell Frank

Sample/Analyses Description

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Client Project #3871-002	Comments B&B SDG Entrix Project #	Oiled rock 07041201 3871-002
	Matrix C	Other
roject ry	Analysis	РАН, ТРН
Geolnsight, Inc. Buzzards Bay Spill Project Sample Inventory	Receive Date	04/12/07
<u>α</u>	Collection Date	03/27/07
	Client Identification	WIF-02-32707
B&B Laboratories Project J03318 Report 07-1857	Laboratory File Number Client Identification Collection Date Receive Date	ETX7207

Product Samples

Total Petroleum Hydrocarbons/ Extractable Organic Matter Concentrations

Geolnsight, Inc. Buzzards Bay Oil Spill Project Total Petroleum Hydrocarbon Data Client Submitted Samples

Sample Name	ETX7207.D	
Client Name	WIF-02-32707	
Matrix	Oiled Rock	
Collection Date	03/27/07	
Received Date	04/12/07	
Extraction Date	04/12/07	
Extraction Batch	ENV 1635	
Date Acquired	04/13/07	
Method	ALI_COMP.M	
Sample Weight (mg)	17.7	
Dilution	NA	
Target Compounds	Su Corrected Conc (µg/mg)	**********
	Done (Hgring)	
Total Petroleum Hydrocarbons		266
Total Resolved Hydrocarbons		69
Unresolved Complex Mixture		197
Surrogate (Su)	Su Recovery (%)	
n-dodecane-d26	102	
n-eicosane-d42	92	
n-triacontane-d62	91	

Qualifiers (Q): J=Below the MDL, U=Not detected, B=In procedural blank > 3x MDL, I=Interference, D=Diluted value, NA=Not applicable, '=Outside QA limits, refer to narrative If n-eicosane-d42 (surrogate) recovery is above 100%, TPH and aliphatic values are surrogate corrected to 100%.

Geolnsight, Inc. Buzzards Bay Oil Spill Project Total Petroleum Hydrocarbon Data Lab Control Standard Report

Sample Name	GC10846B.D					
Client Name	AL-WKDIESEL-1000-005					
Matrix	Solution					
Collection Date	NA					
Received Date	NA					
Extraction Date	NA					
Extraction Batch	EOM 1635					
Date Acquired	04/13/07					
Method	ALI_COMP.M					
Sample Volume (mL)	1.0					
Dilution	NA					
Target Compounds	Su Corrected Conc (µg/mL)	Q	RPD (%)	B&B Average Conc (μg/mL)	-15% Conc (µg/mL)	+15% Conc (µg/mL)
Total Petroleum Hydrocarbons	101		0.5	100	85	115
Surrogate (Su)	Su Recovery (%)					
n-dodecane-d26	94					
n-eicosane-d42	99					
n-triacontane-d62	9 8					

Qualifiers (Q): J=Below the MDL, U=Not detected, B=In procedural blank > 3x MDL, I=Interference, D=Diluted value, NA=Not applicable, *=Outside QA limits, refer to narrative If n-eicosane-d42 (surrogate) recovery is above 100%, TPH values are surrogate corrected to 100%.

Total Petroleum Hydrocarbons Chromatograms

0.00 2.00 4.00 6.00 8.00 10.00 12.00 14.00 16.00 18.00 20.00 22.00 24.00 26.00 30.00 32.00 34.00 36.00 38.00 40.00 42.00 44.00 46.00 48.00 50.00 52.00 View Mode: Integration ETX7207.D\FID2B 160000 150000 140000 130000 120000 110000 100000 00006 80000 70000 00009 50000 40000 30000 20000 10000 0 Response Time

using AcqMethod ALI_COMP.M

U:\2\DATA\GC10846\ETX7207.D TJM

19:40

Acquired: 13 Apr 2007 Instrument: GC#1 Sample Name: WIF-02-32707

Operator

File

56

Misc Info : Vial Number:

Polycyclic Aromatic Hydrocarbon Concentration

Geoinsight, Inc. Buzzards Bay Oil Spill Project Polycyclic Aromatic Hydrocarbon Data Client Submitted Samples

Sample Name Client Name Matrix Collection Date Received Date Extraction Date Extraction Batch	ETX7207.D WIF-02-32707 Oiled Rock 03/27/07 04/12/07 04/12/07 ENV 1835	
Date Acquired Method Sample Weight (mg)	04/13/07 PAH-2002 17.7	
Dilution	NA NA	
Target Compounds	Su Corrected Q Conc. (ng/mg)	
Naphthalene	0.1 J	
C1-Naphthalenes	U. 8.0	
C2-Naphthalenes C3-Naphthalenes	335 1330	
C4-Naphthalenes	1130	
Benzothiophene	<10 U	
C1-Benzothiophenes	1.6 J	
C2-Benzothiophenes	21.6	
C3-Benzothiophenes Biphenyl	75.8 0.1 J	
Acenaphthylene	<10 U	
Acenaphthene	30.9	
Dibenzofuran	18.2	
Fluorene C1-Fluorenes	81.3 390	
C2-Fluorenes	1060	
C3-Fluorenes	1360	
Carbazole	<10 U	
Anthracene Phenanthrene	57.2 317	
C1-Phenanthrene/Anthracenes	2330	
C2-Phenanthrene/Anthracenes	4970	
C3-Phenanthrene/Anthracenes	5240	
C4-Phenanthrene/Anthracenes Dibenzothiophene	2710 75.5	
C1-Dibenzothiophenes	345	
C2-Dibenzothiophenes	737	
C3-Dibenzothiophenes	853	
Fluoranthene Pyrene	88.9 558	
C1-Fluoranthenes/Pyrenes	2490	
C2-Fluoranthenes/Pyrenes	3340	
C3-Fluoranthenes/Pyrenes	2710	
Naphthobenzothiophene C1-Naphthobenzothiophenes	329 799	
C2-Naphthobenzothiophenes	979	
C3-Naphthobenzothiophenes	455	
Benz(a)anthracene	318	
Chrysene C1-Chrysenes	772 2950	
C2-Chrysenes	3230	
C3-Chrysenes	1360	
C4-Chrysenes Benzo(b)fluoranthene	42.8	
Benzo(k)fluoranthene	138 18.4	
Benzo(e)pyrene	148	
Вепzo(а)ругепе	278	
Perylene Indeno(1,2,3-c,d)pyrene	71.2 23.5	
Dibenzo(a,h)anthracene	36.9	
Benzo(g.h.i)perylene	45.8	
Total PAHs	44652	
Individual Alkyl Isomers and Hop	panes	
2-Methylnaphthalene	0.3 J	
1-Methylnaphthalene	1.0 J 63.3	
2,6-Dimethylnaphthalene 1,6,7-Trimethylnaphthalene	206	
1-Methylphenanthrene	326	
C29-Hopane	71.2	
18a-Oleanane C30-Hopane	13.4 85.0	
Surrogate (Su)	Su Recovery (%)	
Nanhthalana da	84	
Naphthalene-d8 Acenaphthene-d10	84 90	
Phenanthrene-d10	91	

Qualifiers (Q): J=Below the MDL, U=Not detected, B=In procedural blank > 3x MDL, I=Interference, D=Diluted value, NA=Not Applicable, *=Outside QA limits, refer to narral

Geolnsight, Inc. Buzzards Bay Oil Spill Project Polycyclic Aromatic Hydrocarbon Data Standard Reference Material Report

Date Acquired DA1307 Method of DA1307 Metho	Sample Name Client Name Matrix Collection Date Received Date	MS30377B.D SRM 1582 Petroleum NA NA					
Sample Weight (g) 1.7	Extraction Batch Date Acquired	ENV 1635 04/13/07					
Conc. (upg)	Sample Weight (g)						
CI-Naghinslenes	Target Compounds			Certified Conc		Conc.	Conc.
CI-Naghhalenes 579 7.2 522 529 7.5 C2-Naghhalenes 1070 10.5 1189 1011 1080 C3-Naghhalenes 1020 1.8 724 641 1087 C3-Naghhalenes 1020 1.8 724 641 1087 C3-Naghhalenes 17.5 1.8 724 641 887 C3-Naghhalenes 17.5 1.8 724 641 887 C3-Benzofilophene 8.3 3 724 C3-Benzofilophenes 17.5 72.5 72.5 C3-Benzofilophenes 17.5 72.5 72.5 C3-Benzofilophenes 17.5 72.5 72.5 C3-Benzofilophenes 17.5 72.5 72.5 72.5 C3-Benzofilophenes 17.5 72.5 72.5 72.5 C3-Benzofilophenes 17.5 72.5 72.5 72.5 72.5 C3-Benzofilophenes 17.5 72.5 72.5 72.5 72.5 C3-Benzofilophenes 17.5 72.5 72.5 72.5 72.5 72.5 72.5 C3-Benzofilophene 17.5 72.	Naphthalene		2.0				
C3-Naghthalenes	C1-Naphthalenes						
Beausthophenes	C3-Naphthalenes	1020	1.7		1037	881	1193
C1-Benzothlophenes	C4-Naphthalenes				754	641	867
C3-Bernothophenes 147 147 34.0 1.5 34.5 29.3 39.7	C1-Benzothiophenes	17.5					
Bignery							
Accessible 1.7 13.8 16.9 16.1 21.7	Siphenyl	34.0			34.5	29.3	39.7
Disentalitarian 12.4	Acenaphthylene				18.9	16 1	21 7
Fluorene	Dibenzofuran	12.4					
C2-Fluorenes	Fluorene C1-Fluorenes						
C3-Fluorenes	C1-Fluorenes C2-Fluorenes	271	5.7		256	218	294
Anthracenee 3.8 J Phenanthrene	C3-Fluorenes				242	206	278
Phenanthrene 100 9.3 100 ± 7.0 110 93.3 126 C2-Phenanthrene/Anthracenes 516 5.1 326 277 375 C2-Phenanthrene/Anthracenes 516 5.1 5.3 402 624 C2-Phenanthrene/Anthracenes 280 1.5 5.1 275 35.5 C2-Phenanthrene/Anthracenes 280 1.5 32.9 ± 1.7 C2-Phenanthrene/Anthracenes 252 3.1 C2-Phenanthrene/Phene 252 3.1 C2-Phenanthrene/Phyrenes 252 3.1 C2-Phenanthrene/Phyrenes 6.2 3.1 C2-Phenanthrenes/Pyrenes 6.2 3.1 C2-Phenanthrenes/Pyrenes 77.2 10.1 95.4 72.5 96.2 C3-Phenanthrenes/Pyrenes 77.2 10.1 95.4 72.5 96.2 C3-Phenanthrenes/Pyrenes 77.2 10.1 95.4 72.5 96.2 C3-Phenanthrenes/Pyrenes 60.7 12.5 58.9 50.1 67.7 C3-Phenanthrenes/Pyrenes 60.7 12.5 75.9 90.1 C3-Phenanthrenes/Pyrenes 60.7 12.5 75.9 90.1 C3-Phenanthrenes/Pyrenes 77.2 10.1 95.4 72.5 96.2 C3-Phenanthrenes/Pyrenes 60.7 12.5 75.9 90.1 C3-Phenanthrenes/Pyrenes 60.7 12.5 75.9 C3-Phenanthrenes/Pyrenes 60.7 12.5 75.0 C3-Phenanthrenes/Pyrenes 60.7 12.5 75.0 C3-Phenanthrenes/Pyrenes 60.7 12.5 75.0 C3-Phenanthrenes/Pyrenes 60.7 12.5 75.0 C3-Phenanthrenes/Pyrenes 60.7	Carbazole Anthracene	3.8	J				
C2-Pinearsthrene/Anthracenes	Phenanthrene			100 ± 7.0			
C2-Phenathrene/Anthracenes	C1-Phenanthrene/Anthracenes C2-Phenanthrene/Anthracenes	516	5.1		543	462	624
Disherosthophene 30,7	C3-Phenanthrene/Anthracenes						
C1-Diberzothiophere	Dibenzothiophene	30.7	14.5	32.9 ± 1.7	35.5	30.2	40.8
C3-Diberozothiophene 252 0.8 250 213 288	C1-Dibenzothiophene						
Fluoranthene 5.2	C3-Dibenzothiophene	252	0.8				
CiFluoranthenes/Pyrenes	Fluoranthene						
C2-Flucranthenes/Pyrenes	C1-Fluoranthenes/Pyrenes	60.7	12.5				
Naphthobenzothiophene	C2-Fluoranthenes/Pyrenes						
C2-Alaphthobenzothiophenes	Naphthobenzothiophene	34.8	13.4		39.8	33.8	45.8
Second Series	C1-Naphthobenzothiophenes						
Benz(a)anthracene	C3-Naphthobenzothiophenes	55.2	0.0				
C1-Chrysenes					216	18 4	24.8
C3-Chrysenes	C1-Chrysenes	68.7	0.4		68.4	58.1	78.7
C4-Chrysenes	C2-Chrysenes						
Benzo(k)fluoranthene	C4-Chrysenes	<10	U		05.0	7-2.11	
Benzo(e)pyrene	Benzo(b)fluoranthene						
Perylene	Benzo(e)pyrene	2.9	J				
Indeno(1,2,3-c,d)pyrene				30.2+1.7	33.5	28.4	38.5
Total PAHs 7829	Indene(1,2,3-c,d)pyrene	2.0	J	50.4.Z 1.1	65.0	20.4	00.0
Total PAHs 7829	Dibenzo(a,h)anthracene						
Selected Ratios D2/P2							
D3/P3	Total PAHs Selected Ratios	7829					
D3/P3	D2/P2	0.483	1.9		0.473	0.402	0.544
D3/C3	D3/P3						
D3/C3	D2/C2	2.008	2.4		2.056		
FI-Py3/C3 0.896 7.5 0.965 0.820 1.110 Individual Aikyl Isomers and Hopane 2-Methylnaphthalene 527 13.3 602 512 692 1-Methylnaphthalene 386 7.2 415 353 477 2,6-Dimethylnaphthalene 545 9.9 602 512 692 1.6,7-Trimethylnaphthalene 188 10.0 152 129 175 1-Methylphenanthrene 88.0 12.8 100 85.0 115 C29-Hopane 198 16a-Oleanane 64.5 C30-Hopane 295 1.4 291 239 323 Surrogate (Su) Su Recovery (%) Naphthalene-d8 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	D3/C3						
FI-Py3/C3 0.896 7.5 0.965 0.820 1.110 Individual Aikyl Isomers and Hopane 2-Methylnaphthalene 527 13.3 602 512 692 1-Methylnaphthalene 386 7.2 415 353 477 2,6-Dimethylnaphthalene 545 9.9 602 512 692 1-Methylphaphthalene 188 10.0 152 129 175 1-Methylphenanthrene 88.0 12.8 100 85.0 115 C29-Hopane 198 18a-Oleanane 64.5 C30-Hopane 295 1.4 291 239 323 Surrogate (Su) Su Recovery (%) Naphthalene-d8 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	FI-Pv2/C2	0.831	1.1		0.840	0.714	0.966
2-Methylnaphthalene 527 13.3 602 512 692 1-Methylnaphthalene 386 7.2 415 353 477 2,6-Dimethylnaphthalene 545 9.9 602 512 692 1.6,7-Trimethylnaphthalene 188 10.0 152 129 175 1.4-Methylphenanthrene 88.0 12.8 100 85.0 115 C29-Hopane 198 18a-Oleanane 64.5 C30-Hopane 295 1.4 291 239 323 Surrogate (Su) Su Recovery (%) Naphthalene-d8 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	FI-Py3/C3						1,110
1-Metrylnaphthalene 386 7.2 415 353 477 2,6-Dimetrylnaphthalene 545 9.9 602 512 692 1,6,7-Timetrylnaphthalene 188 10.0 152 129 175 1-Metrylnaphthalene 88.0 12.8 100 85.0 115 C29-Hopane 198 16a-Oleanane 64.5 C30-Hopane 295 1.4 291 239 323 Surrogate (Su) Su Recovery (%) Naphthalene-dB 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	Individual Alkyl Isomers and Hopane						
1-Methylnaphthalene 386 7.2 415 353 477 2,6-Dimethylnaphthalene 545 9.9 602 512 692 16,7-Timethylnaphthalene 188.0 12.8 100 85.0 115 C29-Hopane 198 18a-Oleanane 64.5 295 1.4 291 239 323 Surrogate (Su) Su Recovery (%) Naphthalene-d8 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	2-Methylnaphthalene						
1.6,7-Trimethylnaphthalene 188 10.0 152 129 175 1-Methylphenanthrene 88.0 12.8 100 85.0 115 C29-Hopane 198 18a-Oleanane 64.5 5 C30-Hopane 295 1.4 291 239 323 Surrogate (Su) Su Recovery (%) Naphthalene-d8 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	1-Methylnaphthalene						
C29-Hopane 198 18a-Oleanane 64.5 C30-Hopane 295 1.4 291 239 323 Surrogate (Su) Su Recovery (%) Naphthalene-d8 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	2,6-Limethylnaphthalene 1,6,7-Trimethylnaphthalene	168	10.0		152	129	175
18a-Oleanane 54.5 C30-Hopane 295 1.4 291 239 323 Surrogate (Su) Su Recovery (%) Naphthalene-d8 94 99 </td <td>1-Methylphenanthrene</td> <td></td> <td>12.8</td> <td></td> <td>100</td> <td>85.0</td> <td>115</td>	1-Methylphenanthrene		12.8		100	85.0	115
Surrogate (Su) Su Recovery (%) Naphthalene-d8 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	18a-Oleanane	64.5					
Naphthalene-dB 94 Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	C30-Hopane		1.4		291	239	323
Acenaphthene-d10 99 Phenanthrene-d10 97 Chrysene-d12 90	Surrogate (Su)	Su Recovery (%)					
Phenanthrene-d10 97 Chrysene-d12 90	Naphthalene-d8						
Chrysene-d12 90							
	Chrysene-d12	90					
Perylene-d12 91	Perylene-d12	91					

Qualifiers (Q): J=Below the MDL, U=Not detected, B=In procedural blank > 3x MDL, I=Interference, D=Diluted value, NA=Not Applicable, *=Outside QA limits, refer to narrative

Geoinsight, Inc. Buzzards Bay Oil Spill Project Polycyclic Aromatic Hydrocarbon Data Laboratory Control Standard Report

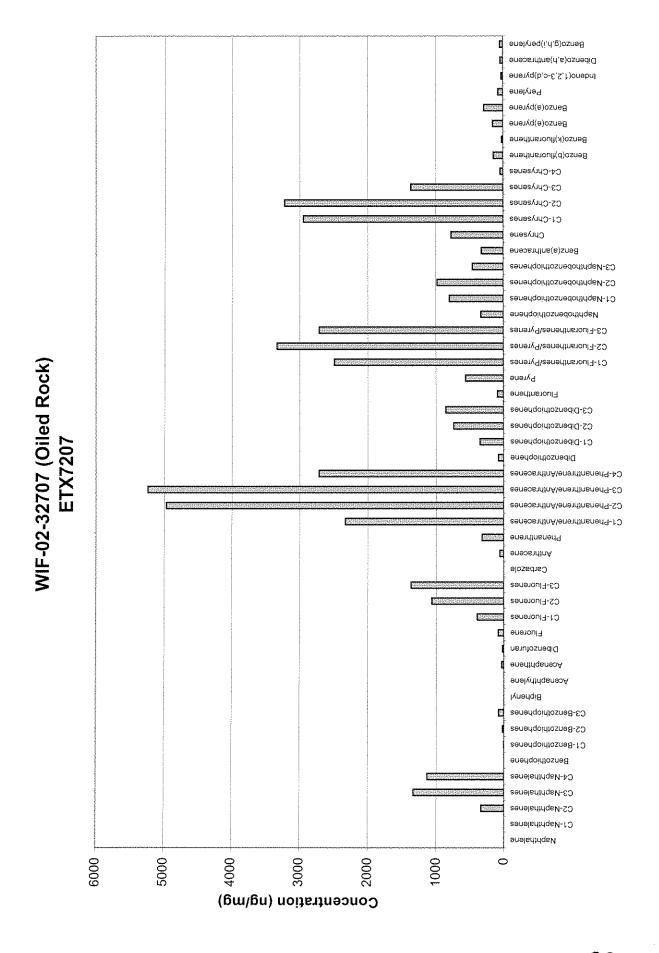
Sample Name	M\$30377J.D
Client Name	AR-WKCC-250-022
Matrix	Solution
Collection Date	NA
Received Date	NA
Extraction Date	NA NA
Extraction Batch	ENV 1635
Date Acquired	04/13/07
Method	PAH-2002
Sample Volume (mL)	1

						······································
Target Compounds	Conc. (ng/ml) C		LCS	-15%	+15%	
		(%)	Certified Conc.	Conc.	Conc.	
			Conc. (ng/ml)	Conc. (ng/ml)	Conc. (ng/ml)	
fanhfhalaga	247	-2.2	959	215	202	
Naphthalene		-2.2	253	215	290	
C1-Naphthalenes	NA					
C2-Naphthalenes	NA					
03-Naphthalenes	NA					
C4-Naphthalenes	NA					
Benzothiophene	241	-3.9	251	213	288	
21-Benzothiophenes	NA					
C2-Benzothiophenes	NA					
23-Benzothiophenes	NA					
Biphenyl	256	2.2	250	213	288	
cenaphthylene	255	1.9	250	213	288	
scenaphthene	247	-1.4	251	213	288	
Dibenzofuran	247					
luorene	253	1.0	251	213	288	
1-Fluorenes	NA					
2-Fluorenes	NΑ					
3-Fluorenes	NA					
arbazole	259	3.4	250	213	288	
nthracene	258	3.0	250	213	288	
henanthrene	253	1.0	251	213	288	
1-Phenanthrene/Anthracenes	NA		***			
2-Phenanthrene/Anthracenes	NA					
3-Phenanthrene/Anthracenes	NA					
4-Phenanthrene/Anthracenes	NA					
ibenzothiophene	257	2.6	250	213	288	
1-Dibenzothiophenes	NA	L .0	200	210	200	
2-Dibenzothiophenes	NA					
C3-Dibenzothiophenes	NA					
luoranthene	239	-4.7	251	213	288	
угеле	259 251	0.2	251	213	288	
		υ.∠	£0;	213	∠88	
21-Fluoranthenes/Pyrenes	NA					
2-Fluoranthenes/Pyrenes	NA ALA					
23-Fluoranthenes/Pyrenes	NA 200		646	0.10	607	
laphthobenzothiophene	260	4.0	250	212	287	
1-Naphthobenzothiophenes	NA					
2-Naphthobenzothiophenes	NA					
3-Naphthobenzothiophenes	NA .					
enz(a)anthracene	269	7.1	251	213	288	
hrysene	260	3.7	251	213	288	
1-Chrysenes	NA					
2-Chrysenes	NA					
3-Chrysenes	NA					
4-Chrysenes	NA					
enzo(b)fluoranthene	237	-5.5	250	213	288	
enzo(k)fluoranthene	239	-4.7	251	213	288	
enzo(e)pyrene	220	-13.0	251	213	288	
enzo(a)pyrene	242	-3.4	250	213	288	
erylene	235	-6.4	250	213	288	
ndeno(1,2,3-c,d)pyrene	265	5.6	251	213	288	
ibenzo(a,h)anthracene	269	7.2	250	213	288	
enzo(g,h,i)perylene	241	-3.9	250	213	288	
dividual Alkyl Isomers and Hopanes						
Methylnaphthalene	253	0.9	251	213	288	
-Methylnaphthalene	242	-3.5	251	213	288	
,6-Dimethylnaphthalene	252	0.6	251	213	288	
6,7-Trimethylnaphthalene	258	3.0	250	213	288	
-Methylphenanthrene	261	4.1	251	213	288	
29-Hopane	NA					
8a-Oleanane	NA					
30-Hopane	234	-6.6	250	213	288	

Surrogate (Su)	Su Recovery (%)
Naphthalene-d8	97
Acenaphthene-d10	100
Phenanthrene-d10	94
Chrysene-d12	108
Perylene-d12	92

Qualifiers (Q): J=Below the MDL, U=Not detected, B=In procedural blank > 3x MDL, (=Interference, D=Diluted value, NA=Not Applicable, *=Outside QA limits, refer to narrative

Polycyclic Aromatic Hydrocarbon Histograms



Polycyclic Aromatic Hydrocarbon Total Ion Chromatograms

54.00 40.00 42.00 44.00 46.00 48.00 50.00 52.00 24.00 26.00 28.00 30.00 32.00 34.00 36.00 38.00 TIC: ETX7207.D 10.00 12.00 14.00 16.00 18.00 20.00 22.00 Abundance 450000 000009 550000 500000 400000 350000 300000 250000 200002 150000 100000 Time-->

pm using AcgMethod PAH-2002

6:43

13 Apr 2007 GC/MS Ins WIF-02-32707

Instrument

Operator Acquired

File

Sample Name: Misc Info : Vial Number:

X:\1\DATA\MS30377\ETX7207.D

Total Petroleum Hydrocarbons/ Aliphatic Hydrocarbons Raw Data

Sequence Name: W:\2\SEQUENCE\GC10846.S

Comment: GeoInsight-Buzzards Bay Spill-Rock

Operator: TJM

Data Path: C:\HPCHEM\2\DATA\gc10846\

Pre-Seq Cmd: Post-Seq Cmd:

Method Sections To Run On A Barcode Mismatch (X) Full Method (X) Inject Anyway () Reprocessing Only () Don't Inject

Line	е Туре	Vial	DataFile	Method	Sample Name
2 3	Sample Sample Sample	52 53	GC10846B GC10846C	ALI_COMP ALI_COMP	Solvent Blank Diesel Std. PEM (Low)
	Sample Sample	96	GC10846E	ALI_COMP	
	Sample Sample		GC10846F GC10846G		
8	Sample	99	GC10846H	ALI_COMP	CS4
	Sample Sample		GC10846I		CS5 AL-WKCC-25-006 (CCC)
11	Sample	56	ETX7207	ALI COMP	
12	Sample	57	GC10846K	ALI_COMP	AL-WKCC-25-006 (CCC)

Evaluate Continuing Calibration Report

Data File : W:\2\DATA\GC10846\GC10846J.D Vial: 55 Acq On : 13 Apr 2007 6:39 pm Sample : AL-WKCC-25-006 (CCC) Operator: TJM Inst : GC#1 Misc Multiplr: 1.00

IntFile : autoint1.e

Method : W:\2\METHODS\C10B414.M (Chemstation Integrator)
Title : C10 - C35 aliphatic

Last Update : Sat Apr 14 11:05:52 2007 Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min

Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	n-hexadecane-d34	1.000	1.000	0.0	105	0.00
2	n-C10	1.016	1.000	1.6	102	0.00
3	n-C11	1.039	1.029	1.0	102	0.00
4 S	n-dodecane-d26	0.865	0.861	0.5	103	0.00
5	n-C12	1.085	1.077	0.7	102	0.00
6	n-C13	1.119	1.115	0.4	101	0.00
7	n-C14	1.164	1.142	1.9	101	0.00
8	n-C15	1.183	1.175	0.7	101	0.00
9	n-C16	1.201	1.183	1.5	101	0.00
10	5a-androstane	1.000	1.000	0.0	104	0.00
11	n-C17	0.929	0.924	0.5	101	0.00
12	Pristane	0.987	0.971	1.6	101	-0.01
13	n-C18	0.933	0.930	0.3	101	-0.01
14	Phytane	0.989	0.970	1.9	101	-0.01
15	n-C19	0.934	0.926	0.9	101	-0.01
16 S	n-eicosane-d42	0.832	0.831	0.1	101	-0.01
17	n-C20	0.928	0.927	0.1	101	-0.01
18	n-C21	0.942	0.947	-0.5	101	-0.01
19	n-C22	0.924	0.915	1.0	101	-0.01
20	n-C23	0.918	0.921	-0.3	101	-0.02
21	n-C24	0.912	0.908	0.4	100	-0.01
22	n-C25	0.906	0.896	1.1	100	-0.01
23	n-C26	0.904	0.899	0.6	100	-0.02
24	n-C27	0.877	0.868	1.0	100	-0.01
25	n-C28	0.859	0.852	0.8	100	-0.02
26	n-C29	0.871	0.854	2.0	100	-0.02
27 S	n-triacontane-d62	0.724	0.714	1.4	100	-0.01
28	n-C30	0.827	0.816	1.3	100	-0.02
29	n-C31	0.803	0.793	1.2	100	-0.01
30	n-C32	0.770	0.765	0.6	101	-0.01
31	n-C33	0.747	0.744	0.4	100	-0.02
32	n-C34	0.730	0.727	0.4	101	-0.02

Evaluate Continuing Calibration Report - Not Founds

Data File : W:\2\DATA\GC10846\GC10846J.D Vial: 55 Acq On : 13 Apr 2007 6:39 pm Operator: TJM Sample : AL-WKCC-25-006 (CCC) Inst : GC#1 Misc Multiplr: 1.00

IntFile : autoint1.e

Method : W:\2\METHODS\C10B414.M (Chemstation Integrator)
Title : C10 - C35 aliphatic
Last Update : Sat Apr 14 11:05:52 2007

Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min

Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev A	Area% Dev(min)
33	ТРН	0.043	0.000	100.0#	0# -22.30#
34	TRH1	0.043	0.000	100.0#	0# -5.85#
35	TRH2	0.043	0.000	100.0#	0# -12.65#
36	TRH3	0.043	0.000	100.0#	0# -19.76#
37	TRH4	0.043	0.000	100.0#	0# -27.15#
38	TRH5	0.043	0.000	100.0#	0# -34.22#
39	TRH6	0.043	0.000	100.0#	0# -42.15#

Quantitation Report (QT Reviewed)

Vial: 55 Data File: W:\2\DATA\GC10846\GC10846J.D Acq On : 13 Apr 2007 6:39 pm Sample : AL-WKCC-25-006 (CCC) Misc : IntFile : autoint1.e Operator: TJM Inst : GC#1 Multiplr: 1.00

Quant Time: Apr 14 11:20 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic

Last Update : Sat Apr 14 11:05:52 2007

Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase: Signal Info :

	Compound	R.T.	Response	Conc Units
				. * *
Inte	ernal Standards			
1)	n-hexadecane-d34	12.85	397397	20.001 ug/mlm
10)	5a-androstane	18.03	516981	20.003 ug/mlm
Syst	tem Monitoring Compounds			
4) S	n-dodecane-d26	8.62	428216	24.909 ug/mlm
16) S	n-eicosane-d42	17.44	537070	24.955 ug/mlm
•	n-triacontane-d62	29.23	462148	24.668 ug/mlm
Taro	get Compounds			
2)	n-C10	6.26	495098	24.511 ug/mlm
3)	n-C11	7.59	509794	24.681 ug/mlm
	n-C12	8.83	533630	24.730 ug/mlm
	n-C13	9.99	551635	24.791 ug/mlm
7)	n-C14	11.07	567676	24.541 ug/mlm
8)	n-C15	12.09	580895	24.701 ug/mlm
9)	n-C16	13.11	587196	24.601 ug/mlm
11)	n-C17	14.20	597091	24.843 ug/mlm
	Pristane	14.31	628118	24.598 ug/mlm
13)	n-C18	15.35	602028	24.929 ug/mlm
14)	Phytane	15.52	617587	24.132 ug/mlm
15)	n-C19	16.57	599261	24.803 ug/mlm
17)	n-C20	17.84	599851	24.964 ug/mlm
18)	n-C21	19.12	611720	25.080 ug/mlm
	n-C22	20.41	580661	24.283 ug/mlm
20)	n-C23	21.68	595299	25.050 ug/mlm
21)	n-C24	22.93	587550	24.881 ug/mlm
22)	n-C25	24.16	580618	24.772 ug/mlm
23)	n-C26	25.35	577646	24.688 ug/mlm
24)	n-C27	26.50	555597	24.479 ug/mlm
25)	n-C28	27.62	549551	24.714 ug/mlm
26)	n-C29	28.70	552158	24.494 ug/mlm
28)	n-C30	29.76	527874	24.663 ug/mlm
29)	n-C31	30.78	512845	24.691 ug/mlm
30)	n-C32	31.77	487349	24.460 ug/mlm
31)	n-C33	32.76	480155	24.853 ug/mlm
32)	n-C34	33.89	468598	24.792 ug/mlm

Data File: W:\2\DATA\GC10846\GC10846J.D Vial: 55 Acq On : 13 Apr 2007 6:39 pm Sample : AL-WKCC-25-006 (CCC) Operator: TJM Inst : GC#1 Misc : IntFile : autointl.e Multiplr: 1.00

Quant Time: Apr 14 11:20 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic Last Update : Sat Apr 14 11:05:52 2007

Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

> R.T. Response Conc Units Compound

 $AL-W\bar{K}CC-25-006$ (CCC)

GC#1 1.00

Inst

Multiplr:

MPL

Vial: Operator:

Misc

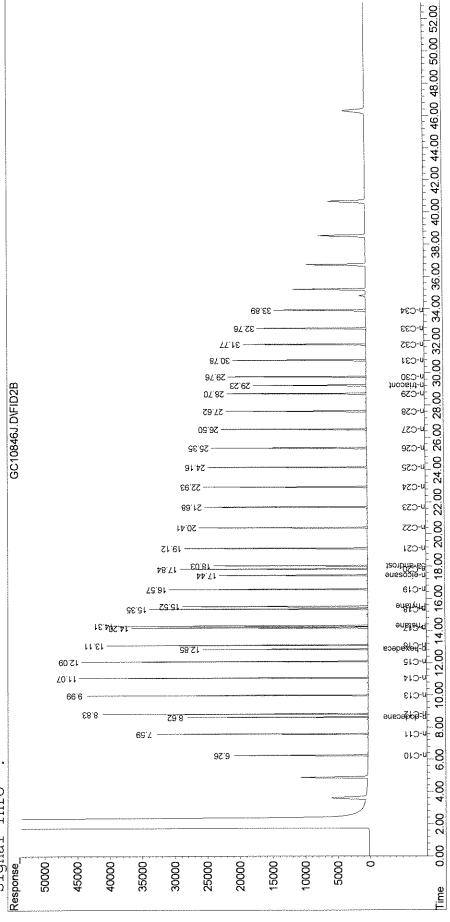
IntFile : autointl.e Quant Time: Apr 14 11:20 2007 Quant Results File: C10B414.RES

C:\GC10846\C10B414.M (Chemstation Integrator) C10 - C35 aliphatic Quant Method Title

: Sat Apr 14 11:05:52 2007 Last Update

Multiple Level Calibration ALI_COMP.M Response via DataAcq Meth

Volume Inj. Siqnal Phase Info Signal Signal



Evaluate Continuing Calibration Report

Data File : W:\2\DATA\GC10846\GC10846K.D Vial: 57 Acq On : 13 Apr 2007 8:40 pm Sample : AL-WKCC-25-006(CCC) Operator: TJM Inst : GC#1 Multiplr: 1.00 Misc

IntFile : autoint1.e

Method : W:\2\METHODS\C10B414.M (Chemstation Integrator)
Title : C10 - C35 aliphatic

Last Update : Sat Apr 14 11:05:52 2007 Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min

Max. RRF Dev: 25% Max. Rel. Area: 150%

	Compound	AvgRF	CCRF	%Dev	Area% Dev(min)
1 2 3 4 5 6 7 8 9	n-hexadecane-d34 n-C10 n-C11 n-dodecane-d26 n-C12 n-C13 n-C14 n-C15 n-C16	1.000 1.016 1.039 0.865 1.085 1.119 1.164 1.183 1.201	1.000 1.035 1.043 0.869 1.094 1.123 1.147 1.167	0.0 -1.9 -0.4 -0.5 -0.8 -0.4 1.5 1.4 2.8	101 0.00 101 0.00 100 0.00 100 0.00 100 0.00 99 0.00 98 0.00 97 0.00 96 0.00
10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37	5a-androstane n-C17 Pristane n-C18 Phytane n-C19 n-eicosane-d42 n-C20 n-C21 n-C22 n-C23 n-C24 n-C25 n-C26 n-C27 n-C28 n-C29 n-triacontane-d62 n-C30 n-C31 n-C32 n-C31 TPH TRH1 TRH1 TRH2 TRH3 TRH4	1.000 0.929 0.987 0.933 0.989 0.934 0.832 0.928 0.942 0.918 0.912 0.906 0.912 0.906 0.877 0.859 0.871 0.724 0.827 0.827 0.827 0.850 0.850 0.850 0.850	1.000 0.958 0.995 0.995 0.987 0.949 0.847 0.943 0.952 0.916 0.907 0.886 0.850 0.850 0.850 0.816 0.773 0.643 0.744 0.688 0.664 0.609 0.606 0.003 0.017 0.003 0.002 0.009	0.186268619291005320385060689 	94
38 39	TRH5 TRH6	0.850 0.850	0.009 0.012	98.9# 98.6#	1# 0.02

Evaluate Continuing Calibration Report - Not Founds

Data File : W:\2\DATA\GC10846\GC10846K.D

Vial: 57 Acq On : 13 Apr 2007 8:40 pm Sample : AL-WKCC-25-006(CCC) Operator: TJM Inst : GC#1 Multiplr: 1.00 Misc

IntFile : autoint1.e

Method : W:\2\METHODS\C10B414.M (Chemstation Integrator)
Title : C10 - C35 aliphatic
Last Update : Sat Apr 14 11:05:52 2007

Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min

Max. RRF Dev : 25% Max. Rel. Area : 150%

AvgRF CCRF %Dev Area% Dev(min) Compound

Vial: 57 Data File: W:\2\DATA\GC10846\GC10846K.D Acq On : 13 Apr 2007 8:40 pm Operator: TJM : AL-WKCC-25-006 (CCC) Sample Inst : GC#1 Misc : Multiplr: 1.00

IntFile : autointl.e

Quant Time: Apr 14 11:20 2007 Quant Results File: C10B414.RES

Quant Method: C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic Last Update : Sat Apr 14 11:05:52 2007

Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

Compound	R.T.	Response	Conc Units
			· · · · · · · · · · · · · · · · · · ·
Internal Standards			
1) n-hexadecane-d34	12.86	383385	20.001 ug/ml
10) 5a-androstane	18.04	470378	20.003 ug/ml
System Monitoring Compounds			
4) S n-dodecane-d26	8.63	416961	25.141 ug/ml
16) S n-eicosane-d42	17.44	497979	25.432 uq/ml
27) S n-triacontane-d62	29.23	378540	22.207 ug/ml
Manage Company de			
Target Compounds 2) n-C10	6.26	494285	25.365 ug/ml
2) n-C10 3) n-C11	7.59	498779	25.030 ug/ml
5) n-C12	8.83	522699	25.030 ug/ml
6) n-C13	9.99	536167	24.976 ug/ml
7) n-C14	11.07	550195	24.655 ug/ml
8) n-C15	12.09	556634	24.635 ug/ml
9) n-C16	13.11	558932	24.273 ug/ml
	14.20	563197	
·	14.20	585240	25.754 ug/ml 25.189 ug/ml
,		564000	25.668 ug/ml
	15.36 15.52	571931	24.562 ug/ml
14) Phytane	16.58		25.416 ug/ml
15) n-C19		558724	
17) n-C20	17.84 19.12	555055	25.388 ug/ml
18) n-C21		559862 539667	25.228 ug/ml
19) n-C22	20.41	528667	24.299 ug/ml
20) n-C23	21.69	533893	24.692 ug/ml
21) n-C24	22.94	521673	24.280 ug/ml
22) n-C25	24.16	506781	23.764 ug/ml
23) n-C26	25.35	496808	23.337 ug/ml
24) n-C27	26.50	475363	23.019 ug/ml
25) n-C28	27.62	466799	23.072 ug/ml
26) n-C29	28.71	454748	22.172 ug/ml
28) n-C30	29.76	438083	22.496 ug/ml
29) n-C31	30.78	404939	21.427 ug/ml
30) n-C32	31.77	385171	21.247 ug/ml
31) n-C33	32.76	357626	20.345 ug/ml
32) n-C34	33.89	355303	20.660 ug/ml

Vial: 57 Data File: W:\2\DATA\GC10846\GC10846K.D Acq On : 13 Apr 2007 8:40 pm Sample : AL-WKCC-25-006 (CCC) Operator: TJM Sample : AL-WKCC-25-Misc : IntFile : autoint1.e Inst : GC#1 Multiplr: 1.00

Quant Time: Apr 14 11:20 2007 Quant Results File: C10B414.RES

Quant Method: C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic

Last Update : Sat Apr 14 11:05:52 2007

Response via : Initial Calibration

DataAcq Meth : ALI_COMP.M

Volume Inj. : Signal Phase: Signal Info :

	Compound	R.T.	Response	Conc Units
33)	ТРН	22.47	1668	0.083 ug/ml
34)	TRH1	5.88	10284	0.514 ug/ml
35)	TRH2	12.62	1541	0.077 ug/ml
36)	TRH3	19.79	998	0.050 ug/ml
37)	TRH4	27.14	5223	0.261 ug/ml
38)	TRH5	34.24	5581	0.279 ug/ml
39)	TRH6	42.11	6891	0.344 ug/ml

Quantitation Report

Multiplr: W:\2\DATA\GC10846\GC10846K.D 13 Apr 2007 8:40 pm AL-WKCC-25-006 (CCC) Data File Acq On Sample Misc

Operator: Vial: Inst

1.00 GC#1 TUM

> autoint1.e IntFile

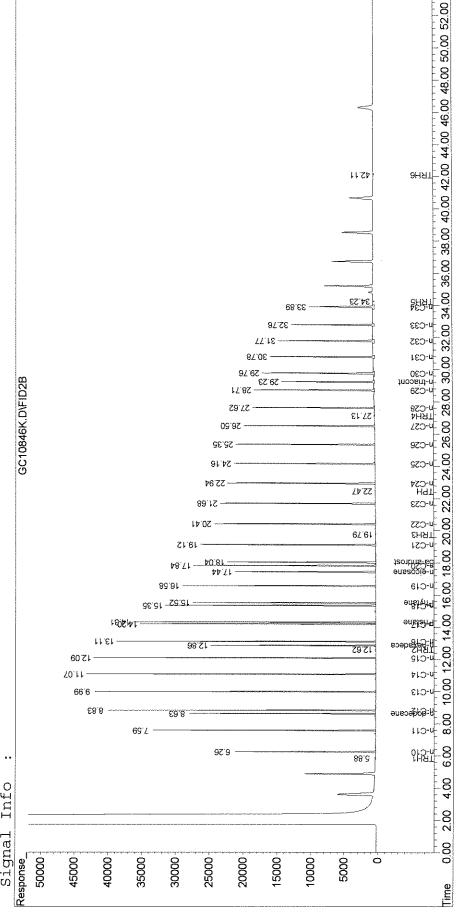
Quant Time: Apr 14 11:20 2007 Quant Results File: C10B414.RES

C:\GC10846\C10B414.M (Chemstation Integrator) C10 - C35 aliphatic Quant Method Title

Multiple Level Calibration Sat Apr 14 11:05:52 2007 Response via Last Update

ALI_COMP.M DataAcq Meth

Phase Info Volume Inj. Signal Signal



000028

U:\2\DATA\GC10846\GC10846B.D

Data File Name GC10846B.D GC10846B Data File Path W:\2\DATA\GC10846\ Diesel S Date Acquired 04/13/20 -1:0: 04/13/20 -1 Sample Name Diesel Std. ALI_COMP.N Sample Multiplier 1				
Name	Amount			20.001
n-hexadecane-d34	20.00			20.003
5a-androstane	20.00			1.870036003
oa androstano	_	Surrogate re	acoverv	1.988539161
n-dodecane-d26	1.87	94	SOOVER Y	1.963768423
n-eicosane-d42	1.99	99		168.8699079
n-triacontane-d62	1.96	98		00.0099079
II-tilacolitalie-uoz	1.90	30		· ·
	c	`urrogoto C	orrootod	0
	-	Surrogate C	onected	0
TOU	400.07	400.04		0
TPH	168.87	169.84		0
TRH1	0.00	0.00		0
TRH2	0.00	0.00		93.50180016
TRH3	0.00	0.00		99.42695807
TRH4	0.00	0.00		98.18842113
TRH5	0.00	0.00		
TRH6	0.00	0.00		

Data File : U:\2\DATA\GC10846\GC10846B.D

Acq On : 13 Apr 2007 10:33 Sample : Diesel Std. Operator: TJM Inst : GC#1

Multiplr: 1.00 Sample Amount: 0.00

Vial: 52

IntFile : autointl.e

Quant Time: Apr 16 7:10 2007 Quant Results File: C10B414.RES

Quant Method: C:\GC10846\C10B414.M (Chemstation Integrator)
Title: C10 - C35 aliphatic
Last Update: Sat Apr 14 11:05:52 2007
Response via: Initial Calibration
DataAcq Meth: ALI_COMP.M

Volume Inj. : Signal Phase : Signal Info :

Misc

	Compound	R.T. Respo		Conc Units
Inte 1) 10)	rnal Standards n-hexadecane-d34 5a-androstane	12.86 18.04	413240 541339	20.001 ug/mlm 20.003 ug/mlm
4) s 16) s	em Monitoring Compounds n-dodecane-d26 n-eicosane-d42 n-triacontane-d62	8.62 17.43 29.21	33430 44812 38525	1.870 ug/mlm 1.989 ug/mlm 1.964 ug/mlm
Targ	et Compounds TPH	12.86	3890476	168.870 ug/mlm

Data File

```
0.00 2.00 4.00 6.00 8.00 10.00 12.00 14.00 16.00 18.00 20.00 22.00 24.00 26.00 28.00 30.00 32.00 34.00 36.00 36.00 40.00 42.00 44.00 46.00 48.00 50.00 52.00
                            GC#1
1.00
0.00
               TUM
                                           Multiplr:
                                                       Sample Amount:
 Vial:
                                                                                              Quant Results File: C10B414.RES
                Operator:
                                                                                                                       C:\GC10846\C10B414.M (Chemstation Integrator)
C10 - C35 aliphatic
Sat Apr 14 11:05:52 2007
Multiple Level Calibration
ALI_COMP.M
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U:\2\DATA\GC10846\GC10846B.D
13 Apr 2007 10:33
Diesel Std.
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                                                                                              7:10 2007
                                                                                                                                                                                                                                                                                                     42.86
                                                                    autoint1.e
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                                                                                                                                                                                                       Volume Inj.
Signal Phase
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Title
                                                                                                                                                                 Response via
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                                                                                                                                                   Last Update
                                                                                                                                                                                                                                     Info
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06:47:26 2007

Tue May 08

C10B414.M

U:\2\DATA\GC10846\GC10846C.D

GC10846C.D

99.95199174

94.73468659

	QQ 100 100.D			
	PEM (Low)			
	04/13/20 -1:1:			
	Samp	le Name Pi	EM (Low)	ALI_COMP.M
	Sample	Multiplier	1	
	·			1
<u>Name</u>	<u>Amount</u>			20.001
n-hexadecane-d34	20.00			20.003
5a-androstane	20.00			1.993664653
	S	urrogate rec	covery	1.999039835
n-dodecane-d26	1.99	100		1.894693732
n-eicosane-d42	2.00	100		68.92844971
n-triacontane-d62	1.89	95		10.48498879
				44.65906223
	S	urrogate Co	rrected	1.605688657
				4.795428516
TPH	68.93	68.96		2.089768265
TRH1	10.48	10.49		0.3763138171
TRH2	44.66	44.68		99.68323263

1.61

4.80

2.09

0.38

TRH3

TRH4

TRH5

TRH6

1.61

4.80

0.38

2.09

Data File Name GC10846C.D

Data File : U:\2\DATA\GC10846\GC10846C.D

Vial: 53 Acq On : 13 Apr 2007 11:34 Sample : PEM (Low) Operator: TJM

Inst : GC#1 Multiplr: 1.00 Misc Sample Amount: 0.00

IntFile : autoint1.e

Quant Time: Apr 14 17:31 2007 Quant Results File: C10B414.RES

Quant Method: C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic

Last Update : Sat Apr 14 11:05:52 2007 Response via : Initial Calibration

DataAcq Meth : ALI_COMP.M

Volume Inj. : Signal Phase : Signal Info :

	Compound	R.T.	Response	Conc Units
Inte:	rnal Standards n-hexadecane-d34 5a-androstane	12.85 18.03	346426 469207	20.001 ug/mlm 20.003 ug/mlm
	em Monitoring Compounds	8.62	29877	1 004 220 /mlm
4) S 16) S	n-dodecane-d26 n-eicosane-d42	17.43	39046	1.994 ug/mlm 1.999 ug/mlm
27) S	n-triacontane-d62	29.21	32217	1.895 ug/mlm
Targ	et Compounds			
33)	TPH	12.85	1376399	68.928 ug/mlm
34)	TRH1	8.62	209370	10.485 ug/mlm
35)	TRH2	12.85	891775	44.659 ug/mlm
36)	TRH3	20.12	32063	1.606 ug/mlm
37)	TRH4	29.21	95758	4.795 ug/mlm
38)	TRH5	38.83	41730	2.090 ug/mlm
39)	TRH6	42.38	7514	0.376 ug/mlm

Quantitation Report

GC#1 1.00 TUM Vial: Multiplr: Sample Amount: Operator: Inst U:\2\DATA\GC10846\GC10846C.D 13 Apr 2007 11:34 PEM (LOW) Data File Acq On Sample Misc

Quant Results File: C10B414.RES Quant Time: Apr 14 17:31 2007 autoint1.e IntFile

C:\GC10846\C10B414.M (Chemstation Integrator)
C10 - C35 aliphatic
Sat Apr 14 11:05:52 2007 Multiple Level Calibration ALI_COMP.M Quant Method DataAcq Meth Response via Last Update Title

Inj. Phase

Volume

Signal

0.00 2.00 4.00 6.00 8.00 10.00 12.00 14.00 16.00 18.00 20.00 22.00 24.00 26.00 30.00 32.00 34.00 36.00 38.00 40.00 42.00 44.00 46.00 48.00 50.00 52.00 8E.S4 9HATснят GC10846C.D\FID2B 12.62 Ino state 20,12 **ЕНЯТ** anseooia-n 5a-androst 18.03 **业外域**Sadeca 12.85 \$3.8 Info Signal Response 26000 16000 8000 0009 2000 -2000 24000 22000 20000 18000 14000 12000 10000 4000 Ó Time 000034

GC10846C.D C10B414.M Tue May

Tue May 08 06:47:54 2007

U:\2\DATA\GC10846\GC10846D.D

	Data Date Samp	File Path \ Acquired (GC10846D.D W:\2\DATA\GC10846\ 04/13/20 -1:2: PEM (High) 0.5657	GC10846D.D PEM (High) 04/13/20 -1:2: ALI_COMP.M
				0.5657
<u>Name</u>	<u>Amount</u>			11.3145657
n-hexadecane-d34	11.31			11.3156971
5a-androstane	11.32			1.094000567
	S	urrogate re	ecovery	1.119919673
n-dodecane-d26	1.09	97		1.11541136
n-eicosane-d42	1.12	99		28.76247931
n-triacontane-d62	1.12	99		1.340272428
				24.69139648
	S	urrogate C	Corrected	0.05657297687
				1.055525332
TPH	28.76	29.06		0.1073976302
TRH1	1.34	1.35		0.03624447017
TRH2	24.69	24.94		96.6944111
TRH3	0.06	0.06		98.98529905
TRH4	1.06	1.07		98.58682698
TRH5	0.11	0.11		
TRH6	0.04	0.04		

Data File : U:\2\DATA\GC10846\GC10846D.D
Acq On : 13 Apr 2007 12:34
Sample : PEM (High) Vial: 54 Operator: TJM

Inst : GC#1 Multiplr: 0.57 Misc Sample Amount: 0.00

IntFile : autointl.e

Quant Time: Apr 16 7:14 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Sat Apr 14 11:05:52 2007
Response via : Initial Calibration
DataAcq Meth : ALI_COMP.M

Volume Inj. : Signal Phase : Signal Info :

	Compound	R.T.	Response	Conc Units
Inte 1) 10)	rnal Standards n-hexadecane-d34 5a-androstane	12.89f 18.10f	3645316 4820418	11.315 ug/mlm 11.316 ug/mlm
4) Ŝ 16) S	em Monitoring Compounds n-dodecane-d26 n-eicosane-d42 n-triacontane-d62	8.62 17.44 29.22	304963 397262 344441	1.094 ug/mlm 1.120 ug/mlm 1.115 ug/mlm
Targ 33) 34) 35) 36) 37) 38)	et Compounds TPH TRH1 TRH2 TRH3 TRH4 TRH5	12.89 8.62 12.89 20.12 29.22 34.47	10430512 486040 8954162 20516 382779 38947	28.762 ug/mlm 1.340 ug/mlm 24.691 ug/mlm 0.057 ug/mlm 1.056 ug/mlm 0.107 ug/mlm
39)	TRH6	41.33	13144	0.036 ug/mlm

Quantitation Report

GC#1 0.57 0.00 TUM 54 Vial: Multiplr: Operator: Inst U:\2\DATA\GC10846\GC10846D.D 13 Apr 2007 12:34 PEM (High) Data File Acq On Sample Misc

Sample Amount:

autointl.e IntFile

Quant Results File: C10B414.RES 7:14 2007 Quant Time: Apr 16

C:\GC10846\C10B414.M (Chemstation Integrator)
C10 - C35 aliphatic
Sat Apr 14 11:05:52 2007 Quant Method Last Update Title

Multiple Level Calibration ALI_COMP.M Response via DataAcq Meth

Volume Inj. Signal Phase

10,00 12.00 14,00 16.00 18.00 20.00 22.00 24.00 26.00 28.00 30.00 32.00 34.00 36.00 38.00 40.00 42.00 44.00 46.00 48.00 50.00 52.00 41.33 энят, ፈ⊅"⊅ይ ТВН5 GC10846D,D\FID2B 25.22 20.12 твнз 01.81 5a-androst ÞÞ.71 12.89 29.8 # Rated ecane 0.00 2.00 4.00 6.00 8.00 Info Signal 140000 120000 1000001 80000 00009 40000 20000 0 Response 180000 160000 000037

C10B414.M GC10846D.D

U:\2\DATA\GC10846\ETX7207.D

	ETX7207.D WIF-02-32707 04/13/20 -1:9: ALI_COMP.M			
				0.5657
<u>Name</u>	<u>Amount</u>			11.3145657
n-hexadecane-d34	11.31			11.3156971
5a-androstane	11.32			1.148877043
	S	Surrogate re	ecovery	1.042315411
n-dodecane-d26	1.15	102		1.033983776
n-eicosane-d42	1.04	92		274.0813551
n-triacontane-d62	1.03	91		2.101523507
				60.73020583
	S	Surrogate C	Corrected	20.14145915
				7.825533995
TPH	274.08	297.51		0
TRH1	2.10	2.28		0
TRH2	60.73	65.92		101.5447272
TRH3	20.14	21.86		92.12616325
TRH4	7.83	8.49		91.38976277
TRH5	0.00	0.00		
TRH6	0.00	0.00		

Data File : U:\2\DATA\GC10846\ETX7207.D

Vial: 56 Acq On : 13 Apr 2007 19:40 Sample : WIF-02-32707 Operator: TJM Inst : GC#1

Multiplr: 0.57 Sample Amount: 0.00

IntFile : autoint1.e

Quant Time: Apr 20 16:20 2007 Quant Results File: C10B414.RES

Quant Method: C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic

Last Update : Sat Apr 14 11:05:52 2007 Response via : Initial Calibration DataAcq Meth : ALI_COMP.M

Volume Inj. : Signal Phase : Signal Info :

Misc

	Compound	R.T.	Response	Conc Units
Inte	rnal Standards			
1)	n-hexadecane-d34	12.89f	3697746	11.315 ug/mlm
10)	5a-androstane	18.11f	5216861	11.316 ug/mlm
Syst	em Monitoring Compounds			
4) S	n-dodecane-d26	8.62	324866	1.149 ug/mlm
16) S	n-eicosane-d42	17.46	400142	1.042 ug/mlm
27) S	n-triacontane-d62	29.24	345556	1.034 ug/mlm
Tarq	et Compounds			
33)	TPH	12.89	107568070	274.081 ug/mlm
34)	TRH1	8.62	824780	2.102 ug/mlm
35)	TRH2	12.89	23834642	60.730 ug/mlm
36)	TRH3	24.67	7904871	20.141 ug/mlm
37)	TRH4	29.24	3071269	7.826 ug/mlm

GC#1 0.57 0.00

Multiplr:

Inst

Sample Amount:

Quant Results File: C10B414.RES

Quant Time: Apr 20 16:20 2007

autoint1.e

IntFile

TUM 26

Operator:

U:\2\DATA\GC10846\ETX7207.D 13 Apr 2007 19:40 WIF-02-32707

Data File

Acq On Sample

Misc

Vial:

Tue May 08 06:48:57 2007

Polycyclic Aromatic Hydrocarbon Raw Data

Sequence Name: G:\1\SEQUENCE\MS30377.S

Comment: GeoInsight-Buzzards Bay Spill-Rock

Operator: TJM

Data Path: C:\HPCHEM\1\data\ms30377\

Pre-Seq Cmd: Post-Seq Cmd:

Method Sections To Run On A Barcode Mismatch (X) Full Method (X) Inject Anyway () Reprocessing Only () Don't Inject

Line	e Type	Vial	DataFile	Method	Sample Name
			MC202777	DAII 2002	Solvent Rinse
	Sample	1.	MOSUSIIA	PAR-ZUUZ	SOTAGUE KTURE
2	Sample	2	MS30377B	PAH-2002	SRM 1582
3	Sample	3	MS30377C	PAH-2002	IS/SU Mixture
4	Sample	41	MS30377D	PAH-2002	Cal Level 1
5	Sample	42	MS30377E	PAH-2002	Cal Level 2
6	Sample	43	MS30377F	PAH-2002	Cal Level 3
7	Sample	44	MS30377G	PAH-2002	Cal Level 4
8	Sample	45	MS30377H	PAH-2002	Cal Level 5
9	Sample	4	MS30377I	PAH-2002	AR-WKCC-250-022
10	Sample	5	ETX7207	PAH-2002	
11	Sample	6	MS30377J	PAH-2002	AR-WKCC-250-022

Evaluate Continuing Calibration Report

Data File : G:\1\DATA\MS30377\MS30377I.D Vial: 4 Acq On : 13 Apr 2007 5:40 pm Sample : AR-WKCC-250-022 Operator: TJM

Inst : GC/MS Ins

Misc Multiplr: 1.00

MS Integration Params: rteint.p

Method : G:\1\METHODS\041507.M (RTE Integrator)
Title : PAH Calibration Table (2002)
Last Update : Sat Apr 14 20:04:44 2007 Response via: Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 25% Max. R.T. Dev 0.50min

Max. RRF Dev: 25% Max. Rel. Area: 200%

		Compound	AvgRF	CCRF	%Dev Area% Dev(min)
7	I	Fluorene-d10	1.000	1.000	0.0 82 0.00
2	S	Naphthalene-d8	1.911	1.995	-4.4 87 0.00
3	T	Decalin	0.404	0.422	-4.5 90 0.00
4	un	C1-Decalin	0.404	0.000	100.0# 0# -12.76#
5	un	C2-Decalin	0.404	0.000	100.0# 0# -14.27#
6	un	C3-Decalin	0.404	0.000	100.0# 0# -15.86#
7	un	C4-Decalin	0.404	0.000	100.0# 0# -19.68#
8	\mathbf{T}	Naphthalene	2.087	2.177	-4.3 87 0.00
9	T	2-Methylnaphthalene	1.365	1.448	-6.1 89 0.00
10	${f T}$	1-Methylnaphthalene	1.263	1.319	-4.4 92 0.00
11		2,6-Dimethylnaphthalene	1.111	1.151	-3.6 88 0.00
12	T	1,6,7-Trimethylnaphthalene	1.026	1.103	-7.5 104 0.00
13	un	C2-Naphthalenes	2.087	0.000	100.0# 0# -17.94#
14	un	C3-Naphthalenes	2.087	0.000	100.0# 0# -20.12#
15	un	C4-Naphthalenes	2.087	0.000	100.0# 0# -22.24#
16	T	Benzothiophene	1.767	1.871	-5.9 87 0.00
17		C1-Benzothiophene	1.767	0.000	100.0# 0# -15.58#
18	un	C2-Benzothiophene	1.767	0.000	100.0# 0# -17.97#
19	un	C3-Benzothiophene	1.767	0.000	100.0# 0# -19.61#
20	S	Acenaphthene-d10	0.921	0.968	-5.1 90 0.00
21		Biphenyl	1.667	1.752	-5.1 87 0.00
22		Acenaphthylene	2.012	2.123	-5.5 87 0.00
23		Acenaphthene	1.141	1.227	-7.5 89 0.00
24		Dibenzofuran	1.850	1.922	-3.9 88 0.00
25	\mathbf{T}	Fluorene	1.328	1.379	-3.8 87 0.00
	un	C1-Fluorenes	1.328	0.000	100.0# 0# -22.65#
27	un	C2-Fluorenes	1.328	0.000	100.0# 0# -24.32#
28	un	C3-Fluorenes	1.328	0.000	100.0# 0# -26.19#
29	I	Pyrene-d10	1.000	1.000	0.0 76 0.00
30	S	Phenanthrene-d10	0.849	0.892	-5.1 81 0.00
31		Pentachlorophenol	0.044	0.000	100.0# 0# -23.30#
32		Carbazole	0.988	1.012	-2.4 78 0.00
33	${f T}$	Dibenzothiophene	1.051	1.155	-9.9 85 0.00
	un	C1-Dibenzothiophene	1.051	0.000	100.0# 0# -25.11#
	un	C2-Dibenzothiophene	1.051	0.000	100.0# 0# -26.47#
	un	C3-Dibenzothiophene	1.051	0.000	100.0# 0# -28.57#
37		Phenanthrene	1.006	1.169	-16.2 93 0.00
38		Anthracene	1.116		-17.5 91 0.00
39		1-Methylphenanthrene	0.859		
40	un	C1-Phenanthrene/Anthracene	1.006	0.000	100.0# 0# -26.20#

^{(#) =} Out of Range

Evaluate Continuing Calibration Report

Vial: 4 Data File : G:\1\DATA\MS30377\MS30377I.D Acq On : 13 Apr 2007 5:40 pm Sample : AR-WKCC-250-022 Operator: TJM

Inst : GC/MS Ins

Multiplr: 1.00 Misc

MS Integration Params: rteint.p

: G:\1\METHODS\041507.M (RTE Integrator)

Method : G:\1\METHODS\04150/.M (KID 1)
Title : PAH Calibration Table (2002) Last Update : Sat Apr 14 20:04:44 2007 Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 25% Max. R.T. Dev 0.50min Max. RRF Dev : 25% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev Ai	rea% Dev(min)
41	un	C2-Phenanthrene/Anthracene	1.006	0.000	100.0#	0# -27.51#
42		C3-Phenanthrene/Anthracene	1.006	0.000	100.0#	0# -30.06#
	un	C4-Phenanthrene/Anthracene	1.006	0.000	100.0#	0# -30.78#
44	T	Naphthobenzothiophene	0.983	1.141	-16.1	94 0.00
45		C1-Naphthobenzothiophene	0.983	0.000	100.0#	0# -33.03#
46	un	C2-Naphthobenzothiophene	0.983	0.000	100.0#	0# -34.70#
47	un	C3-Naphthobenzothiophene	0.983	0.000	100.0#	0# -36.06#
48	T	Fluoranthene	1.299	1.311	-0.9	77 0.00
49	T	Pyrene	1.426	1.587	-11.3	87 0.00
50	un	Cl-Fluoranthenes/Pyrenes	1.299	0.000	100.0#	0# -30.52#
51	un	C2-Fluoranthenes/Pyrenes	1.299	0.000	100.0#	0# -31.88#
52	un	C3-Fluoranthenes/Pyrenes	1.299	0.000	100.0#	0# -32.93#
53	S	Chrysene-d12	1.147	1.362	-18.7	94 0.00
54	T	Benz(a)anthracene	1.324	1.473	-11.3	85 0.00
55	T	Chrysene	1.102	1.135	-3.0	77 0.00
56	un	C1-Chrysenes	1.102	0.000	100.0#	0# -34.25#
57	un	C2-Chrysenes	1.102	0.000	100.0#	0# -35.44#
58	un	C3-Chrysenes	1.102	0.000	100.0#	0# -36.87#
59	un	C4-Chrysenes	1.102	0.000	100.0#	0# -42.39#
60	I	Benzo(a)pyrene-d12	1.000	1.000	0.0	77 0.00
61	un	C29-Hopane	0.679	0.000	100.0#	0# -40.43#
62	un	18a-Oleanane	0.679	0.000	100.0#	0# -42.28#
63	T	C30-Hopane	0.679	0.725	-6.8	87 0.00
64	\mathbf{T}	Benzo(b)fluoranthene	1.787	2.047	-14.5	91 0.00
65	T	Benzo(k)fluoranthene	1.729	1.779	-2.9	78 0.00
66	Τ	Benzo(e)pyrene	1.800	2.026	-12.6	86 0.00
67	T	Benzo(a)pyrene	1.529	1.616	-5.7	83 0.00
68	T	Indeno(1,2,3-c,d)pyrene	1.161	1.162	-0.1	83 0.00
69	T	Dibenzo(a,h)anthracene	1.077	1.059	1.7	79 0.03
	un	C1-Dibenzo(a,h)anthracene	1.077	0.000	100.0#	0# -42.56#
	un	C2-Dibenzo(a,h)anthracene	1.077	0.000	100.0#	0# -44.47#
	un	C3-Dibenzo(a,h)anthracene	1.077	0.000	100.0#	0# -44.94#
73	T	Benzo(g,h,i)perylene	1.298	1.314	-1.2	80 0.00
74	S	Perylene-d12	0.823	0.899	-9.2	88 0.00
75	$^{\mathrm{T}}$	Perylene	1.640	1.813	-10.5	88 0.00

Data File : G:\1\DATA\MS30377\MS30377I.D Vial: 4 Acq On : 13 Apr 2007 5:40 pm Sample : AR-WKCC-250-022 Operator: TJM

Inst : GC/MS Ins Multiplr: 1.00

Misc

MS Integration Params: rteint.p

Quant Results File: 041507.RES Quant Time: Apr 14 20:12 19107

Quant Method : C:\MS30377\041507.M (RTE Integrator)

: PAH Calibration Table (2002) Title Last Update : Sat Apr 14 20:04:44 2007

Response via : Initial Calibration

DataAcq Meth: PAH-2002

Internal Standards	R.T.	QIon	Response	Conc Units Dev	(Min)
 Fluorene-d10 Pyrene-d10 Benzo(a)pyrene-d12 				51.08 ng/ml	0.00
29) Pyrene-d10	28.80	212	3557m	49.98	0.00
60) Benzo(a)pyrene-d12	37.42	264	2109m	45.61	0.00
System Monitoring Compounds					
2) Naphthalene-d8	13.01	136	19435	261.04	0.00
20) Acenaphthene-d10	18.86	164	9425m	262.70 262.70 296.76	0.00
30) Phenanthrene-d10	23.91	188	15868	262.70	0.00
53) Chrysene-d12	33.00	240	24231m	296.76	0.00
74) Perylene-d12	37.70	264	10388m	273.02	0.00
Target Compounds					zalue
3) Decalin	10.39		4120m	261.96 ng/ml	
4) C1-Decalin	0.00	152	0		
5) C2-Decalin	0.00	166	0		
6) C3-Decalin	0.00	180	0	N.D. d	
7) C4-Decalin	0.00		0		
8) Naphthalene	13.09	194 128 142 142 156 170	21243m	261.25	
9) 2-Methylnaphthalene	15.34	142	14147m	266.08	
10) 1-Methylnaphthalene	15.65	142	12869m	261.60	
11) 2,6-Dimethylnaphthalene	17.45	156	11233m	259.49	
12) 1,6,7-Trimethylnaphthalene	20.30	170	10762m		
13) C2-Naphthalenes 14) C3-Naphthalenes 15) C4-Naphthalenes 16) Benzothiophene	0.00	156	U		
14) C3-Naphthalenes	0.00	170			
15) C4-Naphthalenes	0.00	184	0		
16) Benzothiophene	13.24	134	18259m	-	
i'' Cimbelladelladplicite	0.00		0		
18) C2-Benzothiophene	0.00	162	0	N.D. d	
19) C3-Benzothiophene	0.00	176		N.D. d	
21) Biphenyl	16.92	154	17099m	263.21	
22) Acenaphthylene	18.35		20718m	264.32	
23) Acenaphthene	18.95	154	11963m	269.06	
24) Dibenzofuran	19.56	168	18746m	260.16 ng/ml	
25) Fluorene	20.75	700	13462III		
26) C1-Fluorenes	0.00	180	0	N.D. d	
27) C2-Fluorenes	0.00	194	0	N.D. d	
28) C3-Fluorenes	0.00	208	0	N.D. d	
31) Pentachlorophenol	0.00	266	0	N.D. d	
32) Carbazole	24.75		18033	256.48 ng/ml	
33) Dibenzothiophene	23.57		20573m 0	274.94 N.D. d	
34) C1-Dibenzothiophene	0.00	198 	·	N.D. U	

^{(#) =} qualifier out of range (m) = manual integration Mon Apr 16 08:07:51 2007 MS30377I.D 041507.M

Vial: 4 Data File : G:\1\DATA\MS30377\MS30377I.D Acq On : 13 Apr 2007 5:40 pm Sample : AR-WKCC-250-022 Operator: TJM

Inst : GC/MS Ins

Multiplr: 1.00 Misc

MS Integration Params: rteint.p Quant Results File: 041507.RES Ouant Time: Apr 14 20:12 19107

Quant Method : C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002) Last Update : Sat Apr 14 20:04:44 2007

Response via : Initial Calibration

DataAcq Meth: PAH-2002

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35)	C2-Dibenzothiophene	0.00	212	0	N.D. d	
36)	C3-Dibenzothiophene	0.00	226	0	N.D. d	
37)	C3-Dibenzothiophene Phenanthrene	24.01 24.18 26.13	178	20855m	291.28	
38)	Anthracene	24.18	178	23365m	294.16	
39)	Anthracene 1-Methylphenanthrene C1-Phenanthrene/Anthracene	26.13	192	18266m	298.93	
40)	C1-Phenanthrene/Anthracene	0.00	192	0	N.D. d	
	C2-Phenanthrene/Anthracene	0.00	206	0	N.D. d	
	C3-Phenanthrene/Anthracene	0.00	220	0	N.D. d	
43)	C4-Phenanthrene/Anthracene	0.00	234	0	N.D. d	
	Naphthobenzothiophene	32.16	234	20279m	289.73	
	Cl-Naphthobenzothiophene	0.00	248		N.D. d	
	C2-Naphthobenzothiophene	0.00	262	0	N.D. d	
	C3-Naphthobenzothiophene	0.00	276	0	N.D. d	
	Fluoranthene	28.09	202	23392m	253.09	
	Pyrene	28.86	202	28295m	278.85	
	C1-Fluoranthenes/Pyrenes	0.00	216 230	0	N.D. d	
51)	C2-Fluoranthenes/Pyrenes	0.00	230	0	N.D. d	
	C3-Fluoranthenes/Pyrenes	0.00	244	0	N.D. d	
	Benz(a)anthracene	32.96	228	26259m	278.75	
	Chrysene	33.07		20251m	258.19	
56)	C1-Chrysenes	0.00 0.00 0.00	242	0	N.D. d	
57)	C2-Chrysenes C3-Chrysenes C4-Chrysenes C29-Hopane	0.00	256	0	N.D. d	
58)	C3-Chrysenes	0.00	270			
59)	C4 Characana	0 00	201	0	N.D. d	
61)	C29-Hopane 18a-Oleanane C30-Hopane Benzo(b) fluoranthene Benzo(c) pyrene Benzo(a) pyrene	0.00	191	0		
62)	18a-Oleanane	0.00	191	0	N.D. d	
63)	C30-Hopane	41.74	191	8384m	267.12 ng/1	ml
64)	Benzo(b) fluoranthene	36.43	252	23691m	286.68	
65)	Benzo(k)fluoranthene	36.50	252	20613m	257.79	
66)	Benzo(e)pyrene	37.31	252	23486m	282.17	
67)	Benzo(a) pyrene	37.49	252	18715m	264.74	
68)	Indeno(1,2,3-c,d)pyrene	41.84	276	13460m	250.71	
	Dibenzo(a,h)anthracene	41.97	278	12259m	246.24	
70)	C1-Dibenzo(a,h)anthracene	0.00	292	0	N.D. d	
	C2-Dibenzo(a,h)anthracene	0.00	306	0		
	C3-Dibenzo(a,h)anthracene	0.00	320	0	N.D. d	
	Benzo(g,h,i)perylene	43.06	276	15221m	253.62	
	Perylene	37.77	252	20996m	276.93	

^{(#) =} qualifier out of range (m) = manual integration MS30377I.D 041507.M Mon Apr 16 08:07:51 2007

```
Quant Results File: 041507.RES
                           GC/MS
                                        1.00
              E
E
              Operator:
                                       Multiplr:
 Vial:
                           Inst
G:\1\DATA\MS30377\MS30377I.D
             5:40 pm
                                                    MS Integration Params: rteint.p
Quant Time: Apr 14 20:12 19107
            13 Apr 2007 5
AR-WKCC-250-022
Data File
             Acq On
                          Sample
                                        Misc
```

(RTE Integrator) G:\1\METHODS\041507.M
PAH Calibration Table Method

(2002)Sat Apr 14 20:04:44 2007 Initial Calibration Last Update Title

54.00 52.00 50.00 48.00 46.00 44,00 Benzo(g,h,i)perylene, T 45.00 40.00 38.00 Benzo(a)pyrene Herylene, T penzo(a)pyrene-crz, r 36,00 Benzo(b)fluoranthene, T TIC: MS30377I.D 34.00 T,enecentitàşttithracene, T 32.00 Naphthobenzothiophene, T 30,00 Pyrene, T 1,016-enery9 28,00 Fluoranthene, T 26.00 1-Methylphenanthrene, T Carbazole, T 24.00 Dibenzothiophene, T Anthracene, T — Phenanthrene, T Phenanthrene-d10, S 22.00 20.00 Fluorene-d10,1 1,6,7-Trimethyinaphthalene, T Dibenzofuran, T Acenaphthene, T 18.00 Acenaphthylene, T T,enelarthqsnlynaene, T Biphenyl, T 16.00 T ,enelentrylnaphthalene, T ,enelentrylnaphthalene, T 14.00 Response via T , Shell Hatter Handle 18, 17 10,00 12,00 Decalin, T Abundance 22000-2000 20000 18000 16000 14000 12000 10000 8000 9009 4000 00004

S

Page

08:07:53 2007

Mon Apr 16

041507.M

MS30377I.D

Evaluate Continuing Calibration Report

Data File : G:\1\DATA\MS30377\MS30377J.D Vial: 6 Acq On : 13 Apr 2007 7:46 pm Sample : AR-WKCC-250-022 Operator: TJM

Inst : GC/MS Ins

Misc Multiplr: 1.00

MS Integration Params: rteint.p

Method : G:\1\METHODS\041507.M (RTE Integrator)
Title : PAH Calibration Table (2002)
Last Update : Sat Apr 14 20:04:44 2007
Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 25% Max. R.T. Dev 0.50min Max. RRF Dev : 25% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev Area% Dev(min)
1	I	Fluorene-d10	1.000	1.000	0.0 110 0.00
2	S	Naphthalene-d8	1.911	1.861	2.6 109 0.00
3	\mathbf{T}	Decalin	0.404	0.385	4.7 110 0.00
4	un	C1-Decalin	0.404	0.000	100.0# 0# -12.76#
5	un	C2-Decalin	0.404	0.000	100.0# 0# -14.27#
6	un	C3-Decalin	0.404	0.000	100.0# 0# -15.86#
7	un	C4-Decalin	0.404	0.000	100.0# 0# -19.68#
8	\mathbf{T}	Naphthalene	2.087	2.054	1.6 110 0.00
9	${f T}$	2-Methylnaphthalene	1.365	1.379	-1.0 114 0.00
10	T	1-Methylnaphthalene	1.263	1.220	3.4 115 0.00
11		2,6-Dimethylnaphthalene	1.111	1.119	-0.7 114 0.00
12	\mathbf{T}	1,6,7-Trimethylnaphthalene	1.026	1.058	-3.1 134 0.00
13	un	C2-Naphthalenes	2.087	0.000	100.0# 0# -17.94#
14	un	C3-Naphthalenes	2.087	0.000	100.0# 0# -20.12#
15	un	C4-Naphthalenes	2.087	0.000	100.0# 0# -22.24#
16	\mathbf{T}	Benzothiophene	1.767	1.701	3.7 106 0.00
17	un	C1-Benzothiophene	1.767	0.000	100.0# 0# -15.58#
18	un	C2-Benzothiophene	1.767	0.000	100.0# 0# -17.97#
19	un	C3-Benzothiophene	1.767	0.000	100.0# 0# -19.61#
20	S	Acenaphthene-d10	0.921	0.924	-0.3 115 0.00
21		Biphenyl	1.667	1.704	-2.2 114 0.00
22		Acenaphthylene	2.012	2.046	-1.7 113 0.00
23	${f T}$	Acenaphthene	1.141	1.126	1.3 109 0.00
24		Dibenzofuran	1.850	1.825	1.4 112 0.00
25	T	Fluorene	1.328	1.344	-1.2 113 0.00
	un	C1-Fluorenes	1.328	0.000	100.0# 0# -22.65#
27	un	C2-Fluorenes	1.328	0.000	100.0# 0# -24.32#
28	un	C3-Fluorenes	1.328	0.000	100.0# 0# -26.19#
29	I	Pyrene-d10	1.000	1.000	0.0 107 0.00
30	S	Phenanthrene-d10	0.849	0.796	6.2 102 0.00
31		Pentachlorophenol	0.044	0.000	100.0# 0# -23.30#
32	${f T}$	Carbazole	0.988	1.023	-3.5 113 0.00
33	T	Dibenzothiophene	1.051	1.077	-2.5 113 0.00
	un	C1-Dibenzothiophene	1.051	0.000	100.0# 0# -25.11#
	un	C2-Dibenzothiophene	1.051	0.000	100.0# 0# -26.47#
	un	C3-Dibenzothiophene	1.051	0.000	100.0# 0# -28.57#
37		Phenanthrene	1.006		-0.9 115 0.00
38		Anthracene	1.116		-3.2 114 0.00
39		1-Methylphenanthrene	0.859		
40	un	C1-Phenanthrene/Anthracene	1.006	0.000	100.0# 0# -26.20#

^{(#) =} Out of Range

Evaluate Continuing Calibration Report

Vial: 6 Data File : G:\1\DATA\MS30377\MS30377J.D Acq On : 13 Apr 2007 7:46 pm Operator: TJM

Inst : GC/MS Ins Sample : AR-WKCC-250-022

Multiplr: 1.00 Misc

MS Integration Params: rteint.p

: G:\1\METHODS\041507.M (RTE Integrator)

Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 25% Max. R.T. Dev 0.50min Max. RRF Dev : 25% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev Area% Dev(min)
41 un	C2-Phenanthrene/Anthracene	1.006	0.000	100.0# 0# -27.51#
42 un	C3-Phenanthrene/Anthracene	1.006	0.000	100.0# 0# -30.06#
43 un	C4-Phenanthrene/Anthracene	1.006	0.000	100.0# 0# -30.78#
44 T	Naphthobenzothiophene	0.983	1.025	-4.3 119 0.00
45 un	C1-Naphthobenzothiophene	0.983	0.000	100.0# 0# -33.03#
46 un	C2-Naphthobenzothiophene	0.983	0.000	100.0# 0# -34.70#
47 un	C3-Naphthobenzothiophene	0.983	0.000	100.0# 0# -36.06#
48 T	Fluoranthene	1.299	1.236	4.8 103 0.00
49 T	Pyrene	1.426	1.429	-0.2 111 0.00
50 un	C1-Fluoranthenes/Pyrenes	1.299	0.000	100.0# 0# -30.52#
51 un	C2-Fluoranthenes/Pyrenes	1.299	0.000	100.0# 0# -31.88#
52 un	C3-Fluoranthenes/Pyrenes	1.299	0.000	100.0# 0# -32.93#
53 S	Chrysene-d12	1.147	1.234	-7.6 121 0.00
54 T	Benz(a)anthracene	1.324	1.423	-7.5 117 0.00
55 T	Chrysene	1.102	1.142	-3.6 110 0.00
56 un	C1-Chrysenes	1.102	0.000	100.0# 0# -34.25#
57 un	C2-Chrysenes	1.102	0.000	100.0# 0# -35.44#
58 un	C3-Chrysenes	1.102	0.000	100.0# 0# -36.87#
59 un	C4-Chrysenes	1.102	0.000	100.0# 0# -42.39#
60 I	Benzo(a)pyrene-d12	1.000	1.000	0.0 133 0.00
61 un	C29-Hopane	0.679	0.000	100.0# 0# -40.43#
62 un	18a-Oleanane	0.679	0.000	100.0# 0# -42.28#
63 T	C30-Hopane	0.679	0.636	6.3 132 0.00
64 T	Benzo(b) fluoranthene	1.787	1.690	5.4 131 0.00
65 T	Benzo(k)fluoranthene	1.729	1.646	4.8 125 0.00
66 T	Benzo(e)pyrene	1.800	1.581	12.2 116 0.00
67 T	Benzo(a)pyrene	1.529	1.475	3.5 131 0.00
68 T	Indeno(1,2,3-c,d)pyrene	1.161	1.228	-5.8 151 0.00
69 T	Dibenzo(a,h)anthracene	1.077	1.158	-7.5 150 0.00
70 un	C1-Dibenzo(a,h)anthracene	1.077	0.000	100.0# 0# -42.56#
71 un	C2-Dibenzo(a,h)anthracene	1.077	0.000	100.0# 0# -44.47#
72 un	C3-Dibenzo(a,h)anthracene	1.077	0.000	100.0# 0# -44.94#
73 T	Benzo(g,h,i)perylene	1.298	1.251	3.6 131 0.00
74 S	Perylene-d12	0.823	0.758	7.9 129 0.00
75 T	Perylene	1.640	1.537	6.3 128 0.00

Data File : G:\1\DATA\MS30377\MS30377J.D Vial: 6 Acq On : 13 Apr 2007 7:46 pm Sample : AR-WKCC-250-022 Operator: TJM

Inst : GC/MS Ins

Multiplr: 1.00 Misc

MS Integration Params: rteint.p Quant Time: Apr 14 20:20 19107

Quant Results File: 041507.RES

Quant Method : C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002)
Last Update : Sat Apr 14 20:04:44 2007
Response via : Initial Calibration

DataAcq Meth : PAH-2002

Internal Standards	R.T.	QIon	Response	Conc Units De	v(Min)
1) Fluorene-d10 29) Pyrene-d10	20.63	176	2671m		0.00
29) Pyrene-d10	28.80	212	5049m		0.00
60) Benzo(a)pyrene-d12	37.42	264	3645m	45.61	0.00
System Monitoring Compounds					
2) Naphthalene-d8	13.01	136	24333m	243.50	0.00
20) Acenaphthene-d10	18.86			250.73	0.00
	23.91		20094m		0.00
53) Chrysene-d12	33.00	240	31174m	268.97	0.00
74) Perylene-d12	37.70	264	15138		0.00
Target Compounds				0	value
3) Decalin	10.39	138	5036m		
	0.00		0	N.D. d	
5) C2-Decalin	0.00	166	0	N.D. d	
6) C3-Decalin	0.00	180	0	N.D. d	
7) C4-Decalin	0.00	194	0	N.D. d	
8) Naphthalene	0.00 0.00 0.00 13.09 15.34	128		246.52	
9) 2-Methylnaphthalene	15.34	142		253.33	
10) 1-Methylnaphthalene	15.65			241.87	
11) 2,6-Dimethylnaphthalene	17.45			252.32	
12) 1,6,7-Trimethylnaphthalene			13856m		
13) C2-Naphthalenes	0.00	156	0	N.D. d	
14) C3-Naphthalenes	0.00	170	0	N.D. d	
15) C4-Naphthalenes		184	0	N.D. d	
16) Benzothiophene	13.23	134	22282m	241.18 ng/ml	
17) C1-Benzothiophene	0.00	148	0	N.D. d	
18) C2-Benzothiophene	0.00	162	0	N.D. d	
19) C3-Benzothiophene	0.00	176	0	N.D. d	
21) Biphenyl	16.92	154	22314m	255.91	
22) Acenaphthylene	18.35	152	26802m	254.76	
23) Acenaphthene	18.94	154	14732m	246.86	
24) Dibenzofuran	19.56	168	23887m	246.98 ng/ml	
25) Fluorene	20.74	166	17602m	253.44	
26) C1-Fluorenes	0.00	180	0	N.D. d	
27) C2-Fluorenes	0.00	194	0	N.D. d	
28) C3-Fluorenes	0.00	208	0	N.D. d	
31) Pentachlorophenol	0.00	266	0	N.D.	
32) Carbazole	24.75	167	25881m	259.33 ng/ml	
33) Dibenzothiophene	23.57	184	27251m	256.56	
34) C1-Dibenzothiophene	0.00	198	0	N.D. d	

Data File : G:\1\DATA\MS30377\MS30377J.D Vial: 6 Acq On : 13 Apr 2007 7:46 pm Sample : AR-WKCC-250-022 Operator: TJM

Inst : GC/MS Ins

Misc Multiplr: 1.00

MS Integration Params: rteint.p

Quant Time: Apr 14 20:20 19107 Quant Results File: 041507.RES

Quant Method: C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002) Last Update : Sat Apr 14 20:04:44 2007

Response via : Initial Calibration

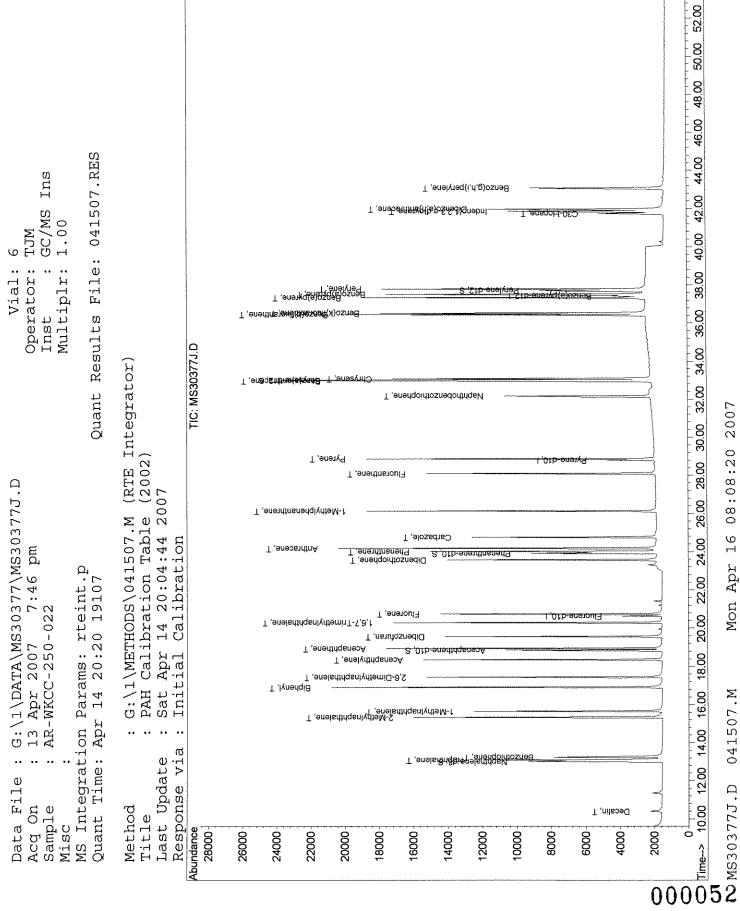
DataAcq Meth : PAH-2002

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35)	C2-Dibenzothiophene	0.00	212	0	N.D. d	
36)	C3-Dibenzothiophene Phenanthrene Anthracene 1-Methylphenanthrene	0.00	226	0	N.D. d	
37)	Phenanthrene	24.01 24.18 26.13	178	25687m	252.75	
38)	Anthracene	24.18	178	29138m	258.44	
39)	1-Methylphenanthrene	26.13	192	22662m	261.28	
40)	C1-Phenanthrene/Anthracene	0.00	192	0	N.D. d	
41)	C2-Phenanthrene/Anthracene	0.00	206	0	N.D. d	
42)	C3-Phenanthrene/Anthracene	0.00	220	0	N.D. d	
43)	C4-Phenanthrene/Anthracene		234	0	N.D. d	
	Naphthobenzothiophene	32.16	234	25862m	260.30	
45)	C1-Naphthobenzothiophene	0.00	248	0	N.D. d	
	C2-Naphthobenzothiophene	0.00	262	0	N.D. d	
	C3-Naphthobenzothiophene	0.00	276		N.D. d	
	Fluoranthene	28.09	202	31298m	238.56	
49)	Pyrene	28.86	202	36166m	251.09	
	C1-Fluoranthenes/Pyrenes	0.00	216	0	N.D. d	
	C2-Fluoranthenes/Pyrenes	0.00	216 230	0	N.D. d	
	C3-Fluoranthenes/Pyrenes	0.00	244	0		
54)	Benz(a)anthracene	32.96	228	36029m	269.44	
55)	Chrysene	33.07	228	28904m	259.62	
56)	C1-Chrysenes	0.00	242	0	N.D. d	
57)	C2-Chrysenes	0.00	256	0	N.D. d	
58)	C3-Chrysenes	0.00	270		N.D. d	
59)	C4-Chrysenes	0.00	284	0	N.D. d	
61)	C29-Hopane	0.00	191	0	N.D. d	
62)	18a-Oleanane	0.00	191	0	N.D. d	
63)	Chrysene C1-Chrysenes C2-Chrysenes C3-Chrysenes C4-Chrysenes C4-Chrysenes C29-Hopane 18a-Oleanane C30-Hopane Benzo(b) fluoranthene Benzo(k) fluoranthene Benzo(e) pyrene Benzo(a) pyrene Indeno(1,2,3-c,d) pyrene	41.74	191	12698	234.08 ng/n	nl
64)	Benzo(b) fluoranthene	36.43	252	12698 33806	236.69	
65)	Benzo(k)fluoranthene	36.50	252	32964	238.53	
66)	Benzo(e)pyrene	37.31	252	31675	220.19	
67)	Benzo(a)pyrene	37.49	252	29509		
68)	Indeno(1,2,3-c,d)pyrene	41.84	276 278	24584	264.95	
69)	Dibenzo(a,h)anthracene	41.95	278	23163	269.20	
		0.00	292	0	N.D. d	
	C2-Dibenzo(a,h)anthracene	0.00	306	0		
	C3-Dibenzo(a,h)anthracene		320		N.D. d	
	Benzo(g,h,i)perylene	43.06			241.41	
	Perylene	37.77		30764		

041507.M

54.0

Quantitation Report



Data File Name MS30377B.D

Data File Path X:11/DATAMMS30377\
Operator TJM

Date Acquired 04/13/20 -1:0:
Method File PAH-2002
Sample Name SRM 1582
Misc Info
Instrument Name GC/MS Ins
Vial Number 2
Sample Multiplier 0.588
Sample Amount 0

Su Amt = 50

MS30377B.D SRM 1582

04/13/20 -1:0: PAH-2002 1.70

Peak#	Compound	Ret Time (min)	Target Response (Area)	Conc. (ng/g or ng/L)	Su. Corrected Conc. (ng/g or ng/L)
3)	Decalin	0.00	0	0.00	0.00
4)	C1-Decalin	0.00	0	0.00	0.00
5)	C2-Decalin	0,00	0	0.00	0.00
6)	C3-Decalin	0.00	0	0.00	0.00
7)	C4-Decalin	0.00	0	0.00	0,00
8)	Naphthalene	13.06	24820	144.66	148.41
9+10) 13)	C1-Naphthalenes	15,49 17,79	96757 178444	563.94 1040.04	578.53 1066.96
14)	C2-Naphthalenes C3-Naphthalenes	19.73	170794	995,45	1021.22
15)	C4-Naphthalenes	22.29	121460	707.91	726.24
16)	Benzothiophene	13.32	1168	8.04	8,25
17)	C1-Benzothiophene	15.34	2484	17.10	17,55
18)	C2-Benzothiophene	17.82	10652	73.34	75.24
19)	C3-Benzothiophene	19.53	20840	143.49	147.20
21)	Biphenyl	16.92	4542	33.14	33.99
22)	Acenaphthylene	0.00	Đ	0.00	0,00
23)	Acenaphthene	18.94	1988	21.19	21.74
24)	Dibenzofuran	19.56	1834	12.06	12.37
25)	Fluorene	20.71	3836	35.13	36.04
26) 27)	C1-Fluorenes C2-Fluorenes	22.71 24.45	13747 28851	125.91	129.17
28)	C3-Fluorenes	26.10	26274	264.25 240.64	271.09 246.87
31)	Pentachlorophenoi	0.00	20274	0.00	0.00
32)	Carbazole	24.82	402	2.14	2.19
38)	Anthracene	24.18	778	3.66	3.76
37)	Phenanthrene	23.98	18683	97.58	100.11
40)	C1-Phenanthrene/Anthracene	25.76	63979	334.16	342.81
41)	C2-Phenanthrene/Anthracene	27.58	96372	503.35	516.38
42)	C3-Phenanthrene/Anthracene	29.17	101432	529.78	543.49
43)	C4-Phenanthrene/Anthracene	31.02	52306	273.19	280.26
33)	Dibenzothiophene	23.57	5987	29.92	30.69
34)	C1-Dibenzothiophene	25.09	25556	127.72	131.02
35)	C2-Dibenzothiophene	26,81	48508	242.42	248.69
36)	C3-Dibenzothiophene	28.02	49064	245.20	251.54
48)	Fluoranthene	28.09	1255	5.08	5.21
49) 50)	Pyrene	28.86	1631	6.01	6.17
51)	C1-Fluoranthenes/Pyrenes C2-Fluoranthenes/Pyrenes	30.68 31.52	14626 24872	59.18 100.63	60.71 103.24
52)	C3-Fluoranthenes/Pyrenes	33.21	18609	75.29	77.24
44)	Naphthobenzothiophene	32.16	6353	33.94	34.82
45)	C1-Naphthobenzothiophene	33.88	11046	59.01	60.54
46)	C2-Naphthobenzothiophene	35.26	14084	75,25	77.19
47)	C3-Naphthobenzothiophene	36.36	10079	53.85	55.24
54)	Benz(a)anthracene	32.96	1086	4.31	4.42
55)	Chrysene	33.07	4852	23.13	23.73
56)	C1-Chrysenes	34.31	14048	66.98	68.71
57)	C2-Chrysenes	36.64	25266	120.46	123.58
58)	C3-Chrysenes	37.13	17614	83.98	86.15
59)	C4-Chrysenes	0.00	0	0.00	0.00
64) 65)	Benzo(b)fluoranthene Benzo(k)fluoranthene	36,43 36,46	420 117	1.62 0.47	1.66 0.48
66)	Benzo(e)pyrene	37,31	740	2.84	2.91
67)	Benzo(a)pyrene	37.49	437	1.97	2.02
75)	Perylene	37.77	8706	36,66	37.61
68)	Indeno(1,2,3-c,d)pyrene	41.82	331	1.97	2.02
69)	Dibenzo(a,h)anthracene	41.92	84	0.54	0.55
70)	C1-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00
71)	C2-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00
72)	C3-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00
73)	Benzo(g,h,i)perylene	43.06	299	1.59	1.63
	Total PAH				7828
	Individual Isomers				
9)	2-Methylnaphthalene	15.34	57656	513.93	527.23
10)	1-Methylnaphthalene	15.65	39101	376.69	386.45
11)	2.6-Dimethylnaphthalene	17.48	48527	531.28	545.04
12)	1,6,7-Trimethylnaphthalene	20.29	13848	164.17	168.42
39)	1-Methylphenanthrene	26.13	14011	85.75	87.97
61)	C29-Hopane	39.79	18934	192.60	197.59
62)	18a-Oleanane	40.83	6183	62.90	64.52
63)	C30-Hopane	41.04	28305	287.93	295.38
	Surrogates (AR-STSU-040-005)				Su Recovery (%)
2)	(AK-S (SU-040-005) Naphthalene-d8	13.01	4339	27.62	94
20)	Acenaphthene-d10	18.83	2201	29.07	99
30)	Phenanthrene-d10	23.91	4629	28.66	97
53)	Chrysene-d12	33.00	5768	26.42	90
74)	Perylene-d12 Internal Stds	37.70	3203	26.88	91
45	(AR-WKIS-0500-007)		***		
1)	Fluorene-d10	20.63	2469	51.08	
29)	Pyrene-d10	28.80	5593	49.98	
60)	Benzo(a)pyrene-d12	37.42	3884	45.61	

Data File : X:\1\DATA\MS30377\MS30377B.D
Acq On : 13 Apr 2007 10:19 am
Sample : SRM 1582
Misc : Vial: 2 Operator: TJM

Inst : GC/MS Ins Multiplr: 0.59

MS Integration Params: rteint.p

Quant Time: Apr 16 7:58 2007 Quant Results File: 041507.RES

Quant Method : C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002)
Last Update : Sat Apr 14 20:04:44 2007
Response via : Initial Calibration
DataAcq Meth : PAH-2002

Internal Standards	R.T.	QIon	Response	Conc Units Dev	(Min)
1) Fluorene-d10	20.63	176	2469m	51.08 ng/ml	0.00
29) Pyrene-d10	28.80	212	5593m	49.98	0.00
60) Benzo(a)pyrene-d12	37.42	264	3884m	45.61	0.00
System Monitoring Compounds	12 01	136	4339m	27.62	0.00
2) Naphthalene-d8	13.01 18.83		2201m		-0.03
20) Acenaphthene-d10	23.91			28.66	0.00
30) Phenanthrene-d10 53) Chrysene-d12	33.00			26.42	0.00
74) Perylene-d12	37.70		3203		0.00
(4) iciyiche diz	J, • , U	201	J J	20.00	
Target Compounds					alue
3) Decalin	0.00		0	N.D.	
4) C1-Decalin	0.00		0	N.D. d	
5) C2-Decalin	0.00		0	N.D. d	
6) C3-Decalin	0.00		0	N.D. d	
7) C4-Decalin	0.00	194	0	N.D. d	
	13.06		24820	144.66	
9) 2-Methylnaphthalene	15.34		57656	513.93	
10) 1-Methylnaphthalene	15.65		39101 48527	376.69	
11) 2,6-Dimethylnaphthalene	17.48				
12) 1,6,7-Trimethylnaphthalene			13848	164.17	
13) C2-Naphthalenes	17.79 19.73		178444 170794	1040.04	
14) C3-Naphthalenes			170794	707 01	
15) C4-Naphthalenes	22.29 13.32		1160	707.91 8.04 ng/ml 17.10 ng/ml	
16) Benzothiophene	15.34		2/8/	17 10 ng/mi	
17) C1-Benzothiophene18) C2-Benzothiophene	17.82		10652	73.34 ng/ml	
19) C3-Benzothiophene	19.53		20840	143.49 ng/ml	
21) Biphenyl	16.92		4542m		
22) Acenaphthylene	0.00		0		
23) Acenaphthene	18.94				
24) Dibenzofuran	19.56		1834	21.19 12.06 ng/ml 35.13	
25) Fluorene	20.71		3836	35.13	
26) C1-Fluorenes	22.71		13747	125.91	
27) C2-Fluorenes	24.45		28851m		
28) C3-Fluorenes	26.10		26274m		
	0.00			N.D.	
32) Carbazole	24.82		402m	2.14 ng/ml	
33) Dibenzothiophene	23.57		5987m	2.14 ng/ml 29.92	
34) Cl-Dibenzothiophene	25.09	198	25556m		
35) C2-Dibenzothiophene	26.81	212	48508m	242.42	
36) C3-Dibenzothiophene	28.02	226	49064m	245.20	
37) Phenanthrene	23.98	178	18683m	97.58	
38) Anthracene	24.18	178	778m	3.66	
39) 1-Methylphenanthrene	26.13		14011m	85.75	
40) C1-Phenanthrene/Anthracene	25.76	192	63979m	334.16	
41) C2-Phenanthrene/Anthracene	27.58		96372m	503.35	
42) C3-Phenanthrene/Anthracene	29.17		101432m	529.78	
43) C4-Phenanthrene/Anthracene	31.02	234	52306m	273.19	
44) Naphthobenzothiophene	32.16	234	6353m	33.94	
45) C1-Naphthobenzothiophene	33.88	248	11046m	59.01	
46) C2-Naphthobenzothiophene	35.26	262	14084m	75.25	

Data File : X:\1\DATA\MS30377\MS30377B.D

Vial: 2 Acq On : 13 Apr 2007 10:19 am Sample : SRM 1582 Misc : Operator: TJM

Inst : GC/MS Ins

Multiplr: 0.59

MS Integration Params: rteint.p Quant Time: Apr 16 7:58 2007

Quant Results File: 041507.RES

Quant Method: C:\MS30377\041507.M (RTE Integrator)
Title: PAH Calibration Table (2002)
Last Update: Sat Apr 14 20:04:44 2007
Response via: Initial Calibration
DataAcq Meth: PAH-2002

	Compound	R.T.	QIon	Response	Conc Unit Qvalue	
47)	C3-Naphthobenzothiophene	36.36	276	10079m	53.85	
48)		28.09	202	1255m	5.08	
49)		28.86	202	1631m	6.01	
50)	C1-Fluoranthenes/Pyrenes	30.68	216	14626m	59.18 ng/mL	
51)	C2-Fluoranthenes/Pyrenes	31.52	230	24872m		
52)	C3-Fluoranthenes/Pyrenes	33.21	244	18609m		
	Benz(a)anthracene	32.96	228	1086m	4.31	
55)	Chrysene	33.07	228	4852m	23.13	
56)	C1-Chrysenes	34.31	242	14048m	66.98 ng/mL	
57)	C2-Chrysenes	36.64	256	25266m	120.46 ng/mL	
58)	C3-Chrysenes	37.13	270	17614m	83.98 ng/mL	
59)	C4-Chrysenes	0.00	284		N.D. d	
61)	C29-Hopane	39.79	191		192.60 ng/ml	
62)	18a-Oleanane	40.83	191	6183	62.90 ng/ml	
63)	C30-Hopane	41.04	191	28305	287.93 ng/ml	
	Benzo(b)fluoranthene	36.43	252	420	1.62	
65)	Benzo(k)fluoranthene	36.46	252	117	0.47	
66)	Benzo(e)pyrene	37.31	252	740	2.84	
67)	Benzo(a)pyrene	37.49	252	437	1.97	
68)	Indeno(1,2,3-c,d)pyrene	41.82	276	331	1.97	
69)	Dibenzo(a,h)anthracene	41.92	278	84	0.54	
70)	C1-Dibenzo(a,h)anthracene	0.00	292	0	N.D. d	
71)	C2-Dibenzo(a,h)anthracene	0.00	306	0	N.D. d	
	C3-Dibenzo(a,h)anthracene	0.00	320	0	N.D. d	
73)		43.06	276	299	1.59	
75)	Perylene	37.77	252	8706m	36.66	

2007

06:52:08

08

May

Tue

041507

MS30377B.D

54.00

52.00

Tissue, Sediment, and Water Sample Report (Use d-10 Phenanthrene only for Surrogate Corrections)

Data File Name MS30377C.D
Data File Path X:\\1\DATA\MS30377\.
Operator TJM
Date Acquired 04/13/20 -1:1:
Method File PAH-2002
Sample Name IS/SU Mixture
Misc Info
Instrument Name GC/MS Ins
Vial Number 3
Sample Multiplier 1
Sample Amount 0

Su Amt = 50

MS30377C.D IS/SU Mixture

04/13/20 -1:1: PAH-2002 1.00

Peak#	Compound	Ret Time (min)	Target Response (Area)	Conc. (ng/g or ng/L)	Su. Corrected Conc. (ng/g or ng/L)		
3)	Decalin	0.00	0	0.00	0.00		
4)	C1-Decalin	0.00	0	0,00	0.00		
5)	C2-Decalin	0.00	0	0.00	0.00		
6)	C3-Decalin	0.00	0	0.00	0.00		
7)	C4-Decalin	0.00	0	0.00	0,00		
8)	Naphthalene	0.00	0	0.00	0.00		
9+10)	C1-Naphthalenes	0.00	0	#DIV/0I	#DIV/01		
13) 14)	C2-Naphthalenes C3-Naphthalenes	0.00	0	0.00 0.00	0.00		
15)	C4-Naphthalenes	0.00	0	0.00	0.00		
16)	Benzothiophene	0.00	0	0.00	0.00		
17)	C1-Benzothiophene	0.00	ő	0.00	0.00		
18)	C2-Benzothiophene	0.00	ō	0.00	0.00		
19)	C3-Benzothiophene	0.00	0	0.00	0.00		
21)	Biphenyl	0.00	0	0.00	0.00		
22)	Acenaphthylene	0.00	0	0.00	0.00		
23)	Acenaphthene	0.00	0	0.00	0.00		
24)	Dibenzofuran	0.00	0	0.00	0.00		
25)	Fluorene	0.00	0	0.00	0.00		
26) 27)	C1-Fluorenes	0.00	0	0.00	0.00		
28)	C2-Fluorenes C3-Fluorenes	0.00	0	0.00	0.00 0.00		
31)	Pentachiorophenol	0.00	0	0.00	0.00		
32)	Carbazole	0.00	0	0.00	0.00		
38)	Anthracene	0.00	ō	0.00	0.00		
37)	Phenanthrene	0.00	0	0.00	0.00		
40)	C1-Phenanthrene/Anthracene	0.00	0	0.00	0.00		
41)	C2-Phenanthrene/Anthracene	0.00	0	0.00	0.00		
42)	C3-Phenanthrene/Anthracene	0.00	0	0.00	0,00		
43)	C4-Phenanthrene/Anthracene	0,00	0	0,00	0.00		
33)	Dibenzothiophene	0.00	0	0.00	0.00		
34)	C1-Dibenzothiophene	0.00	0	0.00	0,00		
35)	C2-Dibenzothiophene	0.00	0	0.00	0.00		
36) 48)	C3-Dibenzothiophene Fluoranthene	0.00	0	0.00	0,00		
49)	Pyrene	0.00	0	0.00 0.00	0.00 0.00		
50)	C1-Fluoranthenes/Pyrenes	0.00	0	0.00	0.00		
51)	C2-Fluoranthenes/Pyrenes	0.00	ő	0.00	0.00		
52)	C3-Fluoranthenes/Pyrenes	0.00	0	0.00	0.00		
44)	Naphthobenzothiophene	0.00	0	0.00	0.00		
45)	C1-Naphthobenzothiophene	0.00	0	0.00	0.00		
46)	C2-Naphthobenzothiophene	0.00	0	0.00	0.00		
47)	C3-Naphthobenzothiophene	0.00	0	0.00	0.00		
54)	Benz(a)anthracene	0.00	0	0.00	0.00		
55)	Chrysene	0.00	0	0.00	0.00		
56)	C1-Chrysenes	0.00	0	0.00	0.00		
57)	C2-Chrysenes	0.00	0	0.00	0.00		
58) 59)	C3-Chrysenes C4-Chrysenes	0.00 0.00	0	0.00	0.00		
64)	Benzo(b)fluoranthene	0.00	0	0.00	0.00		
65)	Benzo(k)fluoranthene	0.00	0	0.00	0.00 0.00		
66)	Benzo(e)pyrene	0.00	ő	0.00	0.00		
67)	Benzo(a)pyrene	0.00	ő	0.00	0.00		
75)	Perylene	0.00	0	0.00	0.00		
68)	Indeno(1,2,3-c,d)pyrene	0.00	0	0.00	0.00		
69)	Dibenzo(a,h)anthracene	0.00	0	0.00	0.00		
70)	C1-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00		
71)	C2-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00		
72)	C3-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00		
73)	Benzo(g,h,i)perylene	0.00	0	0.00	0.00		
	Total PAH				#DIV/0!		
	Individual Isomers						
9)	2-Methylnaphthalene	0.00	0	0.00	0.00		
10)	1-Methylnaphthalene	0.00	o o	0.00	0.00		
11)	2,6-Dimethylnaphthalene	0.00	0	0.00	0.00		
12)	1,6,7-Trimethylnaphthalene	0.00	0	0.00	0.00		
39)	1-Methylphenanthrene	0.00	0	0.00	0.00		
61)	C29-Hopane	0.00	0	0.00	0.00		
62)	18a-Oleanane	0.00	0	0.00	0.00		
63)	C30-Hopane	0.00	0	0.90	0.00		
	Surrogates				Su Recovery (%)		
31	(AR-STSU-040-005)	40.00		***	100		
2)	Naphthalene-d8 Acenaphthene-d10	13.01	4427	50.01	100		
20) 30)	Acenaphthene-d10 Phenanthrene-d10	18.83 23.91	2077	48.69	97		
53)	Chrysene-d12	23.91 33.00	3848 4953	47.59 45.31	95 91		
74)	Perviene-d12	37.70	2567	45.31 46.51	93		
,	Internal Stds	31.70	2001	40.31	24		
	(AR-WKIS-0500-007)						
1)	Fluorene-d10	20.63	2366	51.08			
29)	Pyrene-d10	28.79	4762	49.98			
60)	Benzo(a)pyrene-d12	37.42	3059	45.61			

Quant Results File: 041507.RES

Data File : X:\1\DATA\MS30377\MS30377C.D
Acq On : 13 Apr 2007 11:22 am
Sample : IS/SU Mixture Vial: 3 Operator: TJM

Inst : GC/MS Ins

Misc

Quant Time: Apr 14 20:37 2007

Multiplr: 1.00 MS Integration Params: rteint.p

Quant Method: C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002)
Last Update : Sat Apr 14 20:04:44 2007
Response via : Initial Calibration
DataAcq Meth : PAH-2002

Internal Standards	R.T.	QIon	Response	Conc Units Dev(Min)
1) Fluorene-d10	20.63	176	2366m	51.08 ng/ml 0.00
1) Fluorene-d1029) Pyrene-d1060) Benzo(a)pyrene-d12	28.79	212	4762m	49.98 0.00
60) Benzo(a)pyrene-d12	37.42	264	3059m	45.61 0.00
(
System Monitoring Compounds				
2) Naphthalene-d8 20) Acenaphthene-d10	13.01	136	4427	50.01 0.00
20) Acenaphthene-d10	18.83	164	2077m	
30) Phenanthrene-d10	23.91	188	3848	47.59 0.00
53) Chrysene-d12	33.00	240	4953m	45.31 0.00 46.51 0.00
74) Perylene-d12	37.70	264	2567m	46.51 0.00
Target Compounds				Qvalue
3) Decalin	0.00	138	0	N.D. d
	0.00	152	Ő	N.D. d
5) C2-Decalin	0.00	166	Ö	N.D. d
6) C3-Decalin	0.00	180	ő	N.D. d
7) C4-Decalin	0.00	194	Ő	N.D. d
7) C4-Decalin8) Naphthalene	0.00	128	O	N.D. d
9) 2-Methylnaphthalene 10) 1-Methylnaphthalene	0.00	142	Ö	N.D. d
10) 1-Methylnaphthalene	0.00	142	Ö	N.D. d
11) 2,6-Dimethylnaphthalene	0.00	156	Õ	N.D.
12) 1 6 7-Trimethylnaphthalene	0.00	170	Ö	N.D.
12) 1,6,7-Trimethylnaphthalene 13) C2-Naphthalenes	0.00	156	Ö	N.D. d
14) C3-Naphthalenes	0.00	170	Ō	N.D. d
15) C/-Naphthalenes	0.00	184	Õ	N.D. d
15) C4-Naphthalenes 16) Benzothiophene 17) C1-Benzothiophene	0.00	134		N.D.
17) Cl-Benzothionhene	0.00	148	Ö	N.D. d
18) C2-Benzothiophene	0.00	162	Õ	N.D. d
19) C3-Benzothiophene	0.00	176	Ō	N.D. d
19) C3-Benzothiophene 21) Biphenyl	0.00		Õ	N.D. d
21) Biphenyl 22) Acenaphthylene 23) Acenaphthene 24) Dibenzofuran 25) Fluorene 26) C1-Fluorenes 27) C2-Fluorenes 28) C3-Fluorenes	0.00	152	Ō	N.D.
23) Acenaphthene	0.00	154	0	N.D. d
24) Dibenzofuran	0.00	168	O	N.D. d
25) Fluorene	0.00	166	0	N.D. d
26) C1-Fluorenes	0.00	180	0	N.D. d
27) C2-Fluorenes	0.00	194	0	N.D. d
28) C3-Fluorenes	0.00	208	0	N.D. d
31) Pentachlorophenol	0.00	266	0	N.D.
32) Carbazole 33) Dibenzothiophene	0.00	167	0	N.D. d
33) Dibenzothiophene	0.00	184	0	N.D. d
34) C1-Dibenzothiophene	0.00	198	0	N.D. d
35) C2-Dibenzothiophene	0.00	212	0	N.D. d
36) C3-Dibenzothiophene	0.00	226	0	N.D. d
37) Phenanthrene	0.00	178	0	N.D.
38) Anthracene	0.00	178	0	N.D.
39) 1-Methylphenanthrene	0.00	192	0	N.D.
40) C1-Phenanthrene/Anthracene	0.00	192	0	N.D. d
41) C2-Phenanthrene/Anthracene	0.00	206	0	N.D. d
42) C3-Phenanthrene/Anthracene	0.00	220	0	N.D. d
43) C4-Phenanthrene/Anthracene	0.00	234	0	N.D. d
44) Naphthobenzothiophene	0.00	234	0	N.D.
45) C1-Naphthobenzothiophene	0.00	248	0	N.D. d
46) C2-Naphthobenzothiophene	0.00	262	0	N.D. d

Data File : X:\1\DATA\MS30377\MS30377C.D Vial: 3

Acq On : 13 Apr 2007 11:22 am Sample : IS/SU Mixture Misc : Operator: TJM Inst : GC/MS Ins

Multiplr: 1.00

MS Integration Params: rteint.p

Quant Results File: 041507.RES Quant Time: Apr 14 20:37 2007

Quant Method: C:\MS30377\041507.M (RTE Integrator)
Title: PAH Calibration Table (2002)
Last Update: Sat Apr 14 20:04:44 2007
Response via: Initial Calibration
DataAcq Meth: PAH-2002

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
47)	C3-Naphthobenzothiophene	0.00	276	0	N.D. d	
48)		0.00	202	0	N.D.	
49)	Pyrene	0.00	202	0	N.D.	
50)	C1-Fluoranthenes/Pyrenes	0.00	216	0	N.D. d	
51)	C2-Fluoranthenes/Pyrenes	0.00	230	0	N.D. d	
52)	C3-Fluoranthenes/Pyrenes	0.00	244	0	N.D. d	
54)	Benz(a)anthracene	0.00	228	0	N.D.	
55)	Chrysene	0.00	228	0	N.D.	
56)	Cl-Chrysenes	0.00	242	0	N.D. d	
57)	C2-Chrysenes	0.00		0	N.D. d	
58)	C3-Chrysenes	0.00	270	0	N.D. d	
59)	C4-Chrysenes	0.00	284	0	N.D. d	
61)	C29-Hopane	0.00		0	N.D. d	
62)	18a-Oleanane	0.00	191	0	N.D. d	
63)	C30-Hopane	0.00	191	0	N.D.	
64)	• •	0.00	252	0	N.D.	
65)	Benzo(k)fluoranthene	0.00		0	N.D.	
66)	Benzo(e)pyrene	0.00		0	N.D.	
67)	Benzo(a)pyrene	0.00	252	0	N.D.	
68)	Indeno(1,2,3-c,d)pyrene	0.00		0	N.D. d	
69)	Dibenzo(a,h)anthracene	0.00		0	N.D. d	
70)	Cl-Dibenzo(a,h)anthracene	0.00	292	0	N.D. d	
71)	C2-Dibenzo(a,h)anthracene	0.00	306	0	N.D. d	
72)	C3-Dibenzo(a,h)anthracene	0.00	320	0	N.D. d	
73)		0.00	276	0	N.D. d	
75)	Perylene	0.00	252	0	N.D.	

041507.RES

Quant Results File:

Ins

GC/MS 1.00

Multiplr:

TUM

Operator:

X:\1\DATA\MS30377\MS30377C.D 13 Apr 2007 11:22 am IS/SU Mixture

Data File

Acq On Sample

Misc

MS Integration Params: rteint.p Quant Time: Apr 14 20:37 2007

Method

Inst

Vial:

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Page
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                                                                                                                                                                                                                                                                                                                            24.00 26.00 28.00 30.00 32.00 34.00 36.00 38.00 40.00 42.00 44.00 46.00 48.00
                                                                                         Perylene-d12, gzo(a)pyrene-d12, I
                             TIC: MS30377C.D
X:\l\METHODS\041507.M (RTE Integrator)
PAH Calibration Table (2002)
Sat Apr 14 20:04:44 2007
Initial Calibration
                                                                         Chrysene-d12, 5
                                                                                                                                                                                                                                                                                                                                          2007
                                                                               1,01b-enery9
                                                                                                                                                                                                                                                                                                                                          06:52:53
                                                                                                                                                                                                                                                                                                                                          08
                                                                                                                   Phenanthrene-d10 S
                                                                                                                                                                                                                                                                                                                                          Tue May
                                                                                                                                                                                                                                                                                                                            10.00 12.00 14.00 16.00 18.00 20.00 22.00
                                                                                                      Fluorene-d10, }
                                                                                                                      Acenaphthene-d10, S
                                                                                                                                                                                                                                                                                                                                          041507.M
                       Response via
               Last Update
                                                                                                                              Raphthalene-d8, S
                                                                                                                                                                                                                                                                                                                                         MS30377C.D
       Title
                             Abundance
                                           5500
                                                                   5000
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Tissue, Sediment, and Water Sample Report (Use d-10 Phenanthrene only for Surrogate Corrections)

Data File Name ETX7207.D
Data File Path X:\1\DATA\MS30377\
Operator TJM
Date Acquired 04/13/20 -1:6:
Method File PAH-2002
Sample Name WIF-02-32707
Misc Info
Instrument Name GC/MS Ins
Vial Number 5
Sample Multiplier 1.1314
Sample Amount 0

Su Amt = 50

ETX7207.D WIF-02-32707

04/13/20 -1:6: PAH-2002 0.88

Peak#	Compound	Ret Time (min)	Target Response (Area)	Conc. (ng/g or ng/L)	Su, Corrected Conc. (ng/g or ng/L)
3)	Decalin	0.00	0	0.00	0.00
4)	C1-Decalin	0.00	0	0.00	0.00
5)	C2-Decalin	0.00	0	0.00	0.00
6)	C3-Decalin	0.00	0	0.00	0.00
7)	C4-Decalin	0.00	0	0.00	0.00
8)	Naphthalene	13.10	96	0.11	0.12
9+10)	C1-Naphthalenes	15.50 17.71	597	0.69 304,58	0.76 335,48
13) 14)	C2-Naphthalenes C3-Naphthalenes	19.73	262061 1041965	1211.00	1333.89
15)	C4-Naphthalenes	22.07	881691	1024.73	1128.72
16)	Benzothiophene	13.35	10	0.01	0.02
17)	C1-Benzothiophene	15.63	1050	1.44	1.59
18)	C2-Benzothiophene	17.46	14279	19.60	21.59
19)	C3-Benzothiophene	19.54	50148	68.85	75,84
21)	Biphenyl	16.92	79	0.11	0.13
22)	Acenaphthylene	0.00 18.95	0	0.00	0,00
23) 24)	Acenaphthene Dibenzofuran	19,56	13200 12622	28.06 16.55	30.90 18.23
25)	Fluorene	20.75	40410	73.80	81.29
26)	C1-Fluorenes	22,74	193648	353,68	389.57
27)	C2-Fluorenes	24.38	528750	965.70	1063.70
28)	C3-Fluorenes	26.07	677039	1236.53	1362.01
31)	Pentachlorophenol	0.00	0	0.00	0.00
32)	Carbazole	0.00	0	0.00	0.00
38)	Anthracene	24.18	51366	51.92	57.19
37)	Phenanthrene	24.01	256829	288,01	317.24
40)	C1-Phenanthrene/Anthracene C2-Phenanthrene/Anthracene	25.80	1882469	2111.01	2325.24 4966.81
41) 42)	C3-Phenanthrene/Anthracene	27.62 29.20	4021037 4238955	4509.22 4753.59	4900.81 5235.99
43)	C4-Phenanthrene/Anthracene	31.05	2193555	2459.87	2709.49
33)	Dibenzothiophene	23.57	63908	68,57	75.53
34)	C1-Dibenzothiophene	25.09	291905	313.21	345.00
35)	C2-Dibenzothiophene	26.84	623978	669.52	737.47
36)	C3-Dibenzothiophene	28,05	721721	774.40	852.99
48)	Fluoranthene	28.12	92853	80.66	88.85
49)	Pyrene	28.90	639803	506.25	557.62
50)	C1-Fluoranthenes/Pyrenes	30,41	2606972	2264.69	2494.51
51)	C2-Fluoranthenes/Pyrenes	32.16 33.39	3490858	3032.53	3340.27
52) 44)	C3-Fluoranthenes/Pyrenes Naphthobenzothiophene	32.20	2834421 260280	2462.28 298.57	2712.15 328.87
45)	C1-Naphthobenzothiophene	33.63	632695	725.77	799.42
46)	C2-Naphthobenzothiophene	35.08	774814	888.79	978.99
47)	C3-Naphthobenzothiophene	36.14	359886	412.83	454.72
54)	Benz(a)anthracene	33.00	338217	288.27	317.52
55)	Chrysene	33.14	684544	700.75	771.86
56)	C1-Chrysenes	34.41	2619706	2681.72	2953.86
57)	C2-Chrysenes	35.72	2866092	2933.94	3231.68
58)	C3-Chrysenes	37.20	1208084	1236.68	1362.18
59) 64)	C4-Chrysenes Benzo(b)fluoranthene	40.34 36.46	37956 135753	38.85 125.72	42.80 138.48
65)	Benzo(k)fluoranthene	36.57	17485	16.74	18.43
66)	Benzo(e)pyrene	37.38	146108	134.34	147.98
67)	Benzo(a)pyrene	37.56	233396	252.67	278,31
75)	Perylene	37.84	64000	64.60	71.16
68)	Indeno(1,2,3-c,d)pyrene	41.89	14942	21.30	23.46
69)	Dibenzo(a,h)anthracene	41.97	21789	33.49	36.89
70)	C1-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00
71)	C2-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00
72)	C3-Dibenzo(a,h)anthracene	0.00	0	0.00	0.00
73)	Benzo(g,h,i)perylene	43.11	32574	41,54	45.75
	Total PAH				44663
	Individual Isomers				
9)	2-Methylnaphthalene	15.35	138	0.25	0.27
10)	1-Methylnaphthalene	15.66	459	88.0	0.97
11)	2,6-Dimethylnaphthalene	17.46	26312	57.44	63.27
12)	1,6,7-Trimethylnaphthalene	20.30	79144	187.10	206.09
39)	1-Methylphenanthrene	26.13	225481	296.28	326.34
61)	C29-Hopane	39.86	26503	64.62	71.18
62) 63)	18a-Oleanane C30-Hopane	40.88	4996 31630	12.18	13.42
63)	Сэо-поране	41.09	31000	77.12	84.95
	Surrogates (AR-STSU-040-005)				Su Recovery (%)
2)	Naphthalene-d8	13.01	37610	47.74	84
20)	Acenaphthene-d10	18.86	19352	50.98	90
30)	Phenanthrene-d10	23.94	38637	51.36	91
53)	Chrysene-d12	33.03	52149	51.28	91
74)	Perylene-d12	37.73	27137	54.58	96
	Internal Stds				
41	(AR-WKIS-0500-007)	00.00	nane -	21.00	
1) 29)	Fluorene-d10 Pyrene-d10	20.63 28.83	23824 50123	51.08 49.98	
60)	Benzo(a)pyrene-d12	37.49	31179	45.61	
,	the second second	21.1.0	2	,0.01	

Data File : X:\1\DATA\MS30377\ETX7207.D

Vial: 5 Acq On : 13 Apr 2007 6:43 pm Sample : WIF-02-32707 Misc : Operator: TJM

Inst : GC/MS Ins

Multiplr: 1.13

MS Integration Params: rteint.p

Quant Time: Apr 20 16:23 2007 Quant Results File: 041507.RES

Quant Method: C:\MS30377\041507.M (RTE Integrator)
Title: PAH Calibration Table (2002)
Last Update: Sat Apr 14 20:04:44 2007
Response via: Initial Calibration
DataAcq Meth: PAH-2002

Tituorene-d10			QIon	Response	Conc Ur	nits D	ev(Min)
System Monitoring Compounds 20 Naphthalene=d8 13.01 136 37610m 47.74 0.00 200 Aconaphthene=d10 18.86 164 19352m 50.98 0.00 300 Phenanthrene=d10 23.94 188 38637m 51.36 0.03 531 Chrysene=d12 33.03 240 52149m 51.28 0.03 74) Perylene=d12 37.73 264 27137m 54.58 0.03 0.00 0.	1) Fluorene-d10		176	23824m	51.08	na/ml	0.00
System Monitoring Compounds 2) Naphthalene-d8 13.01 136 37610m 47.74 0.00 20) Aconaphthene-d10 18.86 164 19352m 50.98 0.00 30) Phenanthrene-d10 23.94 188 38637m 51.36 0.03 53) Chrysene-d12 33.03 240 52149m 51.28 0.03 74) Perylene-d12 37.73 264 27137m 54.58 0.03 740 74.74 0.00 166 0.00 0.00 0.	29) Pyrene-d10		212	50123m	49.98	5,	0.03
System Monitoring Compounds 2 Naphthaleneeds	60) Benzo(a)pyrene-d12	37.49	264	31179m	45.61		0.07
20							
200 Acenaphthene-d10							
300 Phenanthrene-d10	2) Naphthalene-d8						
300 Phenanthrene-d10	20) Acenaphthene-d10						
Target Compounds	30) Phenanthrene-d10		188	38637m			
Target Compounds			240	52149m			
3 Decalin	74) Perylene-dl2	3/./3	264	2/13/m	54.58		0.03
3) Decalin 4) Cl-Decalin 0.00 152 0 N.D. d 5) C2-Decalin 0.00 166 0 N.D. d 6) C3-Decalin 0.00 180 0 N.D. d 7) C4-Decalin 0.00 180 0 N.D. d 8) Naphthalene 13.10 128 96m 0.11 9) 2-Methylnaphthalene 15.35 142 138m 0.25 10) 1-Methylnaphthalene 17.46 156 26312m 57.44 12) 1,6-Dimethylnaphthalene 17.46 156 226312m 57.44 12) 1,6-Trimethylnaphthalene 17.46 156 26312m 57.44 12) 1,6,7-Trimethylnaphthalene 18.02-Naphthalenes 19.73 170 1041965m 1211.00 15) C4-Naphthalenes 19.73 170 1041965m 1211.00 15) C4-Naphthalenes 19.73 170 1041965m 1211.00 15) C4-Naphthalenes 19.73 170 1041965m 1211.00 16) Benzothiophene 13.35 134 10m 0.01 ng/ml 17) Cl-Benzothiophene 15.63 148 1050m 1.44 ng/ml 18) C2-Benzothiophene 17.46 162 14279m 19.60 ng/ml 19) C3-Benzothiophene 19.54 176 50148m 68.85 ng/ml 19) Biphenyl 16.92 154 79m 0.11 12) Acenaphthylene 20.00 152 0 N.D. d 23) Acenaphthylene 20.10 152 0 N.D. d 24) Dibenzoturan 19.56 168 1262zm 16.55 ng/ml 25) Fluorene 20.75 166 40410m 73.80 260 C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28 C3-Fluorenes 24.38 194 528750m 965.70 29 C3-Fluorenes 24.38 194 528750m 965.70 29 C3-Fluorenes 26.07 208 677039m 669.52 30 Dibenzothiophene 23.57 184 63908m 68.57 31) Pentachlorophene 24.18 178 51366m 51.92 360 C3-Dibenzothiophene 25.09 198 291905m 313.21 37) Phenanthrene 24.01 178 25682pm 288.01 38) Anthracene 24.18 178 51366m 51.92 24.18 178 51366m 51.92 25781m 296.28 261 C1-Phenanthrene/Anthracene 24.18 178 51366m 51.92 262 C3-Phenanthrene/Anthracene 24.18 178 51366m 51.92 270 298.57 271 21 174 40 272 274 274 277 277 277 277 277 277 277	Target Compounds						Ovalue
C1-Decalin		0.00	138	0	N.D.		Qvarae
S		0 00	152				
8) C3-Decalin 0.00 180 0 N.D. d 8) Naphthalene 13.10 128 96m 0.11 9) 2-Methylnaphthalene 15.35 142 138m 0.25 10) 1-Methylnaphthalene 15.66 142 459m 0.88 11) 2,6-Dimethylnaphthalene 17.46 156 26312m 57.44 12) 1,6,7-Trimethylnaphthalene 17.46 156 26312m 304.58 11) 2,6-Dimethylnaphthalene 17.71 156 262061m 304.58 14) C3-Naphthalenes 19.73 170 1041965m 1211.00 15) C4-Naphthalenes 22.07 184 881691m 1024.73 16) Benzothiophene 13.35 134 1050m 1.44 ng/ml 17) C1-Benzothiophene 15.63 148 1050m 1.44 ng/ml 18) C2-Benzothiophene 17.46 162 14279m 19.60 ng/ml 19) C3-Benzothiophene 19.54 176 50148m 68.85 ng/ml 21) Biphenyl 16.92 154 79m 0.11 22) Acenaphthylene 0.00 152 0 N.D. d 23) Acenaphthene 18.95 154 13200m 28.06 24) Dibenzofuran 19.56 168 12622m 16.55 ng/ml 25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.438 194 528750m 965.70 28) C3-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 3) Pentachlorophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 30) C1-Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 26.13 192 25481m 296.28 40) C1-Phenanthrene/Anthracene 24.18 178 51366m 51.92 20 C3-Phenanthrene/Anthracene 24.02 24 2438955m 4753.59 41) Naphthobenzothiophene 32.09 234 260280m 298.57 45) C1-Naphthobenzothiophene 32.09 234 260280m 298.57 45) C2-Naphthobenzothiophene 32.09 234 260280m 298.57 45) C2-Naphthobenzothiophene 33.63 248 632695m 725.77 45) C2-Naphthobenzothiophene 33.63 248 632695m 725.77 45) C2-Naphthobenzothiophene 33.63 248 632695m 725.77		0.00	166				
Naphthalene		0.00	180				
8) Naphthalene	7) C4-Decalin	0.00					
9) 2-Methylnaphthalene	8) Naphthalene	13.10	128				
10) 1-Methylnaphthalene	9) 2-Methylnaphthalene	15.35	142	138m	0.25		
11				459m	0.88		
12) 1,6,7-Trimethylnaphthalene 20,30 170 79144m 187.10 130 C2-Naphthalenes 19,73 170 1041965m 1211.00 151 C4-Naphthalenes 19,73 170 1041965m 1211.00 151 C4-Naphthalenes 22,07 184 881691m 1024.73 160 Benzothiophene 13,35 134 10m 0.01 ng/ml 170 C1-Benzothiophene 15,63 148 1050m 1,44 ng/ml 180 C2-Benzothiophene 17,46 162 14279m 19,60 ng/ml 190 C3-Benzothiophene 19,54 176 50148m 68.85 ng/ml 190 120 N.D. d 120 N.D. d					57.44		
13) C2-Naphthalenes 17.71 156 262061m 304.58 14) C3-Naphthalenes 19.73 170 1041965m 1211.00 15) C4-Naphthalenes 22.07 184 881691m 1024.73 16) Benzothiophene 13.35 134 10m 0.01 ng/ml 17) C1-Benzothiophene 15.63 148 1050m 1.44 ng/ml 18) C2-Benzothiophene 17.46 162 14279m 19.60 ng/ml 19) C3-Benzothiophene 19.54 176 50148m 68.85 ng/ml 21) Biphenyl 16.92 154 79m 0.11 22) Acenaphthylene 0.00 152 0 N.D. d 23) Acenaphthene 18.95 154 13200m 28.06 24) Dibenzofuran 19.56 168 12622m 16.55 ng/ml 25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenol 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 27.09 198 291905m 313.21 35) C2-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 44.1 Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 35.08 262 774814m 888.79							
14) C3-Naphthalenes 19.73 170 1041965m 1211.00 15) C4-Naphthalenes 22.07 184 881691m 1024.73 16) Benzothiophene 13.35 134 10m 0.01 ng/ml 17) C1-Benzothiophene 15.63 148 1050m 1.44 ng/ml 18) C2-Benzothiophene 17.46 162 14279m 19.60 ng/ml 19. C3-Benzothiophene 17.46 162 14279m 19.60 ng/ml 19. C3-Benzothiophene 19.54 176 50148m 68.85 ng/ml 21) Biphenyl 16.92 154 79m 0.11 22) Acenaphthylene 0.00 152 0 N.D. d 23) Acenaphthene 18.95 154 79m 0.11 22) Acenaphthene 19.56 168 12622m 16.55 ng/ml 23) Acenaphthene 19.56 168 12622m 16.55 ng/ml 24) Dibenzofuran 19.56 168 12622m 16.55 ng/ml 25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 24.38 194 528750m </td <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>							
15) C4-Naphthalenes							
17) C1-Benzothiophene 15.63 148 1050m 1.44 ng/ml 18) C2-Benzothiophene 17.46 162 14279m 19.60 ng/ml 19) C3-Benzothiophene 19.54 176 50148m 68.85 ng/ml 21) Biphenyl 16.92 154 79m 0.11 22) Acenaphthene 18.95 154 13200m 28.06 24) Dibenzofuran 19.56 168 12622m 16.55 ng/ml 25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenel 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene	15) C4-Naphthalenes						
17) C1-Benzothiophene 15.63 148 1050m 1.44 ng/ml 18) C2-Benzothiophene 17.46 162 14279m 19.60 ng/ml 19) C3-Benzothiophene 19.54 176 50148m 68.85 ng/ml 21) Biphenyl 16.92 154 79m 0.11 22) Acenaphthene 18.95 154 13200m 28.06 24) Dibenzofuran 19.56 168 12622m 16.55 ng/ml 25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenel 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene	16) Benzothiophene					ng/ml	
18) C2-Benzothiophene 17.46 162 14279m 19.60 ng/ml 19) C3-Benzothiophene 19.54 176 50148m 68.85 ng/ml 21) Biphenyl 16.92 154 79m 0.11 22) Acenaphthylene 0.00 152 0 N.D. d 23) Acenaphthene 18.95 154 13200m 28.06 24) Dibenzofuran 19.56 168 12622m 16.55 ng/ml 25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenol 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. 33) Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 25.09	17) C1-Benzothiophene		148	1050m	1.44	na/ml	
21) Biphenyl	18) C2-Benzothiophene		162	14279m	19.60	ng/ml	
21) Biphenyl	19) C3-Benzothiophene		176	50148m	68.85	ng/ml	
22) Acenaphthylene 0.00 152 0 N.D. d 23) Acenaphthene 18.95 154 13200m 28.06 24) Dibenzofuran 19.56 168 12622m 16.55 ng/ml 25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenol 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92	21) Biphenyl	16.92	154		0.11		
23) Acenaphthene 24) Dibenzofuran 19.56 168 1262zm 16.55 ng/ml 25) Fluorene 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenol 32) Carbazole 33) Dibenzothiophene 34.37 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 27.02 28.05 226 721721m 774.40 37) Phenanthrene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 39) 1-Methylphenanthrene 40) C1-Phenanthrene/Anthracene 40) C1-Phenanthrene/Anthracene 41) C2-Phenanthrene/Anthracene 42.18 178 51366m 51.92 39) 1-Methylphenanthrene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 42.18 178 51366m 51.92 39) 1-Methylphenanthrene 40) C1-Phenanthrene/Anthracene 41) C2-Phenanthrene/Anthracene 42.20 234 2438955m 4753.59 43) C4-Phenanthrene/Anthracene 43) C4-Phenanthrene/Anthracene 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79	22) Acenaphthylene	0.00	152				
25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenol 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. d 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 25.80 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05 234 21935555m 2459.87 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79				13200m	28.06		
25) Fluorene 20.75 166 40410m 73.80 26) C1-Fluorenes 22.74 180 193648m 353.68 27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenol 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. d 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 25.80 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05 234 21935555m 2459.87 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79	24) Dibenzofuran	19.56	168	12622m	16.55	ng/ml	
27) C2-Fluorenes 24.38 194 528750m 965.70 28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenol 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 26.13 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05	25) Fluorene	20.75	166	40410m	73.80		
28) C3-Fluorenes 26.07 208 677039m 1236.53 31) Pentachlorophenol 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 26.13 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 31.05 234 2193555m 4753.59 43) C4-Phenanthrene/Anthracene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79		22.74	180	193648m	353.68		
31) Pentachlorophenol 0.00 266 0 N.D. 32) Carbazole 0.00 167 0 N.D. d 33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 26.13 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05 234 2193555m 2459.87 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79	27) C2-Fluorenes	24.38	194	528750m	965.70		
32) Carbazole 33) Dibenzothiophene 33, Dibenzothiophene 33, Dibenzothiophene 33, Dibenzothiophene 33, Dibenzothiophene 33, Dibenzothiophene 35, O2-Dibenzothiophene 36, C3-Dibenzothiophene 37) Phenanthrene 38) Anthracene 39) 1-Methylphenanthrene 40) C1-Phenanthrene/Anthracene 40) C1-Phenanthrene/Anthracene 41) C2-Phenanthrene/Anthracene 42) C3-Phenanthrene/Anthracene 43) C4-Phenanthrene/Anthracene 44) Naphthobenzothiophene 35, O2-Dibenzothiophene 36, O3-Dibenzothiophene 36, O3-Dibenzothiophene 37) Phenanthrene 28, O5				677039m	1236.53		
33) Dibenzothiophene 23.57 184 63908m 68.57 34) C1-Dibenzothiophene 25.09 198 291905m 313.21 35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 26.13 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05 234 2193555m 2459.87 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 35.08 262 774814m 888.79	31) Pentachlorophenol	0.00	266	0	N.D.	•	
34) C1-Dibenzothiophene25.09198291905m313.2135) C2-Dibenzothiophene26.84212623978m669.5236) C3-Dibenzothiophene28.05226721721m774.4037) Phenanthrene24.01178256829m288.0138) Anthracene24.1817851366m51.9239) 1-Methylphenanthrene26.13192225481m296.2840) C1-Phenanthrene/Anthracene25.801921882469m2111.0141) C2-Phenanthrene/Anthracene27.622064021037m4509.2242) C3-Phenanthrene/Anthracene29.202204238955m4753.5943) C4-Phenanthrene/Anthracene31.052342193555m2459.8744) Naphthobenzothiophene32.20234260280m298.5745) C1-Naphthobenzothiophene33.63248632695m725.7746) C2-Naphthobenzothiophene35.08262774814m888.79	32) Carbazole	0.00	167	0	N.D.	. d	
35) C2-Dibenzothiophene 26.84 212 623978m 669.52 36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 26.13 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05 234 2193555m 2459.87 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79	33) Dibenzothiophene		184				
36) C3-Dibenzothiophene 28.05 226 721721m 774.40 37) Phenanthrene 24.01 178 256829m 288.01 38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 26.13 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05 234 2193555m 2459.87 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79	34) Cl-Dibenzothiophene						
37) Phenanthrene24.01178256829m288.0138) Anthracene24.1817851366m51.9239) 1-Methylphenanthrene26.13192225481m296.2840) C1-Phenanthrene/Anthracene25.801921882469m2111.0141) C2-Phenanthrene/Anthracene27.622064021037m4509.2242) C3-Phenanthrene/Anthracene29.202204238955m4753.5943) C4-Phenanthrene/Anthracene31.052342193555m2459.8744) Naphthobenzothiophene32.20234260280m298.5745) C1-Naphthobenzothiophene33.63248632695m725.7746) C2-Naphthobenzothiophene35.08262774814m888.79							
38) Anthracene 24.18 178 51366m 51.92 39) 1-Methylphenanthrene 26.13 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05 234 2193555m 2459.87 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79							
39) 1-Methylphenanthrene 26.13 192 225481m 296.28 40) C1-Phenanthrene/Anthracene 25.80 192 1882469m 2111.01 41) C2-Phenanthrene/Anthracene 27.62 206 4021037m 4509.22 42) C3-Phenanthrene/Anthracene 29.20 220 4238955m 4753.59 43) C4-Phenanthrene/Anthracene 31.05 234 2193555m 2459.87 44) Naphthobenzothiophene 32.20 234 260280m 298.57 45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79	· · · · · · · · · · · · · · · · · · ·						
40) C1-Phenanthrene/Anthracene25.801921882469m2111.0141) C2-Phenanthrene/Anthracene27.622064021037m4509.2242) C3-Phenanthrene/Anthracene29.202204238955m4753.5943) C4-Phenanthrene/Anthracene31.052342193555m2459.8744) Naphthobenzothiophene32.20234260280m298.5745) C1-Naphthobenzothiophene33.63248632695m725.7746) C2-Naphthobenzothiophene35.08262774814m888.79	•						
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42) C3-Phenanthrene/Anthracene29.202204238955m4753.5943) C4-Phenanthrene/Anthracene31.052342193555m2459.8744) Naphthobenzothiophene32.20234260280m298.5745) C1-Naphthobenzothiophene33.63248632695m725.7746) C2-Naphthobenzothiophene35.08262774814m888.79							
43) C4-Phenanthrene/Anthracene31.052342193555m2459.8744) Naphthobenzothiophene32.20234260280m298.5745) C1-Naphthobenzothiophene33.63248632695m725.7746) C2-Naphthobenzothiophene35.08262774814m888.79							
44) Naphthobenzothiophene32.20234260280m298.5745) C1-Naphthobenzothiophene33.63248632695m725.7746) C2-Naphthobenzothiophene35.08262774814m888.79							
45) C1-Naphthobenzothiophene 33.63 248 632695m 725.77 46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79	·						
46) C2-Naphthobenzothiophene 35.08 262 774814m 888.79							

Data File : X:\1\DATA\MS30377\ETX7207.D

Vial: 5 Acq On : 13 Apr 2007 Sample : WIF-02-32707 Misc : 6:43 pm Operator: TJM

Inst : GC/MS Ins

Multiplr: 1.13

MS Integration Params: rteint.p

Quant Time: Apr 20 16:23 2007 Quant Results File: 041507.RES

Quant Method: C:\MS30377\041507.M (RTE Integrator)
Title: PAH Calibration Table (2002)
Last Update: Sat Apr 14 20:04:44 2007
Response via: Initial Calibration
DataAcq Meth: PAH-2002

	Compound	R.T.	QIon	Response	Conc Ur	nit Qvalue
47)	C3-Naphthobenzothiophene	36.14	276	359886m	412.83	
48)	Fluoranthene	28.12	202	92853m	80.66	
49)	Pyrene	28.90	202	639803m	506.25	
50)	C1-Fluoranthenes/Pyrenes	30.41	216	2606972m	2264.69	ng/mL
51)	C2-Fluoranthenes/Pyrenes	32.16	230	3490858m	3032.53	ng/mL
52)	C3-Fluoranthenes/Pyrenes	33.39	244	2834421m	2462.28	ng/mL
54)	Benz(a)anthracene	33.00	228	338217m	288.27	
55)	Chrysene	33.14	228	684544m	700.75	
56)	C1-Chrysenes	34.41	242	2619706m	2681.72	ng/mL
57)	C2-Chrysenes	35.72	256	2866092m	2933.94	ng/mL
58)	C3-Chrysenes	37.20	270	1208084m	1236.68	ng/mL
59)	C4-Chrysenes	40.34	284	37956m	38.85	ng/mL
61)	C29-Hopane	39.86	191	26503m	64.62	ng/ml
62)	18a-Oleanane	40.88	191	4996m		
63)	C30-Hopane	41.09	191	31630m	77.12	ng/ml
64)	Benzo(b)fluoranthene	36.46	252	135753m	125.72	
65)	Benzo(k)fluoranthene	36.57	252	17485m	16.74	
66)	Benzo(e)pyrene	37.38	252	146108m	134.34	
67)	Benzo(a)pyrene	37.56	252	233396m	252.67	
68)	Indeno(1,2,3-c,d)pyrene	41.89	276	14942m	21.30	
69)	Dibenzo(a,h)anthracene	41.97	278	21789m	33.49	
70)	C1-Dibenzo(a,h)anthracene	0.00	292	0	N.D.	. d
71)	C2-Dibenzo(a,h)anthracene	0.00	306	0	N.D.	. d
72)	* * *	0.00	320	0	N.D.	. d
73)	Benzo(g,h,i)perylene	43.11	276	32574m	41.54	
75)	Perylene	37.84	252	64000m	64.60	

041507.RES

F110:

Quant Results

GC/MS

Multiplr:

Inst

X:\1\DATA\MS30377\ETX7207.D 13 Apr 2007 6:43 pm WIF-02-32707

Data File

Acq On Sample Params: rteint.p 20 16:23 2007

Apr

Quant Time:

MS Integration

Misc

TUM

Vial: Operator:

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                                                O1-Chrysenes, un
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(RTE Integrator)
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                            C3-Phenanthrene/Anthracene, un
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     PAH Calibration Table (2002)
Sat Apr 14 20:04:44 2007
Initial Calibration
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Total Petroleum Hydrocarbons/ Aliphatic Hydrocarbons Initial Calibration Data

ALI ICAL C10B0414.M

GCFID (ALI-COMP)

Response Factor Report GC#1

Method : W:\2\METHODS\C10B414.M (Chemstation Integrator)
Title : C10 - C35 aliphatic

Last Update : Sat Apr 14 11:05:52 2007

Calibration Files

=GC10846E.D 2 =GC10846F.D 3 =GC10846G.D =GC10846H.D 5 =GC10846I.D

4

		Compound	1	2	3	4	5	Avg	%RSD
1)		n-hexadecane-d34			Т	רויייב			
2)		n-C10	1.032	0.996	1.029	1.027	0.996	1.016.0	1.80
3)		n-C11	1.042	1.020	1.055	1.053	1.024	1.039 0	1.56
	S	n-dodecane-d26	0.867	0.845	0.878	0.881	0.853	0.865 0	1.80
5)		n-C12	1.068	1.069	1.109	1.106	1.075	1.085 0	1.86
6)		n-Cl3						1.119 0	
7)								1.164 0	
8)		n-C15						1.183 0	
9)		n-C16	1.219	1.187	1.225	1.206	1.166	1.201 0	2.03
10)		5a-androstane			IS	STD			_
								0.929 0	
12)		Pristane	1.007	0.972	0.996	0.995	0.964	0.987 0	1.83
13) 14)		n-C18 Phytane	1 021	0.923	0.955	0.957	0.925	0.933 0	2.38
15)		n-C19	0 924	0.970	0.990	0.992	0.901	0.934 0	2.28 1.76
16)		n-eicosane-d42	0.524	0.522	0.251	0.232	0.515	0.934 0	2.04
17)		n-C20	0.021	0.010	0.051	0.017	0.010	0.928 0	2.55
18)								0.942 0	
_								0.924 0	
_								0.918 0	
								0.912 0	
								0.906 0	
23)		n-C26	0.916	0.909	0.927	0.905	0.863	0.904 0	2.71
		n-C27	0.906	0.886	0.895	0.870	0.828	0.877 0	3.48
•								0.859 0	
								0.871 0	
27)	S	n-triacontane-d62	0.768						
28)		n-C30	0.874						
29)		n-C31	0.849					0.803 0	
-		n-C32	0.815					0.770 0	
		n-C33	0.772					0.747 0	
32)		n-C34	0.780					0.730 0	
33) 34)		n-C30 n-C31 n-C32 n-C33 n-C34 TPH TRH1	0.059					0.850 0 0.850 0	2.32
35)		TRH2	0.000					0.850 0	2.32
36)		TRH3	0.859					0.850 0	2.32
37)		TRH4						0.850 0	2.32
38)		TRH5						0.850 0	2.32
39)		TRH6						0.850 0	2.32

Vial: 96 Operator: TJM Data File : W:\2\DATA\GC10846\GC10846E.D Acq On : 13 Apr 2007 1:35 pm Sample : CS1 Sample : CS1
Misc :
IntFile : autoint1.e Inst : GC#1 Multiplr: 1.00

Quant Time: Apr 14 9:26 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI_COMP.M

Volume Inj. : Signal Phase : Signal Info :

Compound	R.T.	Response	Conc Units
Internal Standards			
1) n-hexadecane-d34			20.001 ug/mlm
10) 5a-androstane	18.03	465681	20.003 ug/mlm
System Monitoring Compounds			
4) S n-dodecane-d26	8.62		1.318 ug/mlm
16) S n-eicosane-d42	17.43		1.207 ug/mlm
27) S n-triacontane-d62	29.21f	22392	1.156 ug/mlm
Target Compounds			
2) n-C10	6.26	22172	1.350 ug/mlm
3) n-C11	7.59	22417	1.325 ug/mlm
5) n-C12	8.83	22972	1.286 ug/mlm
6) n-C13	9.98	23397	1.257 ug/mlm
7) n-C14	11.06	25674	1.325 ug/mlm
8) n-C15	12.08	25117	1.244 ug/mlm
9) n-C16	13.10	26268	1.268 ug/mlm
11) n-C17	14.18	26313	1.228 ug/mlm
12) Pristane	14.30f	29330	1.286 ug/mlm
13) $n-C18$	15.34	26427	1.212 ug/mlm
14) Phytane	15.50	29276	1.268 ug/mlm
15) n-C19	16.56f	26944	1.221 ug/mlm
17) n-C20	17.82f	26181	1.177 ug/mlm
18) n-C21	19.10	26351	1.148 ug/mlm
19) n-C22	20.39	26737	1.191 ug/mlm
20) n-C23	21.66	25945	1.134 ug/mlm
21) n-C24	22.91	26559	1.161 ug/mlm
22) n-C25	24.14	26963	1.182 ug/mlm
23) n-C26	25.33f	26515	1.145 ug/mlm
24) n-C27	26.48	26121	1.153 ug/mlm
25) n-C28	27.60	25760	1.142 ug/mlm
26) n-C29	28.69	27376	1.192 ug/mlm
28) n-C30	29.74f	25478	1.149 ug/mlm
29) n-C31	30.76f	24723	1.134 ug/mlm
30) n-C32	31.75	23381	1.109 ug/mlm
31) n-C33	32.74	22445	1.092 ug/mlm
32) n-C34	33.87f	22623	1.115 ug/mlm
,			⊸ .

______ (m) = manual int.

Data File : W:\2\DATA\GC10846\GC10846E.D
Acq On : 13 Apr 2007 1:35 pm
Sample : CS1 Vial: 96 Operator: TJM Inst : GC#1 Misc : IntFile : autoint1.e Multiplr: 1.00

Quant Time: Apr 14 9:26 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

R.T. Response Conc Units Compound ______

m

Quantitation Report

26.00 28.00 30.00 32.00 34.00 36.00 38.00 40.00 42.00 44.00 46.00 48.00 50.00 52.00 1.00 GC#1 TUN Multiplr: Vial: Operator: Quant Results File: C10B414.RES **78.88** (Chemstation Integrator) 32.74 Inst n-C31 31.75 92.08 Lu-C30 Lu-triacont Lu-C29 12.62 – 12.62 – 17.62 – GC10846E.D\FID2B P-C28 u-C5e 22.00 24.00 Multiple Level Calibration W:\2\DATA\GC10846\GC10846E.D 13 Apr 2007 1:35 pm Thu Apr 12 09:52:10 2007 10.00 12.00 14.00 16.00 18.00 20.00 C:\GC10846\C10B414.M C10 - C35 aliphatic 80.81 Ζ8,Ζ<u>Ϳ</u>ͺ 99'91 9:26 2007 ALI_COMP.M autoint1.e 6.8.83 800 Quant Time: Apr 14 69°Z 012-4-80 92.9 Quant Method Response via DataAcq Meth Phase 4.00 Volume Inj. Last Update Info Data File 2.00 IntFile Acq On Signal Sample Signal Title 0.0 Misc Response 26000 24000 18000 1 16000 10000 8000 -0009 4000 2000 -2000 22000 20000 14000-12000-Time

Data File: W:\2\DATA\GC10846\GC10846F.D Vial: 97 Acq On : 13 Apr 2007 2:36 pm Sample : CS2 Operator: TJM Inst : GC#1 Misc Multiplr: 1.00 Misc : IntFile : autoint1.e

Quant Time: Apr 14 9:33 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

	Compound	R.T.	Response	Conc Units
	ernal Standards			
1)	n-hexadecane-d34	12.85	360834	20.001 ug/mlm
10)	5a-androstane	18.03	477652	20.003 ug/mlm
Sys	tem Monitoring Compounds			
4) S	n-dodecane-d26	8.62	152639	10.280 ug/mlm
16) S	n-eicosane-d42	17.43	195478	9.627 ug/mlm
27) S	n-triacontane-d62	29.21f	177124	8.915 ug/mlm
Tar	get Compounds			
2)	n-C10	6.26	179015	10.424 ug/mlm
3)	n-C11	7.58	183624	10.376 ug/mlm
5)	n-C12	8.83	192328	10.297 ug/mlm
6)	n-C13	9.98	198661	10.205 ug/mlm
7)	n-C14	11.06	206908	10.210 ug/mlm
8)	n-C15	12.08	210963	9.992 ug/mlm
9)	n-C16	13.10	213947	9.870 ug/mlm
11)	n-C17	14.19	218486	9.943 ug/mlm
12)	Pristane	14.30	232288	9.933 ug/mlm
13)	n-C18	15.34	220860	9.871 ug/mlm
14)	Phytane	15.50	229648	9.697 ug/mlm
15)	n-C19	16.56	220477	9.741 ug/mlm
17)	n-C20	17.83f	219955	9.643 ug/mlm
18)	n-C21	19.11	224299	9.526 ug/mlm
19)	n-C22	20.40	213351	9.265 ug/mlm
20)	n-C23	21.67	219091	9.333 ug/mlm
21)	n-C24	22.92	216989	9.246 ug/mlm
22)	n-C25	24.14	216899	9.273 ug/mlm
23)	n-C26	25.33	215883	9.093 ug/mlm
24)	n-C27	26.49	209590	9.022 ug/mlm
25)	n-C28	27.61	209021	9.038 ug/mlm
26)	n-C29	28.69	211894	8.996 ug/mlm
28)	n-C30	29.75f	204524	8.996 ug/mlm
29)	n-C31	30.76	199765	8.936 ug/mlm
30)	n-C32	31.75	191369	8.847 ug/mlm
31)	n-C33	32.75	192709	9.144 ug/mlm
32)	n-C34	33.87	186784	8.972 ug/mlm

Vial: 97 Data File: W:\2\DATA\GC10846\GC10846F.D Acq On : 13 Apr 2007 2:36 pm Sample : CS2 Operator: TJM Inst : GC#1 Misc : IntFile : autoint1.e Multiplr: 1.00

Quant Time: Apr 14 9:33 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)
Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI_COMP.M

Volume Inj. : Signal Phase : Signal Info :

R.T. Response Conc Units Compound

Mon Apr 16 06:41:23 2007

C10B414.M

Quantitation Report

Vial:

Operator: Inst W:\2\DATA\GC10846\GC10846F.D 13 Apr 2007 2:36 pm Data File Acq On Sample

Misc

autoint1.e IntFile

9:33 2007 Quant Results File: C10B414.RES Quant Time: Apr 14

1.00 GC#1 MUT

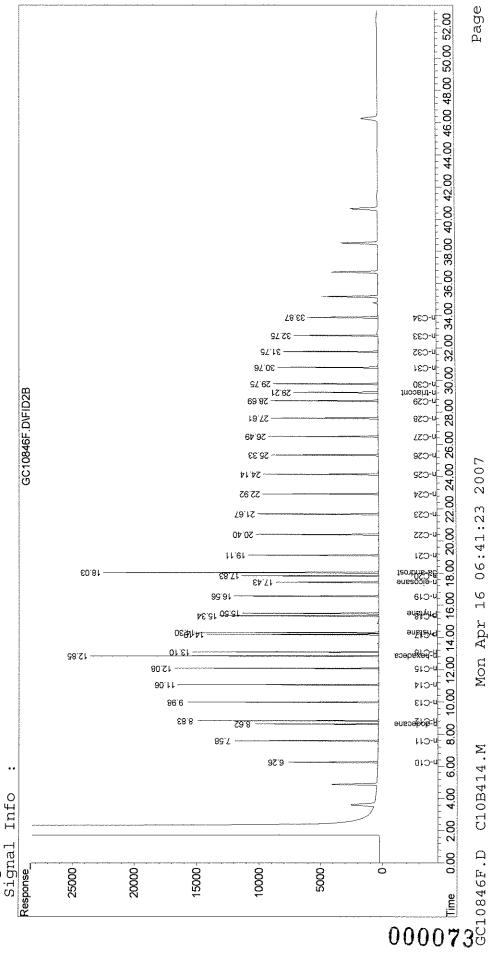
Multiplr:

C:\GC10846\C10B414.M (Chemstation Integrator)
C10 - C35 aliphatic Quant Method Title

Thu Apr 12 09:52:10 2007 Last Update

Multiple Level Calibration ALI_COMP.M DataAcq Meth Response via

Volume Inj. Signal Phase Signal



Data File: W:\2\DATA\GC10846\GC10846G.D Vial: 98 Acq On : 13 Apr 2007 3:37 pm Sample : CS3 Operator: TJM Inst : GC#1 Misc Misc : IntFile : autoint1.e Multiplr: 1.00

Quant Time: Apr 14 9:46 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

	Compound	R.T.	Response	Conc Units
	ernal Standards			
1)	n-hexadecane-d34	12.85	379900	20.001 ug/mlm
10)	5a-androstane	18.03	499362	20.003 ug/mlm
	tem Monitoring Compounds			
4) S	n-dodecane-d26	8.62	417507	26.708 ug/mlm
16) S	n-eicosane-d42	17.44	531251	25.027 ug/mlm
27) S	n-triacontane-d62	29.23	462301	22.256 ug/mlm
Tar	get Compounds			
2)	n-C10	6.26	487020	26.936 ug/mlm
3)	n-C11	7.59	499909	26.830 ug/mlm
5)	n-C12	8.83	524931	26.693 ug/mlm
6)	n-C13	9.99	543956	26.540 ug/mlm
7)	n-C14	11.07	562896	26.383 ug/mlm
8)	n-C15	12.09	573692	25.808 ug/mlm
9)	n-C16	13.11	581309	25.471 ug/mlm
11)	n-C17	14.19	593087	25.817 ug/mlm
12)	Pristane	14.31	622148	25.448 ug/mlm
13)	n-C18	15.35	597649	25.550 ug/mlm
14)	Phytane	15.51	612089	24.723 ug/mlm
15)	n-C19	16.57	594679	25.131 ug/mlm
17)	n-C20	17.84	596383	25.009 ug/mlm
18)	n-C21	19.12	606173	24.624 ug/mlm
19)	n-C22	20.41	577229	23.977 ug/mlm
20)	n-C23	21.68	591849	24.117 ug/mlm
21)	n-C24	22.93	585159	23.850 ug/mlm
22)	n-C25	24.16	578573	23.660 ug/mlm
23)	n-C26	25.35	575451	23.183 ug/mlm
24)	n-C27	26.50	553151	22.775 ug/mlm
25)	n-C28	27.62	547903	22.660 ug/mlm
26)	n-C29	28.71	550574	22.359 ug/mlm
28)	n-C30	29.76	526543	22.153 ug/mlm
29)	n-C31	30.78	511919	21.905 ug/mlm
30)	n-C32	31.77	484744	21.436 ug/mlm
31)	n-C33	32.76	477901	21.691 ug/mlm
32)	n-C34	33.89	465282	21.378 ug/mlm

Data File: W:\2\DATA\GC10846\GC10846G.D Vial: 98 Acq On : 13 Apr 2007 3:37 pm Sample : CS3 Operator: TJM Inst : GC#1 Misc Multiplr: 1.00

Misc : IntFile : autoint1.e

Quant Time: Apr 14 9:46 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

R.T. Response Conc Units Compound

Mon Apr 16 06:41:42 2007

C10B414.M

GC10846G.D

Quantitation Report

1.00 GC#1 TUM Operator: Multiplr:Vial: Quant Time: Apr 14 9:46 2007 Quant Results File: C10B414.RES C:\GC10846\C10B414.M (Chemstation Integrator) C10 - C35 aliphatic Inst GC10846G.DIFID2B Multiple Level Calibration : W:\2\DATA\GC10846\GC10846G.D Thu Apr 12 09:52:10 2007 3:37 pm ALI_COMP.M 13 Apr 2007 15.09 : autointl.e 69.T 0.00 2.00 4.00 6.00 Quant Method: Response via DataAcq Meth Signal Phase Volume Inj. Last Update Info Data File IntFile Acq On Sample Signal Title 50000 25000-20000-15000-10000 5000 45000 40000 35000 30000 Ó Response Time

Data File : W:\2\DATA\GC10846\GC10846H.D Acq On : 13 Apr 2007 4:37 pm Sample : CS4 Vial: 99 Operator: TJM Inst : GC#1 Misc : IntFile : autoint1.e Multiplr: 1.00

Quant Time: Apr 14 10:10 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase: Signal Info :

****	Compound	R.T.	Response	Conc Units
Tnt	ernal Standards			
1)	n-hexadecane-d34	12.86	391299	20.001 ug/mlm
10)	5a-androstane	18.04	503511	20.001 ug/mlm 20.003 ug/mlm
1.07	Ja-androstane	10.04	505511	20.003 dg/mim
Sys	tem Monitoring Compounds			
4) S	n-dodecane-d26	8.63	689809	42.842 ug/mlm
16) S	n-eicosane-d42	17.45	855492	39.970 ug/mlm
27) S	n-triacontane-d62	29.24	717157	34.241 ug/mlm
Тат	get Compounds			
2)	n-C10	6.26	801196	43.021 ug/mlm
3)	n-C11	7.59	822509	42.859 ug/mlm
5)	n-C12	8.83	863068	42.609 ug/mlm
6)	n-C13	9.99	890194	42.168 ug/mlm
7)	n-C14	11.07	913843	41.584 ug/mlm
8)	n-C15	12.09	932506	40.728 ug/mlm
9)	n-C16	13.11	942836	40.109 ug/mlm
11)	n-C17	14.20	957039	41.317 ug/mlm
12)	Pristane	14.32	1002629	40.673 ug/mlm
13)	n-C18	15.36	965746	40.075 ug/mlm 40.947 ug/mlm
14)	Phytane	15.52	984193	39.425 ug/mlm
15)	n-C19	16.58	959630	40.219 ug/mlm
17)	n-C20	17.85	958468	39.862 ug/mlm
18)	n-C21	19.13	972068	39.162 ug/mlm
19)	n-C21	20.42	922591	38.007 ug/mlm
20)	n-C23	21.69	943884	38.144 ug/mlm
20)	n-C24	22.94	931570	37.655 ug/mlm
22)	n-C24 n-C25	24.17	915988	37.149 ug/mlm
23)	n-C26	25.36	906489	36.219 ug/mlm
23)	n-C27	26.51	867199	35.411 uq/mlm
24) 25)		27.63		
	n-C28		855537	35.092 ug/mlm
26)	n-C29	28.72	855268	34.447 ug/mlm
28)	n-C30	29.77	813083	33.926 ug/mlm
29)	n-C31	30.79	785771	33.346 ug/mlm
30)	n-C32	31.78	739603	32.436 ug/mlm
31)	n-C33	32.77	723681	32.577 ug/mlm
32)	n-C34	33.90	699899	31.892 ug/mlm

Data File: W:\2\DATA\GC10846\GC10846H.D Vial: 99 Acq On : 13 Apr 2007 4:37 pm Sample : CS4 Operator: TJM Inst : GC#1 Misc Multiplr: 1.00

Misc : IntFile : autoint1.e

Quant Time: Apr 14 10:10 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

R.T. Response Conc Units Compound _______

Mon Apr 16 06:41:57 2007

C10B414.M

Quantitation Report

8.00 10.00 12.00 14.00 16.00 18.00 20.00 22.00 24.00 26.00 28.00 30.00 32.00 34.00 36.00 38.00 40.00 42.00 44.00 46.00 48.00 50.00 52.00 GC#1 1.00 TUM Operator: Inst : (Multiplr: Vial: IntFile : autoint1.e Quant Time: Apr 14 10:10 2007 Quant Results File: C10B414.RES C:\GC10846\C10B414.M (Chemstation Integrator) n-C36 GC10846H.D\FID2B Multiple Level Calibration W:\2\DATA\GC10846\GC10846H.D 13 Apr 2007 4:37 pm C10 - C35 aliphatic Thu Apr 12 09:52:10 2007 4:37 pm 61O-4 ALI_COMP.M 12.09 69 Z -900 Quant Method: Response vía DataAcq Meth Volume Inj. Signal Phase 2.00 4.00 Info Last Update Data File 0000 CC10846H.D Acq On Sample Signal Signal Title Misc 80000 Response 70000-50000 40000 20000 10000 -00009 30000 ó

Data File : W:\2\DATA\GC10846\GC10846I.D
Acq On : 13 Apr 2007 5:38 pm
Sample : CS5 Vial: 100 Operator: TJM Inst : GC#1 Misc : IntFile : autoint1.e Multiplr: 1.00

Quant Time: Apr 14 10:28 2007 Quant Results File: C10B414.RES

Quant Method: C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

	Compound	R.T.	Response	Conc Units
	nal Standards			
	n-hexadecane-d34	12.86	399244	20.001 ug/mlm
10)	5a-androstane	18.04	511077	20.003 ug/mlm
	m Monitoring Compounds			
	n-dodecane-d26	8.63	851887	51.855 ug/mlm
16) S 1	n-eicosane-d42	17.45	1045910	48.143 ug/mlm
27) S 1	n-triacontane-d62	29.24	845659	39.778 ug/mlm
Target	t Compounds			
2) 1	n-C10	6.26	991150	52.162 ug/mlm
3) 1	n-C11	7.59	1020023	52.093 ug/mlm
5) 1	n-C12	8.84	1070387	51.793 ug/mlm
6) i	n-C13	9.99	1102577	51.190 ug/mlm
7)	n-C14	11.08	1130250	50.408 ug/mlm
8) 1	n-C15	12.09	1152862	49.350 ug/mlm
9) 1	n-C16	13.12	1162863	48.485 ug/mlm
11) 1	n-C17	14.20	1179039	50.148 ug/ml
12)	Pristane	14.32	1232615	49.263 ug/mlm
13) 1	n-C18	15.36	1184312	49.470 ug/mlm
14)	Phytane	15.53	1209131	47.718 ug/mlm
15) ı	n-C19	16.59	1175615	48.542 ug/mlm
17) ı	n-C20	17.85	1175269	48.155 ug/mlm
18) 1	n-C21	19.14	1190151	47.239 ug/mlm
19) ı	n-C22	20.42	1128786	45.813 ug/mlm
20) 1	n-C23	21.70	1149860	45.780 ug/mlm
21) 1	n-C24	22.95	1132079	45.083 ug/mlm
22)	n-C25	24.17	1110968	44.390 ug/mlm
23) 1	n-C26	25.36	1096425	43.160 ug/mlm
	n-C27	26.51	1047632	42.146 ug/mlm
25) I	n-C28	27.63	1026060	41.463 ug/mlm
	n-C29	28.72	1021478	40.532 ug/mlm
	n-C30	29.77	968992	39.833 ug/mlm
	n-C31	30.79	930890	38.919 ug/mlm
	n-C32	31.78	870519	37.613 ug/mlm
	n-C33	32.78	850247	37.707 ug/mlm
·	n-C34	33.91	818059	36.724 ug/mlm
				<u>.</u>

Data File : W:\2\DATA\GC10846\GC10846I.D Acq On : 13 Apr 2007 5:38 pm Sample : CS5 Vial: 100 Operator: TJM Inst : GC#1 Misc Multiplr: 1.00

Misc : IntFile : autoint1.e

Quant Time: Apr 14 10:28 2007 Quant Results File: C10B414.RES

Quant Method : C:\GC10846\C10B414.M (Chemstation Integrator)

Title : C10 - C35 aliphatic
Last Update : Thu Apr 12 09:52:10 2007
Response via : Initial Calibration

DataAcq Meth : ALI COMP.M

Volume Inj. : Signal Phase : Signal Info :

R.T. Response Conc Units Compound

Quantitation Report

26.00 28.00 30.00 32.00 34.00 36.00 38.00 40.00 42.00 44.00 46.00 48.00 50.00 52.00 GC#1 TUM 100 Vial: Operator: Inst Multiplr: Quant Time: Apr 14 10:28 2007 Quant Results File: C10B414.RES C:\GC10846\C10B414.M (Chemstation Integrator)
C10 - C35 aliphatic n-C21 n-C31 n-C31 n-C31 n-C31 n-C31 GC10846I.D\FID2B Multiple Level Calibration P-C53 W:\2\DATA\GC10846\GC10846I.D Thu Apr 12 09:52:10 2007 0.00 2.00 4.00 6.00 8.00 10.00 12.00 14.00 16.00 18.00 20.00 n-eicosane Carantost Carantost 5:38 pm 14.50.32 ALI COMP.M 13 Apr 2007 autoint1.e 18.8 69°Z Volume Inj. Signal Phase Quant Method Response via DataAcq Meth Info Last Update Data File IntFile Signal 28000 2 Time 0.00 2 2 C108461.D Acg On Sample Misc 00006 80000 70000 40000 30000 20000 10000 Ó Response 00009 50000

Mon Apr 16 06:43:05 2007

C10B414.M

Polycyclic Aromatic Hydrocarbon Initial Calibration Data

PAH ICAL 041507.M

GC/MS 3 (PAH 2002)

Response Factor Report GC/MS Ins

Method : G:\1\METHODS\041507.M (RTE Integrator)
Title : PAH Calibration Table (2002) Last Update : Sat Apr 14 20:04:44 2007

Response via : Initial Calibration

Calibration Files

1 =MS30377D.D 2 =MS30377E.D 3 =MS30377F.D 4 =MS30377G.D 5 =MS30377H.D

		Compound	1	2	3	4	5	Avg	%RSD
1)	I	Fluorene-d10			т с	רוחי			
2)	S	Naphthalene-d8				1.878			
	S T	Decalin				0.377			6.24
3) 4)		C1-Decalin				0.377			6.95
_	un	C2-Decalin				0.377			6.95
5) 6)	un	C3-Decalin				0.377			6.95
7)	un	C4-Decalin				0.377			6.95
-	un								6.95
8)	T	Naphthalene				2.040			6.82
9)	T	2-Methylnaphthalene				1.352			8.43
10)	T	1-Methylnaphthalene				1.271			8.14
11)	T	2,6-Dimethylnaphthale				1.130			9.80
12)	T	1,6,7-Trimethylnaphth				1.092			14.22
13)	un	C2-Naphthalenes				2.040			6.82
14)	un	C3-Naphthalenes				2.040			6.82
15)	un	C4-Naphthalenes				2.040			6.82
16)	T	Benzothiophene				1.743			5.49
17)	un	C1-Benzothiophene				1.743			5.49
18)	un	C2-Benzothiophene				1.743			5.49
19)	un	C3-Benzothiophene				1.743			5.49
20)	S	Acenaphthene-d10				0.893			7.93
21)	T	Biphenyl				1.657			7.23
22)	T	Acenaphthylene				2.006			5.94
23)	T	Acenaphthene				1.152			7.14
24)	T	Dibenzofuran				1.924			11.46
25)	\mathbf{T}	Fluorene				1.335			9.28
26)	un	C1-Fluorenes				1.335			9.28
27)	un	C2-Fluorenes				1.335			9.28
28)	un	C3-Fluorenes	1.141	1.381	1.306	1.335	1.478	1.328	9.28
29)	I	Pyrene-d10			I	STD			
30)	S	Phenanthrene-d10	0.811	0.824	0.835	0.882	0.891	0.849	4.21
31)	T	Pentachlorophenol	0.048	0.040	0.040	0.043	0.049	0.044	10.23
32)	T	Carbazole	0.882	0.917	0.974	1.022	1.146	0.988	10.44
33)	T	Dibenzothiophene	1.058	0.994	1.026	1.026	1.153	1.051	5.83
34)	un	C1-Dibenzothiophene	1.058	0.994	1.026	1.026	1.153	1.051	5.83
35)	un	C2-Dibenzothiophene	1.058	0.994	1.026	1.026	1.153	1.051	5.83
36)	un	C3-Dibenzothiophene	1.058	0.994	1.026	1.026	1.153	1.051	5.83
37)	${f T}$	Phenanthrene	0.987	0.949	0.947	1.009	1.139	1.006	7.81
38)	T	Anthracene				1.181			11.40
39)	T	1-Methylphenanthrene	0.783	0.802	0.820	0.888	1.000	0.859	10.29
40)	un	C1-Phenanthrene/Anthr				1.009			7.81
41)	un	C2-Phenanthrene/Anthr				1.009			7.81
42)	un	C3-Phenanthrene/Anthr	0.987	0.949	0.947	1.009	1.139	1.006	7.81

Response Factor Report GC/MS Ins

Method : G:\1\METHODS\041507.M (RTE Integrator)
Title : PAH Calibration Table (2002) Last Update : Sat Apr 14 20:04:44 2007 Response via : Initial Calibration

Calibration Files

=MS30377D.D 2 =MS30377E.D 3 =MS30377F.D =MS30377G.D 5 =MS30377H.D 1

4

		Compound	1	2	3	4	5	Avg	%RSD
43)	un	C4-Phenanthrene/Anthr	0.987	0 949	0 947	1 009	1.139	1.006	7.81
44)	T	Naphthobenzothiophene							8.17
45)	un	C1-Naphthobenzothioph							8.17
46)	un	C2-Naphthobenzothioph							8.17
47)	un	C3-Naphthobenzothioph							8.17
48)	T	Fluoranthene				1.386			7.07
49)	T	Pyrene	1.304	1.350	1.381	1.506	1.588	1.426	8.24
50)	un	C1-Fluoranthenes/Pyre	1.165	1.274	1.285	1.386	1.384	1.299	7.07
51)	un	C2-Fluoranthenes/Pyre	1.165	1.274	1.285	1.386	1.384	1.299	7.07
52)	un	C3-Fluoranthenes/Pyre	1.165	1.274	1.285	1.386	1.384	1.299	7.07
53)	S	Chrysene-d12				1.258			11.35
54)	T	Benz(a)anthracene				1.405			12.67
55)	T	Chrysene				1.206		1.102	9.11
56)	un	C1-Chrysenes		1.065		1.206		1.102	9.11
57)	un	C2-Chrysenes		1.065		1.206			9.11
58)	un	C3-Chrysenes		1.065		1.206		1.102	9.11
59)	un	C4-Chrysenes	0.952	1.065	1.112	1.206	1.176	1.102	9.11
60)	Ι	Benzo(a)pyrene-d12			I	STD	··· ··· ··· ··· ··· ··· ··· ··· ··· ··		
61)	un	C29-Hopane	0.760	0.646	0.641	0.676	0.671	0.679	7.03
62)	un	18a-Oleanane	0.760	0.646	0.641	0.676	0.671	0.679	7.03
63)	${f T}$	C30-Hopane			0.641		0.671	0.679	7.03
64)	T	Benzo(b)fluoranthene	1.722	1.642	1.717	1.904	1.951	1.787	7.45
65)	T	Benzo(k)fluoranthene			1.747			1.729	5.11
66)	${f T}$	Benzo(e)pyrene				1.899		1.800	4.11
67)	T	Benzo(a)pyrene				1.688		1.529	9.89
68)	\mathbf{T}	Indeno(1,2,3-c,d)pyre		1.057		1.252		1.161	11.59
69)	${f T}$	Dibenzo(a,h)anthracen				1.183		1.077	12.33
70)	un	C1-Dibenzo(a,h)anthra		0.958		1.183		1.077	12.33
71)	un	C2-Dibenzo(a,h)anthra	0.967			1.183		1.077	12.33
72)	un	C3-Dibenzo(a,h)anthra				1.183	1.251		12.33
73)	${ m T}$	Benzo(g,h,i)perylene				1.408		1.298	11.36
74)	S	Perylene-d12				0.859		0.823	8.72
75)	T	Perylene	1.516	1.555	1.588	1.751	1.788	1.640	7.44

Data File : G:\1\DATA\MS30377\MS30377D.D
Acq On : 13 Apr 2007 12:25 pm Vial: 41 Operator: TJM

Sample : Cal Level 1 Inst : GC/MS Ins

Misc Multiplr: 1.00

MS Integration Params: rteint.p

Quant Results File: 041507.RES Quant Time: Apr 14 20:04 19107

Quant Method : C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002) Last Update : Sat Apr 07 22:34:26 2007

Response via : Initial Calibration

DataAcq Meth : PAH-2002

Internal Standards	R.T.	QIon	Response	Conc Units De	v(Min)
 Fluorene-d10 Pyrene-d10 Benzo(a)pyrene-d12 					
System Monitoring Compounds 2) Naphthalene-d8 20) Acenaphthene-d10 30) Phenanthrene-d10	13.01 18.83	136 164	2113m 1006m 1726	18.36	0.03 0.03 0.03
30) Phenanthrene-d10 53) Chrysene-d12 74) Perylene-d12	33.00 37.70	240 264	1726 2151 961	17.68 18.78	0.03
Target Compounds 3) Decalin 4) C1-Decalin 5) C2-Decalin	10.37	138 152 166	528m 0	21.35 ng/ml N.D. d	value
4) C1-Decalin 5) C2-Decalin 6) C3-Decalin 7) C4-Decalin 8) Naphthalene 9) 2-Methylnaphthalene	0.00 0.00 13.10	180 194 128	2295m	N.D. a 17.70	
11) 2,6-Dimethylnaphthalene	17.46	156	1399m 1137m	18.18 18.29	
13) C2-Naphthalenes 14) C3-Naphthalenes 15) C4-Naphthalenes 16) Benzothiophene 17) C1-Benzothiophene	0.00	156 170 184	0 0	N.D. d N.D. d	
16) Benzothiophene 17) C1-Benzothiophene 18) C2-Benzothiophene	13.24 0.00 0.00	134 148 162 176	0 0	N.D. d N.D. d	
23) Acenaphthene	18.95	154	1787m 2206m 1217m	N.D. d 19.94 18.96 17.19	
24) Dibenzofuran25) Fluorene26) C1-Fluorenes	19.57 20.72 0.00	168 166 180	1814m 1377m 0	18.09 ng/ml 17.58 N.D. d	
27) C2-Fluorenes28) C3-Fluorenes31) Pentachlorophenol32) Carbazole	0.00 0.00 23.34 24.79	194 208 266 167	0 0 103m 1878m	N.D. d N.D. d 25.57 ng/ml	
32) Carbazore 33) Dibenzothiophene 34) C1-Dibenzothiophene	23.57	184 198	2254 0	16.62 ng/ml 16.91 N.D. d	

^{(#) =} qualifier out of range (m) = manual integration MS30377D.D 041507.M Mon Apr 16 08:05:57 2007

Vial: 41 Data File : G:\1\DATA\MS30377\MS30377D.D Operator: TJM

Acq On : 13 Apr 2007 12:25 pm Sample : Cal Level 1 Inst : GC/MS Ins

Multiplr: 1.00 Misc

MS Integration Params: rteint.p Quant Time: Apr 14 20:04 19107 Quant Results File: 041507.RES

Quant Method : C:\MS30377\041507.M (RTE Integrator)
Title : PAH Calibration Table (2002) Last Update : Sat Apr 07 22:34:26 2007 Response via : Initial Calibration

DataAcq Meth: PAH-2002

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35)	C2-Dibenzothiophene	0.00	212	0	N.D. d	
	C3-Dibenzothiophene	0.00	226	0	N.D. d	
	Phenanthrene	24.01	178	2104	16.41	
	Anthracene	24.18	178	2060	14.75	
39)	1-Methylphenanthrene	26.13	192	1670	15.26	
40)	C1-Phenanthrene/Anthracene	0.00	192	0	N.D. d	
	C2-Phenanthrene/Anthracene	0.00	206	0	N.D. d	
42)	C3-Phenanthrene/Anthracene	0.00	220	0	N.D. d	
43)	C4-Phenanthrene/Anthracene	0.00	234	0	N.D. d	
44)	Naphthobenzothiophene	32.16	234	1983	16.19	
45)	C1-Naphthobenzothiophene	0.00	248	0	N.D. d	
	C2-Naphthobenzothiophene	0.00	262	0	N.D. d	
47)	C3-Naphthobenzothiophene	0.00	276	0	N.D. d	
48)	Fluoranthene	28.09		2484	18.21	
49)	Pyrene	28.86		2781	16.30	
	C1-Fluoranthenes/Pyrenes	0.00	216	0	N.D. d	
51)	C2-Fluoranthenes/Pyrenes	0.00		0	N.D. d	
52)	C3-Fluoranthenes/Pyrenes	0.00		0	N.D. d	
54)	Benz(a)anthracene	32.96		2393	17.18	
	Chrysene	33.07		2030	17.28	
56)	C1-Chrysenes	0.00		0	N.D. d	
	C2-Chrysenes	0.00		0	N.D. d	
58)	C3-Chrysenes	0.00		0	N.D. d	
	C4-Chrysenes	0.00		0	N.D. d	
	C29-Hopane	0.00		0	N.D. d	
	18a-Oleanane	0.00		0	N.D. d	_
	C30-Hopane	41.74		958	21.92 ng/m	ιL
	Benzo(b) fluoranthene	36.43		2174	19.50	
	Benzo(k)fluoranthene	36.50		2064	22.00	
	Benzo(e)pyrene	37.31		2232	21.36	
	Benzo(a)pyrene	37.49		1748		
	Indeno(1,2,3-c,d)pyrene	41.84		1347m		
	Dibenzo(a,h)anthracene	41.95		1221m		
	C1-Dibenzo(a,h)anthracene	0.00		0	N.D. d	
	C2-Dibenzo(a,h)anthracene	0.00		0	N.D. d	
	C3-Dibenzo(a,h)anthracene	0.00		0	N.D. d	
	Benzo(g,h,i)perylene	43.06		1468	18.31	
75)	Perylene	37.77	252	1916	19.43	

^{(#) =} qualifier out of range (m) = manual integration MS30377D.D 041507.M Mon Apr 16 08:05:58 2007

041507.RES

Quant Results File:

(RTE Integrator)

G:\1\METHODS\041507.M

Last Update

Method Title

Params: rteint.p

MS Integration

Quant Time: Apr 14 20:04 19107

Ins

GC/MS 1.00

Multiplr

Inst

TUM

Vial: Operator:

G:\1\DATA\MS30377\MS30377D.D

Data File

Acq On

Sample

Misc

12:25 pm

13 Apr 2007 Cal Level 1

Data File : G:\1\DATA\MS30377\MS30377E.D Vial: 42 Acq On : 13 Apr 2007 Sample : Cal Level 2 Misc : Operator: TJM 1:28 pm

Inst : GC/MS Ins Multiplr: 1.00 Misc

MS Integration Params: rteint.p

Quant Time: Apr 14 19:58 19107 Quant Results File: 041507.RES

Quant Method: G:\1\METHODS\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002)
Last Update : Sat Apr 07 22:34:26 2007

Response via: Initial Calibration

DataAcq Meth : PAH-2002

Internal Standards	R.T.	QIon	Response	Conc Units De	v(Min)
 Fluorene-d10 Pyrene-d10 Benzo(a)pyrene-d12 					
29) Pyrene-d10	28.80	212	4817m	49.98	0.03
60) Benzo(a)pyrene-d12	37.42	264	2747m	45.61	0.04
System Monitoring Compounds					
	13.01	136	9306m	110.32	0.02
20) Acenaphthene-d10	18.86	164	4400m	107.74	0.05
30) Phenanthrene-d10	23.91	188		92.58	0.03
53) Chrysene-d12 74) Perylene-d12	33.00	240	10238m	92.86	0.04
74) Perylene-d12	37.70	264	4693m	96.02	0.04
Target Compounds				0	value
3) Decalin	10.39	138	1932m		
4) C1-Decalin	0.00	152	0	N.D. d	
4) C1-Decalin 5) C2-Decalin 6) C3-Decalin 7) C4-Decalin 8) Naphthalene 9) 2-Methylnaphthalene	0.00	166	0	N.D. d	
6) C3-Decalin	0.00	180	0		
7) C4-Decalin	0.00	194	0		
8) Naphthalene	13.09	128		105.92	
9) 2-Methylnaphthalene	15.34	142		112.81	
10) 1-Methylnaphthalene	15.65	142	6336m	110.44	
11) 2,6-Dimethylnaphthalene	17.45		5314m	114.67	
12) 1.6.7-Trimethylnaphthalene	20.29	170	5063m		
13) C2-Naphthalenes 14) C3-Naphthalenes 15) C4-Naphthalenes 16) Benzothiophene	0.00	156	0	N.D. d	
14) C3-Naphthalenes	0.00	170	0	N.D. d	
15) C4-Naphthalenes	0.00	184	0	N.D. d	
16) Benzothiophene	13.23	134	8480m	109.00 ng/ml	
1// Ci-benzoenrophene	0.00	T. 4 O	0	N.D. d	
18) C2-Benzothiophene	0.00	162	0	N.D. d	
19) C3-Benzothiophene	0.00	176	0		
21) Biphenyl	16.92	154	7963m		
22) Acenaphthylene	18.35	152	9618m		
23) Acenaphthene	18.94	154	5503m		
24) Dibenzofuran 25) Fluorene 26) C1-Fluorenes	19.56	168			
25) Fluorene	20.74	166	6210m		
•			0	N.D. d	
27) C2-Fluorenes	0.00	194	0	N.D. d	
28) C3-Fluorenes	0.00	208	0	N.D. d	
31) Pentachlorophenol	23.30	266	382m	104.66 ng/ml	
32) Carbazole	24.75	167	8849m	86.42 ng/ml	
33) Dibenzothiophene	23.57	184	9596m	79.46	
34) C1-Dibenzothiophene	0.00	198	0	N.D. d	

^{(#) =} qualifier out of range (m) = manual integration MS30377E.D 041507.M Mon Apr 16 08:06:14 2007

Data File : G:\1\DATA\MS30377\MS30377E.D
Acq On : 13 Apr 2007 1:28 pm Vial: 42 Operator: TJM

Sample : Cal Level 2 Inst : GC/MS Ins

Multiplr: 1.00 Misc

MS Integration Params: rteint.p Quant Time: Apr 14 19:58 19107

Quant Results File: 041507.RES

Quant Method: G:\1\METHODS\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002)
Last Update : Sat Apr 07 22:34:26 2007

Response via: Initial Calibration

DataAcq Meth : PAH-2002

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35)	C2-Dibenzothiophene	0.00	212	0	N.D. d	
		0.00	226	0	N.D. d	
37)	C3-Dibenzothiophene Phenanthrene	24.01	178		78.92	
38)	Anthracene	24.18	178	10119m	79.99	
39)	Anthracene 1-Methylphenanthrene C1-Phenanthrene/Anthracene	24.18 26.13	192	7742m	78.11	
40)	C1-Phenanthrene/Anthracene	0.00	192	0	N.D. d	
	C2-Phenanthrene/Anthracene		206	0	N.D. d	
	C3-Phenanthrene/Anthracene	0.00	220	0	N.D. d	
	C4-Phenanthrene/Anthracene	0.00	234	0	N.D. d	
	Naphthobenzothiophene	32.16		8952m	80.67	
	Cl-Naphthobenzothiophene	0.00	248	0	N.D. d	
	C2-Naphthobenzothiophene	0.00	262	0	N.D. d	
	C3-Naphthobenzothiophene	0.00	276	0	N.D. d	
	Fluoranthene	28.09	202	12305m	99.57	
	Pyrene	28.86	202	13035m	84.34	
	C1-Fluoranthenes/Pyrenes	0.00	216	0	N.D. d	
	C2-Fluoranthenes/Pyrenes	0.00	216 230	0	N.D. d	
	C3-Fluoranthenes/Pyrenes	32.93			N.D.	
	Benz(a)anthracene	32.96	228	11821m	93.68	
55)		33.07	228	10287m	96.67	
	C1-Chrysenes	0.00	242	0	N.D. d	
57)	C2-Chrysenes	0.00 0.00 0.00	256			
	C3-Chrysenes	0.00	270	0	N.D. d N.D. d	
	C4-Chrysenes	0.00	284	0	N.D. d	
	**			0	N.D. d	
62)	18a-Oleanane	0 00	191	0	N.D. d	
63)	C30-Hopane	41.74	191	3892m	93.25 ng/m	ıl
64)	Benzo(b) fluoranthene Benzo(k) fluoranthene Benzo(e) pyrene	36.43	252	9900m		
65)	Benzo(k)fluoranthene	36.50	252	9951m	111.04	
66)	Benzo(e)pyrene	37.31	252	10265m	102.83	
67)	Benzo(a)pyrene	37.49	252	8386m	96.02	
68)		41.84	276	6381m	94.65	
69)	Dibenzo(a,h)anthracene	41.94	278	5776m	91.35	
	C1-Dibenzo(a,h)anthracene	0.00	292	0	N.D. d	
	C2-Dibenzo(a,h)anthracene		306		N.D. d	
	C3-Dibenzo(a,h)anthracene	0.00	320			
	Benzo(g,h,i)perylene	43.06	276	7037m	91.89	
75)	Perylene	37.77	252	9381m	99.58	

^{(#) =} qualifier out of range (m) = manual integration MS30377E.D 041507.M Mon Apr 16 08:06:15 2007

Quant Results File: 041507.RES

G:\1\METHODS\041507.M (RTE Integrator)
PAH Calibration Table (2002)

Sat Apr 14 20:04:44 2007

Calibration

Initial

Response via

Last Update

Method Title

Ins

GC/MS

TUM

Operator:

G:\1\DATA\MS30377\MS30377E.D 13 Apr 2007 1:28 pm Cal Level 2

Data File

Acq On Sample

Misc

MS Integration Params: rteint.p Quant Time: Apr 14 19:58 19107

Vial:

1.00

Multiplr:

Inst

Multiplr: 1.00

Vial: 43 Data File : G:\1\DATA\MS30377\MS30377F.D Acq On : 13 Apr 2007 2:31 pm Operator: TJM

Sample : Cal Level 3
Misc : Inst : GC/MS Ins

Misc MS Integration Params: rteint.p

Quant Time: Apr 14 19:59 19107 Quant Results File: 041507.RES

Quant Method: C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002)
Last Update : Sat Apr 07 22:34:26 2007
Response via : Initial Calibration

DataAcq Meth: PAH-2002

1) Fluorene-d10	Internal Standards		QIon	Response	Conc Units Dev	/(Min)
29 Pyrene-d10	1) Fluorene-d10	20.63	176	2427m	51.08 ng/ml	0.03
System Monitoring Compounds 2) Naphthalene-d8	29) Pyrene-d10	28.79				
2) Naphthalene-d8	60) Benzo(a)pyrene-d12	37.42				
2) Naphthalene-d8	System Monitoring Compounds					
200 Acenaphthene-d10		13.01	136	22285m	249.50	0.03
53) Chrysene-d12 33.00 240 25782m 239.12 0.04 74) Perylene-d12 37.70 264 11778m 240.80 0.04 Target Compounds Qvalue 3) Decalin 10.39 138 4577m 234.47 ng/ml 4) C1-Decalin 0.00 166 0 N.D. d 5) C2-Decalin 0.00 166 0 N.D. d 6) C3-Decalin 0.00 180 0 N.D. d 7) C4-Decalin 0.00 180 0 N.D. d 8) Naphthalene 13.09 128 24356m 238.00 9) 2-Methylnaphthalene 15.34 142 15813m 255.17 10) 1-Methylnaphthalene 15.65 142 13928m 229.27 11) 2,6-Dimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 0.00 156 0 N.D. d 13) C2-Naphthalenes 0.00 156 0 N.D. d 14) C3-Naphthal		18.83	164	10491m		
53) Chrysene-d12 33.00 240 25782m 239.12 0.04 74) Perylene-d12 37.70 264 11778m 240.80 0.04 Target Compounds Qvalue 3) Decalin 10.39 138 4577m 234.47 ng/ml 4) C1-Decalin 0.00 166 0 N.D. d 5) C2-Decalin 0.00 166 0 N.D. d 6) C3-Decalin 0.00 180 0 N.D. d 7) C4-Decalin 0.00 180 0 N.D. d 8) Naphthalene 13.09 128 24356m 238.00 9) 2-Methylnaphthalene 15.34 142 15813m 255.17 10) 1-Methylnaphthalene 15.65 142 13928m 229.27 11) 2,6-Dimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 0.00 156 0 N.D. d 13) C2-Naphthalenes 0.00 156 0 N.D. d 14) C3-Naphthal	30) Phenanthrene-d10	23.91	188	19681	234.60	0.03
Target Compounds 3) Decalin 10.39 138 4577m 234.47 ng/ml 4) C1-Decalin 0.00 152 0 N.D. d 5) C2-Decalin 0.00 166 0 N.D. d 6) C3-Decalin 0.00 180 0 N.D. d 7) C4-Decalin 0.00 194 0 N.D. d 7) C4-Decalin 0.00 194 0 N.D. d 8) Naphthalene 13.09 128 24356m 238.00 9) 2-Methylnaphthalene 15.34 142 15813m 255.17 10) 1-Methylnaphthalene 15.65 142 13928m 229.27 11) 2,6-Dimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 13.09 120 10355m 26.60 13) C2-Naphthalenes 0.00 156 0 N.D. d 14) C3-Naphthalenes 0.00 156 0 N.D. d 15) C4-Naphthalenes 0.00 170 0 N.D. d 15) C4-Naphthalenes 0.00 184 0 N.D. d 16) Benzothiophene 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 0.00 148 0 N.D. d 18) C2-Benzothiophene 0.00 162 0 N.D. d 19) C3-Benzothiophene 0.00 162 0 N.D. d 19) C3-Benzothiophene 18.35 152 23701m 258.12 23) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 277.49 22) Acenaphthene 18.94 154 13464m 277.49 22) Acenaphthene 18.95 Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 194 0 N.D. d 29) C3-Fluorenes 0.00 194 0 N.D. d 20) N.D. d 21) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 23) Dibenzothiophene 23.57 184 24212m 204.99	53) Chrysene-d12	33.00	240		239.12	0.04
3) Decalin 4) C1-Decalin 6) C2-Decalin 7) C2-Decalin 8) Naphthalene 13.09 128 24356m 9) 2-Methylnaphthalene 15.34 142 15813m 255.17 10) 1-Methylnaphthalene 15.65 142 13928m 229.27 11) 2,6-Dimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 0 N.D. d 13) C2-Naphthalenes 10.00 156 0 N.D. d 14) C3-Naphthalenes 10.00 156 0 N.D. d 15) C4-Naphthalenes 10.00 170 0 N.D. d 15) C4-Naphthalenes 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 13.25 152 23701m 258.12 23 Acenaphthylene 18.35 152 23701m 258.12 23 Acenaphthene 18.35 152 23701m 258.12 23 Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 20.00 194 0 N.D. d 27) C2-Fluorenes 20.00 194 0 N.D. d 28) C3-Fluorenes 20.00 194 0 N.D. d 29 C2-Fluorenes 20.00 194 0 N.D. d 20 C2-Fluorenes 20.00 194 0 N.D. d 21) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 23) Dibenzothiophene 23.57 184 24212m 204.99	74) Perylene-d12	37.70	264		240.80	0.04
3) Decalin 4) C1-Decalin 6) C2-Decalin 7) C2-Decalin 8) Naphthalene 13.09 128 24356m 9) 2-Methylnaphthalene 15.34 142 15813m 255.17 10) 1-Methylnaphthalene 15.65 142 13928m 229.27 11) 2,6-Dimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 0 N.D. d 13) C2-Naphthalenes 10.00 156 0 N.D. d 14) C3-Naphthalenes 10.00 156 0 N.D. d 15) C4-Naphthalenes 10.00 170 0 N.D. d 15) C4-Naphthalenes 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 18.35 152 23701m 258.12 23 Acenaphthylene 18.35 152 23701m 258.12 23 Acenaphthene 18.35 152 23701m 258.12 23 Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 20.00 194 0 N.D. d 27) C2-Fluorenes 20.01 180 0 N.D. d 28) C3-Fluorenes 20.00 194 0 N.D. d 29) C3-Banzothiophenol 21.30 266 934m 261.65 ng/ml 22.31 Pentachlorophenol 23.30 266 934m 261.65 ng/ml 23.30 Dibenzothiophene 23.57 184 24212m 204.99	Target Compounds				r <u>O</u>	value
5) C2-Decalin		10.39	138	4577m		
6) C3-Decalin 7) C4-Decalin 8) Naphthalene 9) 2-Methylnaphthalene 13.09 128 24356m 238.00 9) 2-Methylnaphthalene 15.34 142 15813m 255.17 10) 1-Methylnaphthalene 15.65 142 13928m 229.27 11) 2,6-Dimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 17.45 156 0 N.D. d 13) C2-Naphthalenes 0.00 156 0 N.D. d 14) C3-Naphthalenes 0.00 170 0 N.D. d 15) C4-Naphthalenes 0.00 184 0 N.D. d 15) C4-Naphthalenes 0.00 184 0 N.D. d 16) Benzothiophene 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 0.00 148 0 N.D. d 18) C2-Benzothiophene 0.00 162 0 N.D. d 19) C3-Benzothiophene 0.00 162 0 N.D. d 19) C3-Benzothiophene 18.35 152 23701m 258.12 23) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 194 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 194 0 N.D. d 29) C3-Benzothiophene 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene	4) C1-Decalin	0.00	152	0	-	
7) C4-Decalin 8) Naphthalene 13.09 128 24356m 238.00 9) 2-Methylnaphthalene 15.34 142 158813m 255.17 10) 1-Methylnaphthalene 15.65 142 13928m 229.27 11) 2,6-Dimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalene 20.29 170 10355m 226.60 13) C2-Naphthalenes 0.00 156 0 N.D. d 14) C3-Naphthalenes 0.00 170 0 N.D. d 15) C4-Naphthalenes 0.00 170 0 N.D. d 15) C4-Naphthalenes 0.00 184 0 N.D. d 16) Benzothiophene 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 0.00 148 0 N.D. d 18) C2-Benzothiophene 0.00 162 0 N.D. d 19) C3-Benzothiophene 0.00 176 0 N.D. d 19) C3-Benzothiophene 16.92 154 19624m 277.49 22) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 23.57 184 24212m 204.99	5) C2-Decalin	0.00	166	0	N.D. d	
8) Naphthalene	6) C3-Decalin	0.00	180	0	N.D. d	
8) Naphthalene				0	N.D. d	
10) 1-Methylnaphthalene	8) Naphthalene	13.09	128	24356m	238.00	
10) 1-Methylnaphthalene	9) 2-Methylnaphthalene	15.34	142	15813m	255.17	
11) 2,6-Dimethylnaphthalene 17.45 156 12835m 261.55 12) 1,6,7-Trimethylnaphthalenes 20.29 170 10355m 226.60 13) C2-Naphthalenes 0.00 156 0 N.D. d 14) C3-Naphthalenes 0.00 170 0 N.D. d 15) C4-Naphthalenes 0.00 184 0 N.D. d 16) Benzothiophene 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 0.00 148 0 N.D. d 18) C2-Benzothiophene 0.00 162 0 N.D. d 19) C3-Benzothiophene 0.00 176 0 N.D. d 19) C3-Benzothiophene 0.00 176 0 N.D. d 21) Biphenyl 16.92 154 19624m 277.49 22) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 0.00 180 0 N.D. d	10) 1-Methylnaphthalene	15.65	142	13928m	229.27	
13) C2-Naphthalenes	11) 2,6-Dimethylnaphthalene	17.45	156	12835m	261.55	
13) C2-Naphthalenes	12) 1,6,7-Trimethylnaphthalene	20.29	170	10355m	226.60	
14) C3-Naphthalenes			156	0	N.D. d	
16) Benzothiophene 13.23 134 20939m 254.18 ng/ml 17) C1-Benzothiophene 0.00 148 0 N.D. d 18) C2-Benzothiophene 0.00 162 0 N.D. d 19) C3-Benzothiophene 0.00 176 0 N.D. d 21) Biphenyl 16.92 154 19624m 277.49 22) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99 <td>14) C3-Naphthalenes</td> <td>0.00</td> <td>170</td> <td>0</td> <td>N.D. d</td> <td></td>	14) C3-Naphthalenes	0.00	170	0	N.D. d	
17) C1-Benzothiophene	15) C4-Naphthalenes	0.00	184	0	N.D. d	
18) C2-Benzothiophene 0.00 162 0 N.D. d 19) C3-Benzothiophene 0.00 176 0 N.D. d 21) Biphenyl 16.92 154 19624m 277.49 22) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99				20939m	254.18 ng/ml	
19) C3-Benzothiophene 0.00 176 0 N.D. d 21) Biphenyl 16.92 154 19624m 277.49 22) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	17) C1-Benzothiophene	0.00	148	0	N.D. d	
21) Biphenyl 16.92 154 19624m 277.49 22) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	18) C2-Benzothiophene	0.00	162	0	N.D. d	
22) Acenaphthylene 18.35 152 23701m 258.12 23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	19) C3-Benzothiophene	0.00		0	N.D. d	
23) Acenaphthene 18.94 154 13464m 240.89 24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	21) Biphenyl	16.92	154	19624m	277.49	
24) Dibenzofuran 19.56 168 21406m 270.41 ng/ml 25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99						
25) Fluorene 20.74 166 15547m 251.41 26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99						
26) C1-Fluorenes 0.00 180 0 N.D. d 27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	24) Dibenzofuran	19.56				
27) C2-Fluorenes 0.00 194 0 N.D. d 28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	25) Fluorene					
28) C3-Fluorenes 0.00 208 0 N.D. d 31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	26) C1-Fluorenes					
31) Pentachlorophenol 23.30 266 934m 261.65 ng/ml 32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	27) C2-Fluorenes	0.00	194	0	N.D. d	
32) Carbazole 24.75 167 22976m 229.45 ng/ml 33) Dibenzothiophene 23.57 184 24212m 204.99	28) C3-Fluorenes	0.00		0		
33) Dibenzothiophene 23.57 184 24212m 204.99	31) Pentachlorophenol				- .	
				22976m		
24) C1 Dibanathianhana 0 00 100 0 ND d						
34) CI-DIDenzoumiophene 0.00 198 0 N.D. d	34) C1-Dibenzothiophene	0.00	198	0	N.D. d	

^{(#) =} qualifier out of range (m) = manual integration MS30377F.D 041507.M Mon Apr 16 08:06:31 2007

Vial: 43 Operator: TJM

Data File : G:\1\DATA\MS30377\MS30377F.D
Acq On : 13 Apr 2007 2:31 pm
Sample : Cal Level 3 Inst : GC/MS Ins Misc Multiplr: 1.00

MS Integration Params: rteint.p

Quant Time: Apr 14 19:59 19107 Quant Results File: 041507.RES

Quant Method : C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002)
Last Update : Sat Apr 07 22:34:26 2007
Response via : Initial Calibration

DataAcq Meth : PAH-2002

ha	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35)	C2-Dibenzothiophene	0.00	212	0	N.D. d	
36)	C3-Dibenzothiophene Phenanthrene Anthracene 1-Methylphenanthrene	0.00	226	0	N.D. d	
37)	Phenanthrene	24.01	178	22363m 25658m 19359m 0		
38)	Anthracene	24.18	178	25658m		
39)	1-Methylphenanthrene	26.13	192	19359m	199.71	
40)	C1-Phenanthrene/Anthracene	0.00	192	0	N.D. d	
	C2-Phenanthrene/Anthracene	0.00	206	0	N.D. d	
	C3-Phenanthrene/Anthracene	0.00	220	0	N.D. d	
	C4 Dhonanthrono/Anthrogono	0 00	224	0	N.D. d	
	Naphthobenzothiophene	32.16 0.00 0.00	234	21666m	199.64	
	Cl-Naphthobenzothiophene	0.00	248	0	N.D. d	
	C2-Naphthobenzothiophene	0.00	262	0	N.D. d	
	C3-Naphthobenzothiophene	0.00	276	0	N.D. d	
	Fluoranthene	28.09	202	30359m	251.19	
	Pyrene	28.86 0.00 0.00	202	32614m	215.76	
	C1-Fluoranthenes/Pyrenes	0.00	216	0	N.D. d	
	C2-Fluoranthenes/Pyrenes	0.00	230	0	N.D. d	
52)	C3-Fluoranthenes/Pyrenes	0.00	244	0	N.D. d	
E4\	Pong (a) anthracene	32.96	228	30910m	250.46	
55)	Chrysene C1-Chrysenes C2-Chrysenes C3-Chrysenes C4-Chrysenes C4-Chrysenes C29-Hopane 18a-Oleanane C30-Hopane	33.07	228	26268m		
56)	C1-Chrysenes	0.00	242	0	N.D. d	
57)	C2-Chrysenes	0.00	242 256	0	N.D. d	
58)	C3-Chrysenes	0.00	270	0	N.D. d	
59)	C4-Chrysenes	0.00	284	0	N.D. d	
61)	C29-Hopane	0.00	191	0	N.D. d	
62)	18a-Oleanane	0.00	191	0	N.D. d	
63)	C30-Hopane	41.74	191	0 9656 25904 26394	231.19 ng/m	ıl
64)	Benzo(b) fluoranthene	36.43	252	25904	243.14	
65)	Benzo(b) fluoranthene Benzo(k) fluoranthene Benzo(e) pyrene Benzo(a) pyrene Indeno(1,2,3-c,d) pyrene Dibenzo(a,h) anthracene C1-Dibenzo(a,h) anthracene	36.50	252	26394	294.32	
66)	Benzo(e)pyrene	37.31	252	27283	273.10	
67)	Benzo(a)pyrene	37.49	252	22576	258.31	
68)	Indeno(1,2,3-c,d)pyrene	41.84	276	16253		
69)	Dibenzo(a,h)anthracene	41.94	278	15476m	244.58	
70)	C1-Dibenzo(a,h)anthracene	0.00	292	0	N.D. d	
71)	C2-Dibenzo(a,n)anthracene	0.00	306	Ü	N.D. d	
72)	C3-Dibenzo(a,h)anthracene		320		N.D. d	
73)	Benzo(g,h,i)perylene	43.06	276	19049m	248.57	
75)	Perylene	37.77	252	23975m	254.31	

^{(#) =} qualifier out of range (m) = manual integration MS30377F.D 041507.M Mon Apr 16 08:06:32 2007

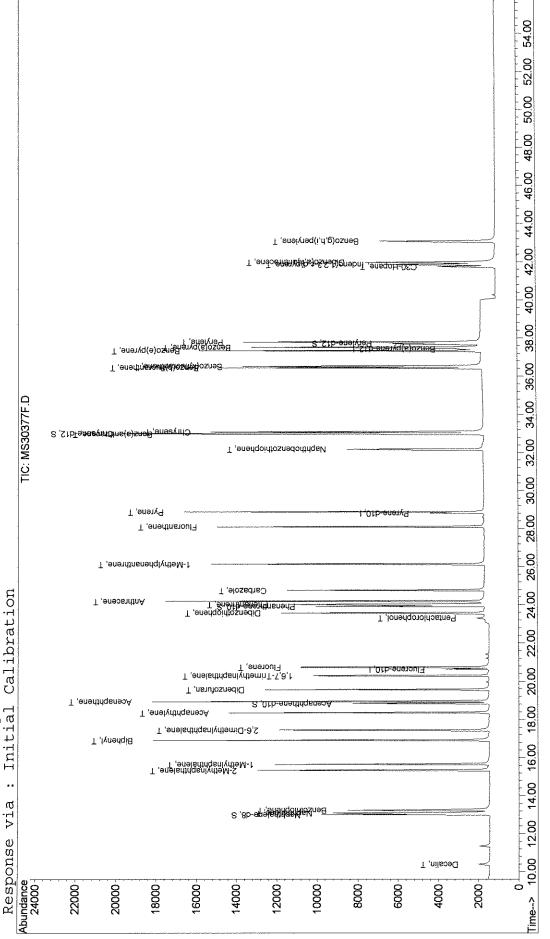
Quantitation Report

Quant Results File: 041507.RES Multiplr:Vial: Operator: Inst G:\1\DATA\MS30377\MS30377F.D 13 Apr 2007 2:31 pm Cal Level 3 MS Integration Params: rteint.p Quant Time: Apr 14 19:59 19107 Data File Acq On Sample Misc

GC/MS ΜĊΙ

Ins

G:\1\METHODS\041507.M (RTE Integrator) (2002)Sat Apr 14 20:04:44 2007 Initial Calibration PAH Calibration Table Last Update Method Title



08:06:33 2007

Quantitation Report (QT Reviewed)

Vial: 44 Data File : G:\1\DATA\MS30377\MS30377G.D Operator: TJM 3:34 pm

Acq On : 13 Apr 2007 Sample : Cal Level 4 Inst : GC/MS Ins

Misc Multiplr: 1.00

MS Integration Params: rteint.p

Quant Results File: 041507.RES Quant Time: Apr 14 20:03 19107

Quant Method: C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002)
Last Update : Sat Apr 07 22:34:26 2007

Response via: Initial Calibration

DataAcq Meth : PAH-2002

Internal Standards	R.T.	QIon	Response	Conc Units Dev	(Min)
1) Fluorene-d10 29) Pyrene-d10 60) Benzo(a)pyrene-d12	20.63 28.80 37.42	176 212 264	2331m 4362m 2667m	51.08 ng/ml 49.98 45.61	0.02 0.03 0.04
System Monitoring Compounds 2) Naphthalene-d8 20) Acenaphthene-d10 30) Phenanthrene-d10 53) Chrysene-d12 74) Perylene-d12	13.01 18.86 23.91 33.00 37.70	136 164 188 240 264	42856m 20383m 38505m 54883m 25119m	499.57 490.77 495.71 549.75 529.34	0.02 0.05 0.03 0.04 0.04
Target Compounds 3) Decalin 4) C1-Decalin 5) C2-Decalin 6) C3-Decalin 7) C4-Decalin 8) Naphthalene 9) 2-Methylnaphthalene 10) 1-Methylnaphthalene 11) 2,6-Dimethylnaphthalene 12) 1,6,7-Trimethylnaphthalene 13) C2-Naphthalenes 14) C3-Naphthalenes 15) C4-Naphthalenes 16) Benzothiophene 17) C1-Benzothiophene 18) C2-Benzothiophene 19) C3-Benzothiophene 19) C3-Benzothiophene 21) Biphenyl 22) Acenaphthylene 23) Acenaphthene 24) Dibenzofuran 25) Fluorene 26) C1-Fluorenes 27) C2-Fluorenes 28) C3-Fluorenes 31) Pentachlorophenol 32) Carbazole	0.00 0.00 0.00 13.09 15.34 15.65 17.45 20.29 0.00 0.00 0.00	152 166 180 194 128 142 156 170 156 170 184 148 162 176 154 154 168 166 180 194 208 266	0 0 0 46653m 30943m 29039m 25845m 24953m 0 0	459.50 ng/ml N.D. d N.D. d N.D. d N.D. d 474.65 519.87 497.71 548.36 568.55 N.D. d N.D. d N.D. d N.D. d N.D. d N.D. d 503.69 ng/ml N.D. d N.D. d 557.68 520.16 490.30 578.27 ng/ml 513.83 N.D. d N.D. d N.D. d N.D. d N.D. d	<i>r</i> alue
33) Dibenzothiophene 34) C1-Dibenzothiophene	23.57		44830m 0	409.93 N.D. d	

^{(#) =} qualifier out of range (m) = manual integration MS30377G.D 041507.M Mon Apr 16 08:06:47 2007

Quantitation Report (QT Reviewed)

Vial: 44 Data File: G:\1\DATA\MS30377\MS30377G.D Operator: TJM 3:34 pm

Acq On : 13 Apr 2007
Sample : Cal Level 4 Inst : GC/MS Ins

Multiplr: 1.00 Misc

MS Integration Params: rteint.p Quant Results File: 041507.RES Quant Time: Apr 14 20:03 19107

Quant Method : C:\MS30377\041507.M (RTE Integrator)
Title : PAH Calibration Table (2002) Last Update : Sat Apr 07 22:34:26 2007

Response via : Initial Calibration

DataAcq Meth: PAH-2002

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
35)	C2-Dibenzothiophene	0.00	212	0	N.D. d	
36)	C3-Dibenzothiophene		226		N.D. d	
37)	C3-Dibenzothiophene Phenanthrene	24.01	178	44125m	419.43	
38)	Anthracene	24.18	178	51595m	450.42	
	1-Methylphenanthrene	26.13	178 192 192	38834m	432.67	
40)	C1-Phenanthrene/Anthracene	0.00	192	0	N.D. d	
41)	C2-Phenanthrene/Anthracene	0.00	206	0	N.D. d	
	C3-Phenanthrene/Anthracene	0.00	220	0	N.D. d	
43)	The state of the s	0.00	234	0	N.D. d	
	Naphthobenzothiophene	32.16	234	45498m	452.79	
	C1-Naphthobenzothiophene	0.00	248	0	N.D. d	
	C2-Naphthobenzothiophene	0.00	262	0	N.D. d	
	C3-Naphthobenzothiophene	0.00	276	0	N.D. d	
	Fluoranthene	28.09	202	60628m	541.76	
-	Pyrene	28.86	202	65862m	470.58	
	C1-Fluoranthenes/Pyrenes	0.00	216	0	N.D. d	
51)	C2-Fluoranthenes/Pyrenes	0.00	230	0	N.D. d	
52)	C3-Fluoranthenes/Pyrenes	0.00	244	0	N.D. d	
	Benz(a) anthracene	32.96	228	61436m	537.64	
	Chrysene	33.07	228	52749m	547.41	
		0.00	242	0	N.D. d	
	C2-Chrysenes	0.00	256	0	N.D. d	
		0.00	270	0	N.D. d	
	C3-Chrysenes C4-Chrysenes C29-Hopane 18a-Oleanane C30-Hopane	0.00	284	0	N.D. d	
	C29-Hopane	0.00	191	0	N.D. d	
	18a-Oleanane	0.00	191	0	N.D. d	
63)	C30-Hopane				487.92 ng/r	nl
64)	Benzo (b) fluoranthene	36.43	252	55736m	539.24	
65)	Benzo(k) fluoranthene Benzo(e) pyrene Benzo(a) pyrene	36.50	252			
66)	Benzo(e)pyrene	37.31	252		574.28	
67)	Benzo(a) pyrene	37.49	252	49425m	582.89	
	Indeno(1,2,3-c,d)pyrene	41.84		36686m	560.49	
69)	Dibenzo(a,h)anthracene	41.95	278	34626m	564.06	
	C1-Dibenzo(a,h)anthracene	0.00	292	0	N.D. d	
	C2-Dibenzo(a,h)anthracene	0.00	306	0	N.D. d	
	C3-Dibenzo(a,h)anthracene		320		N.D. d	
	Benzo(g,h,i)perylene	43.06	276	41233m	554.60	
	Perylene	37.77		51302m	560.91	
,	-					

G:\1\DATA\MS30377\MS30377G.D 13 Apr 2007 3:34 pm Cal Level 4 Data File Acq On Sample Misc

MS Integration Params: rteint.p Quant Time: Apr 14 20:03 19107

Ins GC/MS Inst

ΣŊL

Vial: Operator: Multiplr:

Quant Results File: 041507.RES

G:\1\METHODS\041507.M (RTE Integrator)

(2002)Sat Apr 14 20:04:44 2007 Initial Calibration PAH Calibration Table Response via Last Update Method Title

54.00 52.00 20.00 48.00 46.00 44.00 Denzo(g,h,i)perylene, T 42.00 Indeno(128868499)(19881117 acene, T 40.00 38.00 Henzo(a)ozneRq(a)oyrene, T 36.00 Senzo(k)fluora Barara (p)fluoranthene, T TIC: MS30377G.D 34.00 T ,enelogenbran(skyne) ,enekynt 32.00 T ,enertqointosnedorthqeM 30,00 Pyrene, T 1,0tb-елетуЧ 28,00 Fluoranthene, T 26.00 1-Methylphenanthrene, T Carbazole, T 22.00 24.00 Anthracene, T Dibenzothiophene, T Pentachlorophenol, T 20 00 T leneneul F T ,enelshindsnlydteminT-7,8,1 Acenaphthene, T 10.00 12.00 14.00 16.00 18.00 Acenaphthylene, T T ,enelerthylnaphthalene, T Biphenyl, T T ,enelentinaphthalene, T ,enelene, T ,enelene, T ,enelentinaphthalene, T T , e Rei Bhith a gent de la Chinos na El Decalin, T 20002 15000 10000 5000 45000 40000 35000 30000 25000 Ó Time-->

Vial: 45 Data File: G:\1\DATA\MS30377\MS30377H.D Acq On : 13 Apr 2007 Sample : Cal Level 5 Operator: TJM 4:37 pm

Inst : GC/MS Ins

Multiplr: 1.00 Misc

MS Integration Params: rteint.p

Quant Time: Apr 14 20:03 19107 Quant Results File: 041507.RES

Quant Method : C:\MS30377\041507.M (RTE Integrator)

Title : PAH Calibration Table (2002) Last Update : Sat Apr 07 22:34:26 2007 Response via : Initial Calibration

DataAcq Meth: PAH-2002

Internal Standards	R.T.	QIon	Response	Conc Units D	ev(Min)
1) Fluorene-d10 29) Pyrene-d10 60) Benzo(a)pyrene-d12					
29) Pyrene-d10	28.80	212	4129m	49.98	0.03
60) Benzo(a)pyrene-d12	37.42	264	2785m	45.61	0.04
System Monitoring Compounds					
2) Naphthalene-d8 20) Acenaphthene-d10 30) Phenanthrene-d10	13.01	136	83349m	1048.99 1111.81 1000.99 1146.52	0.03
20) Acenaphthene-d10	18.86	164	42769m	1111.81	0.05
30) Phenanthrene-d10	23.91	188	73600m	1000.99	0.03
53) Chrysene-d12	33.00	240	108347m	1146.52	0.04
2) Naphthalene-d8 20) Acenaphthene-d10 30) Phenanthrene-d10 53) Chrysene-d12 74) Perylene-d12	37.70	264	56926m	1148.79	0.04
Target Compounds					Qvalue
Target Compounds 3) Decalin 4) C1-Decalin 5) C2-Decalin 6) C3-Decalin 7) C4-Decalin 8) Naphthalene 9) 2-Methylnaphthalene 10) 1-Methylnaphthalene 11) 2,6-Dimethylnaphthalene 12) 1,6,7-Trimethylnaphthalene 13) C2-Naphthalenes 14) C3-Naphthalenes	10.39	138	16456m	947.65 ng/ml	
4) C1-Decalin	0.00	152	0	N.D. d	
5) C2-Decalin	0.00	166	0	N.D. d	
6) C3-Decalin	0.00	180	0	N.D. d	
7) C4-Decalin	0.00	194	0	N.D. d	
8) Naphthalene	13.09	128	91937m	1009.88	
9) 2-Methylnaphthalene	15.34	142	62685m	1137.07	
10) 1-Methylnaphthalene	15.65	142	55086m	1019.35	
11) 2,6-Dimethylnaphthalene	17.45	156	51796m	1186.53	
12) 1,6,7-Trimethylnaphthalene	20.29	170	49694m	1222.47	
13) C2-Naphthalenes 14) C3-Naphthalenes 15) C4-Naphthalenes 16) Benzothiophene	0.00	156	0	N.D. d	
14) C3-Naphthalenes	0.00	170	0	N.D. d	
	0.00	184	0		
16) Benzothiophene	13.23	134	77082m	1051.85 ng/ml	
16) Benzothiophene 17) C1-Benzothiophene 18) C2-Benzothiophene	0.00	Tt ()	0	N.D. d	
18) C2-Benzothiophene	(1) (1) (1)	1 (-1)	0	N.D. d	
19) C3-Benzothiophene	0.00	176	0	N.D. d	
18) C2-Benzothrophene 19) C3-Benzothrophene 21) Biphenyl 22) Acenaphthylene 23) Acenaphthene 24) Dibenzofuran 25) Fluorene 26) C1-Fluorenes	16.92	154	75291m	1196.79	
22) Acenaphthylene	18.35	152	88733m	1086.31	
23) Acenaphthene	18.94	154	50208m	1009.82	
24) Dibenzofuran	19.56	168	86010m	1221.40 ng/ml	
25) Fluorene	20.74	166	62621m	1138.35	
26) C1-Fluorenes	0.00	180	0	N.D. d	
27) C2-Fluorenes	0.00	194	0	N.D. a	
28) C3-Fluorenes	0.00		0	N.D. d	
31) Pentachlorophenol	23.30			1289.35 ng/ml	
32) Carbazole	24.75	167		1079.67 ng/ml	
33) Dibenzothiophene	23.57		95431m	921.87	
34) C1-Dibenzothiophene	0.00	198	0	N.D. d	which place while while comparing

^{(#) =} qualifier out of range (m) = manual integration MS30377H.D 041507.M Mon Apr 16 08:07:04 2007

Vial: 45 Data File: G:\1\DATA\MS30377\MS30377H.D Operator: TJM 4:37 pm

Acq On : 13 Apr 2007 Sample : Cal Level 5 Inst : GC/MS Ins

Multiplr: 1.00 Misc

MS Integration Params: rteint.p Quant Results File: 041507.RES Quant Time: Apr 14 20:03 19107

Quant Method : C:\MS30377\041507.M (RTE Integrator)

: PAH Calibration Table (2002) Title Last Update : Sat Apr 07 22:34:26 2007

Response via : Initial Calibration DataAcq Meth : PAH-2002

	Compound	R.T.	QIon	Response	Conc Unit	t Qvalue
35)	C2-Dibenzothiophene	0.00	212	0	N.D.	d
36)	C3-Dibenzothiophene	0.00	226	0	N.D.	
37)	C3-Dibenzothiophene Phenanthrene	24.01	178	94287m		
38)	Anthracene	24.18	178	107343m		
39)	1-Methylphenanthrene	26.13	192	82770m		
40)	Anthracene 1-Methylphenanthrene C1-Phenanthrene/Anthracene	0.00	192	107343m 82770m 0	N.D.	d
	C2-Phenanthrene/Anthracene	0.00	206	0	N.D.	
	C3-Phenanthrene/Anthracene		220	0	N.D.	d
	C4-Phenanthrene/Anthracene		234	0	N.D.	
	Naphthobenzothiophene	32.16	234	0 90540m 0	951.89	
	C1-Naphthobenzothiophene	0.00	248	0	N.D.	đ
	C2-Naphthobenzothiophene	0.00	262	0	N.D.	d
	C3-Naphthobenzothiophene	0.00	276	0	N.D.	
	Fluoranthene	28.09		114644m	1082.25	
	Pyrene	28.86		131442m	992.14	
	C1-Fluoranthenes/Pyrenes	0.00	216	0	N.D.	d
	C2-Fluoranthenes/Pyrenes	0.00	230	0	N.D.	d
52)	C3-Fluoranthenes/Pyrenes	0.00 0.00 0.00	244	0	N.D.	d
54)	Benz(a)anthracene	32.96	228	129040m	1192.98	
55)	Chrysene	33.07	228	97417m	1068.02	
56)	Chrysene C1-Chrysenes C2-Chrysenes C3-Chrysenes C4-Chrysenes C29-Hopane 18a-Oleanane C30-Hopane Benzo(b) fluoranthene	0.00	242	0	N.D.	d
57)	C2-Chrysenes	0.00	256	0	N.D. N.D. N.D.	d
58)	C3-Chrysenes	0.00	270	0	N.D.	d
59)	C4-Chrysenes	0.00	284	0	N.D.	d
61)	C29-Hopane	0.00	284 191	0		
62)	18a-Oleanane	0.00	191	0	N.D.	d
63)	C30-Hopane	41.74	191	40974m	968.34 n	g/ml
64)	Benzo(b) fluoranthene	36.43 36.50 37.31 37.49	252	119283m	1105.15	
65)	Benzo(k)fluoranthene	36.50	252	108634m	1195.70	
66)	Benzo(k) fluoranthene Benzo(e) pyrene Benzo(a) pyrene	37.31	252	112035m	1106.95	
67)	Benzo(a)pyrene	37.49	252	103137m	1164.80	
68)	Indeno(1,2,3-c,d)pyrene	41.84	276	82832m	1211.90	
69)	Dibenzo(a,h)anthracene	41.95	278	76502m	1193.42	
70)	C1-Dibenzo(a,h)anthracene	0.00	292	0	N.D.	d
	C2-Dibenzo(a,h)anthracene	0.00	306	0	N.D.	d
	C3-Dibenzo(a,h)anthracene	0.00	320	0	N.D.	d
	Benzo(g,h,i)perylene	43.06	276	91278m	1175.71	
	Perylene	37.77	252	0 91278m 109377m	1145.20	

08:07:06 2007

Mon Apr 16

041507.M

MS30377H.D

54.00

Quantitation Report

52.00 50.00 48.00 46.00 Quant Results File: 041507.RES 8.4 Benzo(g,h,i)perylene, T 42.00 GC/MS T ,ene-Js:181181(4),695,661;Clonebrit C30-Hobsue I TUM 40.00 Operator: Vial: 38.00 Multiplr Benzogang Tener, T. Benzogan T. Benzagan T venzo(a)bkrana q12,1 36.00 Denzo(k)fluoranthene, T Inst Benzo(b)fluoranthene, T TIC: MS30377H.D 34.00 (RTE Integrator) eunité)auquisceue 1 32.00 Naphthobenzothiophene, T 30.00 (2002)Pyrene, T 28,00 Fluoranthene, T G:\1\DATA\MS30377\MS30377H.D Sat Apr 14 20:04:44 2007 Initial Calibration 26.00 1-Methylphenanthrene, T G:\1\METHODS\041507.M PAH Calibration Table T ,elozedas T Calibration 24.00 Anthracene, T Dibenzothicphene, T Phenanthrene, T - Phenanthrene, Pentachlorophenol, T MS Integration Params: rteint.p 22.00 Quant Time: Apr 14 20:03 19107 20.00 Fluorene-d10, L 1, anelathylnaphthalene, T Dibenzofuran, T 13 Apr 2007 Cal Level 5 Acenaphthenewaldshaghthene, T 16.00 18.00 Acenaphthylene, T T ,lynently Biphenyl, T .3.6-Dimethyinsphthshene, T T ,enelsritriqsnivrite M_T ,enelsritriqsnivriteM-t14.00 Response via Last Update Tenzoul/sphinal-graphsalene, T 10.00 12.00 Data File Acq On Sample Method Title Misc Abundance 00006 85000 80000 75000 70000 -00009 5000 65000 55000 50000 45000 40000 35000 30000 25000 20000 10000 15000 Time--> 000101

Supporting Documents

Shipping, Sample Receiving, and Project Initiation Documents

B&B LABORATORIES RECEIVING/INTEGRITY REPORT

703318
Job: 300134 Date Received: 04 12 07 SDG#: 07041201
Job: 500134 Date Received: 04/12/07 SDG#: 07041201 Sender: GeoIns: pht - Keun Traner Buzzarde bay
1. Number of Shipping Containers:
Comments:
2. Airbill Present? (Yes)/No Shipping Company: Fedky Airbill Number: Comments:
3. Custody Seals on Container? No Yes Intact Not Intact Comments:
Chain of Custody Records? Comments No Yes
5. General Sample Conditions: Frozen Cool Unrefrigerated Dry Ice Blue Ice Ice Ice Ice
6. List of Broken Containers:
7. Number of Samples Expected: Number of Samples Received: Voc Voc Voc
8. Problems/Discrepancies:
9. Resolutions:
10. Checked in by: Jurel frank Date: 04/12/07
10. Checked in by: Wurld trank Date: 04/12/07

. 020	4a Express Package Service Packages up to 150 lbs. Packages up to 150 lbs. PedEx Priority Divernight Next business absence. PedEx First Opernight PedEx Fi	FedEx 2)3ay Second buseress day The training the training that the	Service Delivery	And in roof it regists to the control business day in the	Packaging 'Outranduraturaturaturaturaturaturaturaturaturat	Largo Prick, and Fred Ex. Startop Paik Suecial Handlinn	SATURDAY Delivery Available volve feeds. Provy Available volve feeds. Zevy Available v	Signer Statement of Stateme	Payment Bill to: Button Feeting Augs, New or Credit Card No. ballow. Sender Act No. Third Party Credit Card Credit Card Cash/Check Act No.		fotal Packages Total Weight Fotal Declared Value [†] Total Charges	1 Our habitity is limited to Shift unless you decline a higher value. See back for details.	Release Signature Sign to authorize delevery without obtaining signature.	By signing you authorize us to dehave this shipment without obtaining a signishing a signishing and ague to indemnify soft but shamless from any resulting claims. The signing you authorize us to dehave this signing the state of fedex.com or call tabol Go. Factic-8 8004863.7389 Res. Due 10011-Pint st Signit-8-2004-6-015-40041EB IN 12 A. WCS. 033
Express USA Airbill rects B4180527446 tom.		Name 166 HM HRS OV Phone 970 692 HM Soon	Company COSTUS GWT AV.	Address S A C (VE	State MA ZIP 0183(Your Internal Billing Reference 2011 CO CO 150	CONTRACTOR GAY LOLL LY LINE		FedEs location, print FodEs address. Was connect definer to P.D. boxes or P.D. Zip codes.	Days Floors's interfacent Contractions of the Contraction of the Contr	101 100 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	8418 0952 7446	8 8	

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By eighing you authorize us to defear this shipment without obtaining a signature and adgress to indemnity with those than these share in resulting cleans. Out and signed to be seen to the state of factor of the state o	or cell 1800.50 Foot F 800 A65.233. Rec. Date 1901 - Four 4 (Shift F \$1994-200 Foot F RRVIED IN U.S.A. VCSL U.S.	

Phone: 413-323-2332 FRX -419 000 6406

Emails info

CHAIN OF CUSTODY RECORD

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Page / of /

# of containers					æ	5 0				l iii	ii ii	[ii]	ا ا ا ا	ii) ii ii ii ii	ii] u b	id i		ii ii ii ii ii ii ii ii					
	~Cont.Code	Cont. Code:	G=glass P=plastic	ST=stentle V= viet	S=summa can	T=tedlar bag	0=Other		Comments:	Contack		The state of the s	978 672 4144		\$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$ \$	3			Codes:		X = Na hydroxide	X = Na hydroxide T = Na thiosulfate	X = Na hydroxid T = Na thiosulfat
		STED														-			**Preservation Codes:		l = lced	I = Iced H = HCL	I = Iced H = HCL M = Methanol
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<i>5</i> 111-7	200	(00)	check one): DWEBSITE CLIENT		CI PDF CI GIS KEY			Comp- osite Grab "Matrix Code	/ Cowore										Detection Limit Requirements	ordatione?	regulations	guidanoria:	Data Enhancement Project? TY
カルトでもり (死b) ::auduale1	Project # 3871-002	Client PO# 3371-001	DATA DELIVERY (check one): OFAX CEMAIL CIWEBSITI	Fax#;Email:	O EXCEL		Date Sampled	Start Stop Cc Date/Time Date/Time o	40/12/5										rnaround	☐ 24 Hour* Re		48 Hour*	48 Hour*
						J		Lab#											Date/Time:		1	1	1 .
	ddress: 5 (an Down S. L. 200)	Westford, MA 01886	Kevia Traine	2 Suy MA	- CHILLEN -	Billing purposes)	proposal date		WIF-02-32707								The second secon		Communication of the second				0
ompany Name:		Wester	ttention: <u> ke いい</u>	roject Location: Bo 22ads Say	ed by. The left	roposal Provided? (For Billing purposes)	U yes	Sample De	WIF-03										elinquished by: (signature)		****	(signature	experied by: (signature)

Con-Test Laboratory is the ONLY independent laboratory in all of New England with both prestigious AIHA and NELAC Certifications! Received by: (signature)

B = Sodium bisulfate S = Sulfuric Acid

> S = soil/solid SL = sludge 0 = other_

Special Requirements or DL's:

Date needed**

Date/Time:

*Require lab approval.

o = Other_

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7 7 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	Cilent Project #	3871-002
0000	500000	07041201
STIME	COMMENTS	Oiled rock
MATON	VEL SE	OTHER
SISVINA	20017000	PAH, TPH
UNUSA CANOSA	3	04/12/07
DATE:		03/27/07
S		WIF-02-32/07
FILENAME	***************************************	E17/20/
CLIENT NAME	Coolneight Dissessed Dog. Call	Georgiasgus-Duzzalus Day Spill

B&B LABORATORIES SAMPLE INITIATION FORM-ENV

Job#: JO33 (S Number of Samples:
SDG: 07041201 Matrix: 0,1 ed voct
Client. Go Ing. tht - Buzzards Bay Due Date:
Initiation Date: Ot 10 OF Soul Comments:
Initiation Date.
Analyses OCS/PCRS Aliphatics/TPH EOM
PARIS — OCSTEDS (C TAPATAN)
Dry Wt.
Short Columns Long Columns
Requested QA/QC (per batch of Client Samples) NO QA
□ Blank Spike □ Blank Spike Duplicate
Duplicate Matrix Spike
Matrix Spike Duplicate SRM
SEE BACK FOR SPECIFIC STANDARDS TO USE
Surrogate(s): High AL, AR Volume(s): 100ml
Spike Standard(s): Volume(s):
Internal Standard(s): High AL, AR Volume(s):
Final Extract Volume (ml): Final Solvent: K
Comments:
Sample Custodian Signature: Date Date Date Date Date Date Date Date
Project Administrator Signature: Date: 4//3/07 cc: COC Book
Spl Initiation.doc cc: COC Book Extraction Lab 000108

Extraction Standard Inventory

Organophosphates (OPs)			
□ OP-WKSU-0050-006 (Surrogate)			•
☐ OP-WKSK-TPT-006 (Surrogate)			
□ OP-WKIS-0100-006 (Int STD)			
Aliphatic Hydrocarbons (ALI/TPH)			
☐ AL-WKSU-20-009 (Surrogate)			
☐ AL-WKSK-100-010 (Spike)			
☐ AL-WKIS-200-007 (Int STD)			
AL-STSU-200-007 (High Surrogate)		4	
AL-STIS-2000-008 (High Int STD)			
Polycyclic Aromatic Hydrocarbons (PAHs)]	
☐ AR-WKSU-0500-016 (Surrogate)			
☐ AR-WKSK-1000-013 (Spike)			
☐ AR-WKIS-0500-011 (Int STD)			
AR-STSU-5000-009 (High Surrogate)		-	
AR-STIS-5000-008 (High Int STD)			·
Organochlorine Pesticides/PCBs (OCs/PCBs)			
OC-WKSU-1000-010 (Surrogate)			
] OC-WKSK-0400-008 (Spike)			
] PEST-WKSU-1000-003 (Surrogate)(Epsilon-HCH)			
OC-WKIS-1000-008 (Int STD)			
Polychlorinated Biphenyls (PCBs by GC/MS)			
] PCB-WKSU-008-008 (Surrogate)	\Box P	CB-WKIS-010-007 (Int STD)	
] PCB-WKSK-300-003 (Spike)			
olybrominated Diphenyl Ether (PBDEs) and Poly	bron	ninated Biphenyls (PBBs)	
PBDE-WKSU-1.0-005 (Surrogate)	\Box P	BB-WKSU-5.0-004 (Surrogate	e) .
PBDE-WKSK-1-006 (Spike)	□ P	BB-WKSK-2500-001 (Spike)	
PBDE-WKIS-0.5-005 (Int STD)	\Box P	BB-WKIS-5.0-004 (Int STD)	
Linear Alkylbenzenes (LABs)			

Laboratory Bench Sheet Logs

B&B LABORATORIES ENVIRONMENTAL EXTRACTION LOG

	MATRIX	10/10/101		Surrogate: / € μL	Spike: HL
<u>I</u> A.	区 OTHER	-t y	Lipids Y/K	PAH: A12-515V-520-621	PAH:
	□ WATER	الحا ا	Copper 2/M	Pest/PCB:	Pest/PCB:
<u></u>	SEDIMENT	***************************************	EOM (V)N	Aliphatic: AL-575 V-202 DA	Aliphatic:
	☐ TISSUE	Extraction Solvent: DCM Lot # 9222		Other	Other:
	(rock)	Final Solvent: のCA Lot # ソインコン Final Vol:	(m)	GC Int Std:	Turbo Vap II
8	General Comments:		IIM pe	58	, , , , , , , , , , , , , , , , , , ,
	N OA	A 10 ms & short colons summate	10/21/2		baun I (c):
	3		TA 12	Pest/PCB:	Pressure (>20psf):
		11.5h Suraste + latine Spike:			Check Weter Level:
		Internal:		Other:	Turbo Vap Date:
	Sample Name	Client ID Wet Wt. (g or L)	. Dry Wt. Dry Wt. % (9)	Hydromatrix (g) Extraction Comments	ments Internal Chain of Custody
_	toztk13	WIF - 02-32787			Extracti
2	ļ				(u/oz
က				1	Freight Tol que
4				1, ,	Extra
5					2/07
ဖ					Fight: To: 4M
7					Concentration
8					4-13-03 43-03
6					From Ell Er
<u></u>					
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<u> ₽</u> (rog /
001		Sample shaller with 300 ml of OCH	cont passed	that sodiu sulfat	کار 1635 Page 1 of 2
.11		ENVIROLOG3_D2_Working.xls Env. Extraction Log			

B&B LABORATORIES ENVIRONMENTAL EXTRACTION LOG

000112

B&B LABORATORIES EOM LOGBOOK

	Take I wil to short column. cone.	to 1 ml. DAG > 1ml ~ Which IS, SU	ar will lotigh IS, Su			Comments																	EOM 261 Page 1 of 2
ents:	at to 56	in the property	二年學			EOM (Dry Wt. Basis)		17.677															
General comm	Take In	to Iml	CDH: D	Date/Int:		EOM (Wet Wt. Basis)		7 AVE.	->									E					
			•		13-03	Wt. of 100 µl EOM Wt. (mg)	17.265	17.947	812-21														
		***		Date/Int:	11-13-03	Final Extract Vol (mL)	W	W	8														***************************************
201/206	Br. Dr	, R ()			ENVISE	Dry Wt. (%)			/														
SDG #: 0704/20/	Buzzede Bu Still			Transferred by Date/Int;	From ENV Pg: ピールレーとろ From DRY Pg:	Smpl Wt./Vol (g/L) Wet Wt. Dry Wt.			/									Verential de la companya de la compa					
Job#: 503318	Galasiah) ((a) (a) (a) (a) (a) (a) (a) (a) (a)		Lab Manager	1/3/06 gK	Client ID	40428-20-3/M	CU112-02-82707	WIF-02-5223														
MATRIX	OTHER	\ _		WAIEK		Sample Name	ETX 2207	6TX7207	ETX7207														FOM Logs
<u>L</u>	·····	********************************		***************************************				2	က	4	5	9	7	8	6	10	=	12	13	14	15	<u>ه</u>	00113

Smpl Wt./Vol (9/L) Dry Wt. (%) Extract Vol (mL) Lol (mL) Extract Dry wt.								
Smpl Wt./Vol (g/L) (g/L) Dry Wet Wt. Dry Dry Wt.								
Sample Name	17	18	19	20	21	22	23	24

The Relative Percent Difference (RPD) between duplicates must be ≤ 25%.

| Date/Int: RPD | RPD | Sample: | Duplicate: | Du

x 100%

 $(EOM_1 + EOM_2) \times 0.5$

 $(EOM_1 - EOM_2)$

%RPD=____

× 1000%

(EOM Wt. (mg)) (Final Extract Vol. (ml))
(Smpl Wt/Vol. (g/L)) (0.10 ml)

11

EOM

EOM 261Page 2 of 2

000114

Last Page



APPENDIX C

LEISURE SHORES TEST PIT LOCATIONS AND DESCRIPTIONS (SEPTEMBER 20, 2006)

Transect A

GPS Coordinates for Start Point (0,0): 41.62870° N 70.82366°

	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	
A-30	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	30
A-40	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	40
A-50	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	50
A-60	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 40% coverage		0	60
A-70	12	9	Approximately 80 particles of oil each less than 1 mm in diameter, 1 particle of oil 3 mm in diameter, and approximately 10% silver sheen coverage, fractured, breaks up, may be weathered	Approximately 6 particles of oil each less than 1 mm in diameter with approximately 15% coverage	0	70
A-80	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 40% coverage	Evidence of oil not observed in bucket	0	80
A-90	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	90
A-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	100
A-110	12	9	3 particles of oil each less than 1 mm in diameter and approximately 5% silver sheen coverage, breaks up when touched	2 particles of oil each less than 1 mm in diameter and approximately 10% silver sheen coverage, breaks up, may be from weathering	0	110
A-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	120
A-130	12	9	Evidence of oil not observed in test pit, orange/brown sediment observed on the bottom of the test pit	Evidence of oil not observed in bucket	0	130
A-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	140
A-150	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	150
A-160	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 80% coverage	Evidence of oil not observed in bucket	0	160
A-170	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 80% coverage	Evidence of oil not observed in bucket	0	170

Notes:

- 1. Test pits were excavated on September 20, 2006.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed
Sheen
1 to 5 particles of oil
6 to 10 particles of oil

11 to 20 particles of oil

21 to 100 particles of oil

Greater than 100 particles of oil

Transect B

Test Pit	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
B-30	12	9	Evidence of oil not observed in test pit	3 mm particle of oil observed on cobble	15	30
			•	sticking out above water line,		
				approximately 20 particles of oil each less		
				than 1 mm in diameter, approximately 5%		
				sheen coverage with the sheen radiating		
				from particles, largest sheen is		
				approximately 1 cm in diameter		
B-40	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	40
B-50	12	9	Evidence of oil not observed in test pit,	Evidence of oil not observed in bucket,	15	50
			organic "scum" observed on surface of	organic "scum" observed on surface of		
			test pit with approximately 5% coverage	water covering spoils with approximately		
D 60				20 % coverage		
B-60 B-70	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15 15	60 70
B-70	12	9	1 particle of oil less than 1 mm in	4 particles of oil each less than 1 mm in diameter, 1 particle of oil approximately 1	15	70
			long, less than 5% sheen and oil coverage,			
			organic "scum" observed on surface of	coverage, organic "scum" observed on		
			test pit with nearly 100% coverage	surface of water covering spoils with		
			less pit with hearly 100% coverage	approximately 75% coverage		
B-80	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	80
B-90	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	90
B-100	12	9	Evidence of oil not observed in test pit,	Evidence of oil not observed in bucket,	15	100
			organic "scum" observed on surface of	organic "scum" observed on surface of		
			test pit with approximately 70% coverage	water covering spoils with approximately		
				70% coverage		
B-110	12	9	Evidence of oil not observed in test pit,	Evidence of oil not observed in bucket	15	110
			organic "scum" observed on surface of			
			test pit with nearly 100% coverage			
B-120	12	9	Evidence of oil not observed in test pit,	Evidence of oil not observed in bucket	15	120
			organic "scum" observed on surface of			
			test pit with approximately 80% coverage			
D 420	12	9	B.:1 6.21 4.1 12.4 4.2		1.5	120
B-130	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of	Evidence of oil not observed in bucket, organic "scum" observed on surface of	15	130
			test pit with approximately 90% coverage	water covering spoils with approximately		
			test pit with approximatery 90% coverage	50% coverage		
B-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	140
B-150	12	9	Evidence of oil not observed in test pit.	Evidence of oil not observed in bucket.	15	150
2 100			organic "scum" observed on surface of	organic "scum" observed on surface of	15	150
			test pit with nearly 100% coverage	water covering spoils with nearly 100%		
				coverage		
B-160	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	160
B-170	12	9	Evidence of oil not observed in test pit,	Evidence of oil not observed in bucket	15	170
			organic "scum" observed on surface of			
			test pit with approximately 90% coverage			
D 100						
B-180	12	9	Evidence of oil not observed in test pit,	Evidence of oil not observed in bucket	15	180
			organic "scum" observed on surface of			
			test pit with approximately 20% coverage			
B-190	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	190
B-200	12	9	Evidence of oil not observed in test pit.	Evidence of oil not observed in bucket	15	200
2 200		,	organic "scum" observed in test pit,	or or not observed in bucket	1.5	230
			test pit with approximately 90% coverage			

Notes:

- 1. Test pits were excavated on September 20, 2006.
- Depths are approximate values.
 Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:
- Evidence of oil not observed Sheen
- 1 to 5 particles of oil 6 to 10 particles of oil
 - 11 to 20 particles of oil 21 to 100 particles of oil

 - Greater than 100 particles of oil

Transect C

Test Pit ID	Diameter (inches)	Depth (inches)	Field Team Observation - Test Pit	Field Team Observation - Spoils in Bucket	X Coordinate	Y Coordinate
C-20	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	10 particles of oil each 1 to 2 mm in diameter with sheen emanating from them, 1 particle of oil 3 mm long, approximately 25 % sheen coverage	30	20
C-30	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	30
C-40	12	9	Evidence of oil not observed in test pit	I particle of oil with sheen emanating from it, sheen approximately 1.5 cm in diameter.	30	40
C-50	12	9	Evidence of oil not observed in test pit	1 silver sheen approximately 1.5 inches long, no particles of oil observed	30	50
C-60	12	9	Evidence of oil not observed in test pit	1 particle of oil approximately 1 mm in diameter.	30	60
C-70	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	70
C-80	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	80
C-90	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with approximately 10 % coverage	30	90
C-100	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 90 % coverage	Evidence of oil not observed in bucket	30	100
C-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	110
C-120	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 5 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with approximately 5 % coverage	30	120
C-130	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 5 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with approximately 10 % coverage	30	130
C-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	140
C-150	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	150
C-160	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	160
C-170	12	9	Evidence of oil not observed in test pit, black sand particles observed, organic "scum" observed on surface of test pit with approximately 80% coverage	3 particles of oil each less than 1 mm in diameter, less than 5 % sheen coverage with sheen radiating from particles up to 1 cm	30	170
C-180	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	180
C-190	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 5 % coverage	water covering spoils with less than 5 % coverage	30	190
C-200	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 80 % coverage	Evidence of oil not observed in bucket	30	200

Notes:

- 1. Test pits were excavated on September 20, 2006.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed Sheen

1 to 5 particles of oil 6 to 10 particles of oil

11 to 20 particles of oil 21 to 100 particles of oil

Greater than 100 particles of oil

Transect D

Test Pit ID	(inches)	Depth (inches)	Field Team Observation - Test Pit	Field Team Observation - Spoils in Bucket	X Coordinate	Y Coordinate
D-30	12	9	Evidence of oil not observed in test pit, test pit was slightly underwater due to advancing tide	Evidence of oil not observed in bucket	45	30
D-40	12	9	Evidence of oil not observed in test pit, test pit was under less than 1 inch of water due to the advancing tide	Particles of oil not observed, less than 5 % silver sheen coverage	45	40
D-50	12	9	15 cm long streamer 1 to 2 cm wide	6 sheens 1.5 to 5 cm in diameter, 1 sheen 2 cm in diameter, 1 sheen 5 cm by 2.5 cm	45	50
D-60	12	9	1 particle of oil 2 mm in diameter, 2 sheens each 5 mm in diameter	25 particles of oil each less than 1 mm, 10 sheens some with slight rainbow each 5 mm in diameter, 5 silver streamers 3 to 5 cm long	45	60
D-70	12	9	Evidence of oil not observed in test pit	8 particles of oil each less than 1 mm with sheen radiating out 0.5 cm, 1 silver streamer 2.5 cm long and 3 mm wide	45	70
D-80	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage	45	80
D-90	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket,	45	90
D-100	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage.	45	100
D-110	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage.	45	110
D-120	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage.	45	120
D-130	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage.	45	130
D-140	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage.	45	140
D-150	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage.	45	150
D-160	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage.	45	160
D-170	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket,	45	170
D-180	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100% coverage	45	180
D-190	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100%	45	190
D-200	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with nearly 100%	45	200

Notes:

- Notes:

 1. Test pits were excavated on September 20, 2006.

 2. Depths are approximate values.

 3. Coordinates are reported in feet.

 4. Trenches were excavated at 10-foot intervals along the transect.

 5. mm = millimeters

 6. cm = centimeters

 7. Observations:

 Evidence of oil not observed

 Sheen

Evidence of oil not observed Sheen 1 to 5 particles of oil 6 to 10 particles of oil 11 to 20 particles of oil 21 to 100 particles of oil Greater than 100 particles of oil

Transect E

Test Pit	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
E-10	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	10
E-20	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	20
E-30	12	9	Approximately 100 particles of oil each	5 % silver to dull rainbow sheen	60	30
			less than 1 mm in diameter, 1 particle of	coverage, particles of oil not observed		
			oil approximately 3 mm in diameter, 1			
			stringer approximately 4 cm long, less			
			than 5 % silver to dull rainbow sheen			
E-40	12	9	1 particle of oil 3 mm in diameter, 7	7 particles of oil less than 1 mm in	60	40
			particles of oil each 2 mm in diameter,	diameter, approximately 25 % dull		
			approximately 200 particles of oil less	rainbow sheen coverage		
			than 1 mm in diameter, silver sheen			
			surrounding particles, approximately 5 %			
			total sheen coverage			
E-50	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	50
E-60	12	9	1 brown particle and sheen 1.5 cm	5% sheen coverage, 6 sheens 1 cm in	60	60
				diameter, 1 sheen 1.5 cm by 2 cm silver		
				sheen, 1 silver streamer 5 cm long		
				radiating from particles in center		
E-70	12	9	2 silver streamers of sheen 3 cm long and		60	70
			2 cm wide	organic "scum" observed on surface of		
				water covering spoils with nearly 100 %		
				coverage		
E-80	12	9	Evidence of oil not observed in test pit	1 particle of oil with sheen radiating	60	80
				around it 2 cm in diameter, 1 particle		
				with sheen around it approximately 5 cm		
				long and 2 cm wide		
E-90	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	90
E-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	100
E-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	110
E-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	120
E-130	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	130
E-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	140
E-150	12	9	Evidence of oil not observed in test pit	2 particles of oil surrounded by sheen	60	150
				radiating 5 mm, 6 silver streamers, 2 5		
				mm sheens with brown middle, 5% sheen		
F 160	12	0	Poids and of the state of the s	coverage	CO	1.00
E-160	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	160
E-170	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	170
E-190	12 12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	190
E-200	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	200

Notes:

- 1. Test pits were excavated on September 20, 2006.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed

Sheen

1 to 5 particles of oil

6 to 10 particles of oil

11 to 20 particles of oil

21 to 100 particles of oil

Greater than 100 particles of oil

Transect F

	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID To 10	(inches)	(inches)		Bucket	Coordinate	Coordinate
F-10	12	10.5	Evidence of oil not observed in test pit	1 sheen approximately 2 cm long surrounding <1 mm particle, 1 particle of oil with sheen 3 mm in diameter	75	10
F-20	12	10.5	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	20
F-30	12	10.5	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	30
F-40	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	40
F-50	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	50
F-60	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	60
F-70	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	70
F-80	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	80
F-90	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	90
F-100	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	100
F-110	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	110
F-120	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	75	120
F-130	12	10.5	1 long streamer of sheen approximately 1 foot long and 0.5 inches wide	1 particle of oil 1 mm in diameter, sheen 2 cm long radiating out from particle	75	130
F-140	12	10.5	Evidence of oil not observed in test pit, only approximately 1 inch of water in test pit	8 sheen clumps approximately 1 cm in diameter with brown centers composed of particles of oil each less than 1 mm in diameter, approximately 5 % silver sheen coverage, sheen stringers also	75	140
F-150	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 90 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with approximately 90 % coverage	75	150
F-160	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed	Evidence of oil not observed in bucket, organic "scum" observed	75	160
F-170	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with approximately 50 % coverage	Evidence of oil not observed in bucket	75	170
F-190	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with 90 % coverage	Evidence of oil not observed in bucket	75	190
F-200	12	10.5	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with 90 % coverage	Evidence of oil not observed in bucket	75	200

Notes:

- Notes:

 1. Test pits were excavated on September 20, 2006.

 2. Depths are approximate values.

 3. Coordinates are reported in feet.

 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters

7. Observations:
Evidence of oil not observed

Sheen

- 1 to 5 particles of oil 6 to 10 particles of oil

- 11 to 20 particles of oil 21 to 100 particles of oil
- Greater than 100 particles of oil

Transect G

Test Pit	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
G-70	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	90	70
G-80	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils with approximately 20 % coverage	90	80
G-90	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	90	90
G-100	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	1 particle of oil with sheen 1 mm in diameter	90	100
G-110	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit with nearly 100 % coverage	Evidence of oil not observed in bucket	90	110
G-120	12	9	Silver sheen 3 cm wide and 8 inches long around the edges of the test pit	Evidence of oil not observed in bucket	90	120
G-130	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	130
G-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	140
G-160	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	160
G-170	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	170
G-180	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils	90	180
G-190	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils	90	190
G-200	12	9	Evidence of oil not observed in test pit, organic "scum" observed on surface of test pit	Evidence of oil not observed in bucket, organic "scum" observed on surface of water covering spoils	90	200

Notes:

- 1. Test pits were excavated on September 20, 2006.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed

Sheen
1 to 5 particles of oil
6 to 10 particles of oil
11 to 20 particles of oil
21 to 100 particles of oil
Greater than 100 particles of oil

Vertical Distribution

Test Pit ID	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in Bucket	X	Y
	(inches)	(inches)			Coordinate	
A-70	12	2	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	70
A-70	12	4	1 dull rainbow sheen approximately 7 cm long and 5 cm wide seeping out of the sidewall and 1 sheen 2 cm long and 8 mm wide, approximately 15 % total sheen coverage	2 sheens, 1 sheen 1 cm diameter, 1 sheen 12 cm long and 1.5 cm wide, less than 5 % total sheen coverage	0	70
A-70	12	6	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	70
A-110	12	2	10 particles of oil 2 mm in diameter, 4 cm streamer, 1 sheen 2 cm in diameter, 1 sheen 4 cm in diameter, less than 5 % total sheen coverage	Evidence of oil not observed in bucket	0	110
A-110	12	4		4 cm rainbow streamer, 9 sheens 0.5 in diameter, 1 sheen 1 cm in diameter, less than 5 % total sheen coverage	0	110
A-110	12	6	7 sheens 0.5 to 1 cm in diameter, less than 5 % total sheen coverage	10 small sheens 3 mm to 1 cm in diameter	0	110
E-40	12	2	7 particles of oil each less than 1 mm, less than 5 % total sheen coverage	Evidence of oil not observed in bucket	60	40
E-40	12	4	20 particles of oil, less than 5 % total sheen coverage	Evidence of oil not observed in bucket	60	40
E-40	12	6	6 particles of oil, less than 5 % total sheen coverage.	Evidence of oil not observed in bucket	60	40

Notes:

- 1. Test pits were excavated on September 20, 2006.
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- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed

Sheen
1 to 5 particles of oil
6 to 10 particles of oil
11 to 20 particles of oil
21 to 100 particles of oil
Greater than 100 particles of oil



APPENDIX D

LEISURE SHORES TEST PIT LOCATIONS AND DESCRIPTIONS (JULY 10, 2007)

Transect A

GPS Coordinates for Start Point (0,0): 41.62870° N 70.82368°

	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
A-30	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 10% coverage	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with approximately 10% coverage	0	30
A-40	12	9	5 particles of oil up to 1 mm in diameter with silver sheen radiating out up to 1.5 cm, less than 5% sheen coverage	Evidence of oil not observed in bucket	0	40
A-50	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	50
A-60	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 60% coverage	Evidence of oil not observed in bucket	0	60
A-70	12	9	Approximately 40 particles of oil up to 2 mm in diameter, rainbow sheen up to 4 cm in diameter, approximately 5% rainbow sheen coverage	Approximately 40 particles of oil up to 5 mm in diameter, rainbow sheen up to 9 cm in diameter, approximately 20% sheen coverage	0	70
A-80	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	80
A-90	12	9	Evidence of oil not observed in test pit; red brown sediment settled to the bottom of the test pit	Evidence of oil not observed in bucket; red brown sediment settled to the top of the spoils in the bucket	0	90
A-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	100
A-110	12	9	1 sheen approximately 1.5 cm in diameter, less than 5% sheen coverage	Sheens up to approximately 2 cm in diameter, less than 5% sheen coverage	0	110
A-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	120
A-130	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	130
A-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	140
A-150	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	0	150
A-160	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 60% coverage		0	160
A-170	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 60% coverage	Evidence of oil not observed in bucket	0	170

Notes:

- 1. Test pits were excavated on July 10, 2007.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- $4.\ Trenches$ were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed
Sheen
1 to 5 particles of oil
6 to 10 particles of oil
11 to 20 particles of oil
21 to 100 particles of oil

Greater than 100 particles of oil

Transect B

Test Pit ID	Diameter (inches)	Depth (inches)	Field Team Observation - Test Pit	Field Team Observation - Spoils in Bucket	X Coordinate	Y Coordinate
B-30	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	30
B-40	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	40
B-50	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	50
B-60	12	9	1 sheen 1.5 cm in diameter, less than 5% sheen coverage	Evidence of oil not observed in bucket	15	60
B-70	12	9	Evidence of oil not observed in test pit	3 particles of oil up to 1 mm in diameter, silver sheen up to 6.5 cm in diameter, approximately 10% sheen coverage	15	70
B-80	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	80
B-90	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	90
B-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	100
B-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	110
B-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	120
B-130	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	130
B-140	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 10% coverage	Evidence of oil not observed in bucket	15	140
B-150	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	150
B-160	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	160
B-170	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	170
B-180	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	15	180
B-190	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 30% coverage	Evidence of oil not observed in bucket	15	190
B-200	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 40% coverage	Evidence of oil not observed in bucket	15	200

Notes:

- 1. Test pits were excavated on July 10, 2007.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed
Sheen

1 to 5 particles of oil

6 to 10 particles of oil

11 to 20 particles of oil

21 to 100 particles of oil

Greater than 100 particles of oil

Transect C

Test Pit	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
C-10	12	9	Evidence of oil not observed in test pi	Evidence of oil not observed in bucket	30	10
C-20	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	20
C-30	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	30
C-40	12	9	Evidence of oil not observed in test pit,	1 silver sheen approximately 7.5 cm in	30	40
			surface runoff flowing over test pit during evaluation	diameter, less than 5% sheen coverage		
C-50	12	9	Evidence of oil not observed in test pit	3 silver sheens up to 5 mm in diameter, less than 5% sheen coverage	30	50
C-60	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	60
C-70	12	9	Evidence of oil not observed in test pit,	1 particle of oil 1 mm in diameter with	30	70
			surface runoff flowing over test pit during evaluation	silver sheen radiating out up to 5 mm long and less than 2 mm wide, less than 5% sheen coverage		
C-80	12	9	Evidence of oil not observed in test pit	1 silver sheen approximately 7.5 cm wide and 10 cm long, less than 5% sheen coverage	30	80
C-90	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	90
C-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	100
C-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	110
C-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	120
C-130	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	130
C-140	12	9	Evidence of oil not observed in test pit;	Evidence of oil not observed in bucket;	30	140
			organic "scum" observed on surface of	organic "scum" observed on surface of		
			test pit with approximately 25% coverage			
				25% coverage		
C-150	12	9	Evidence of oil not observed in test pit;	Evidence of oil not observed in bucket;	30	150
			organic "scum" observed on surface of	organic "scum" observed on surface of		
			test pit with approximately 20% coverage	water covering spoils with approximately 10% coverage		
C-160	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	160
C-170	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	170
C-180	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	180
C-190	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	190
C-200	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	30	200

Notes:

- 1. Test pits were excavated on July 10, 2007.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed

Sheen
1 to 5 particles of oil
6 to 10 particles of oil
11 to 20 particles of oil
21 to 100 particles of oil
Greater than 100 particles of oil

Transect D

	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
D-20	12	9	Evidence of oil not observed in test pit	1 particle of oil 1 mm in diameter with sheen radiating out up to 2 mm in diameter, less than 5% sheen coverage	45	20
D-30	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 75% coverage	2 particles of oil less than 1 mm in diameter with sheen radiating out approximately 4 mm, less than 5% sheen coverage	45	30
D-40	12	9	Evidence of oil not observed in test pit	Approximately 25 particles of oil up to 2 mm in diameter with sheen radiating out up to 2.5 cm, approximately 5% sheen coverage	45	40
D-50	12	9	2 particles of oil 2 mm in diameter with sheen radiating out up to 2 cm in diameter	1 particle of oil 2 mm in diameter, silver sheen up to 5 cm in diameter, 10% sheen coverage	45	50
D-60	12	9	2 particles of oil less than 1 mm in diameter with silver sheen emanating off of them up to 5 mm in diameter	Approximately 40 particles of oil up to 3 mm in diameter with rainbow sheen radiating out up to 5 mm in diameter, approximately 20% sheen coverage	45	60
D-70	12	9	2 particles of oil less than 1 mm in diameter with silver sheen radiating out of them up to 3 mm in diameter	Evidence of oil not observed, possibly very weathered or inorganic (breaks up to the touch) sheen observed up to 1.5 cm in diameter, approximately 5% inorganic sheen coverage		70
D-80	12	9	2 particles of oil less than 2 mm in diameter with sheen radiating out approximately 12.5 cm long and 1.5 cm wide, less than 5% sheen coverage	Evidence of oil not observed in bucket	45	80
D-90	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	45	90
D-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with approximately 25% coverage	45	100
D-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	45	110
D-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	45	120
D-130	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	45	130
D-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	45	140
D-150 D-160	12	9	Evidence of oil not observed in test pit Evidence of oil not observed in test pit	Evidence of oil not observed in bucket Evidence of oil not observed in bucket	45 45	150 160
D-160 D-170	12	9	Evidence of oil not observed in test pit Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	45	170
D-170 D-180	12	9	Evidence of oil not observed in test pit Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	45	180
D-190	12	9	Evidence of oil not observed in test pit, surface runoff flowing over test pit during evaluation	Evidence of oil not observed in bucket	45	190

Notes:

- 1. Test pits were excavated on July 10, 2007.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters

7. Observations:
Evidence of oil not observed Sheen 1 to 5 particles of oil 6 to 10 particles of oil 11 to 20 particles of oil 21 to 100 particles of oil Greater than 100 particles of oil

Transect E

Test Pit ID	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
	(inches)	(inches)		Bucket	Coordinate	
E-10	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with approximately 5% coverage	60	10
E-20	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	20
E-30	12	9	Evidence of oil not observed in test pit, surface runoff flowing over test pit during evaluation	Sheen up to 2 mm in diameter, less than 5% sheen coverage	60	30
E-40	12	9	Evidence of oil not observed in test pit, surface runoff flowing over test pit during evaluation	1 particle of oil 2 mm in diameter, sheens up to 1.5 cm in diameter, less than 5% silver sheen coverage	60	40
E-50	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with approximately 10% coverage	60	50
E-60	12	9	2 particles of oil each 2 mm in diameter with rainbow sheen radiating out 4 cm, less than 5% sheen coverage	Evidence of oil not observed in bucket	60	60
E-70	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	70
E-80	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	80
E-90	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	90
E-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	100
E-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	110
E-120 E-130	12 12	9	Evidence of oil not observed in test pit Evidence of oil not observed in test pit	Evidence of oil not observed in bucket Evidence of oil not observed in bucket	60	120 130
E-140	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 10% coverage	Evidence of oil not observed in bucket	60	140
E-150	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with approximately 40% coverage	60	150
E-160	12	9	Test pit dry, evidence of oil not observed on sediment in test pit	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with approximately 40% coverage	60	160
E-170	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	170
E-200	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with approximately 50% coverage	60	200

Notes:

- $1. \ Test \ pits \ were \ excavated \ on \ July \ 10, \ 2007.$
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed Sheen

1 to 5 particles of oil 6 to 10 particles of oil

11 to 20 particles of oil

21 to 100 particles of oil

Greater than 100 particles of oil

Transect F

Test Pit ID	Diameter (inches)	Depth (inches)	Field Team Observation - Test Pit	Field Team Observation - Spoils in Bucket	X Coordinate	Y Coordinate
F-10	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 50% coverage	Evidence of oil not observed in bucket	75	10
F-20	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 25% coverage	Evidence of oil not observed in bucket	75	20
F-30	12	9	Evidence of oil not observed in test pit, surface runoff flowing over test pit during evaluation	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with less than 5% coverage	75	30
F-40	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with less than 5% coverage	75	40
F-50	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	50
F-60	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	60
F-70	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket, less than 5% inorganic (breaks up to the touch) sheen up to 5 cm in diameter	75	70
F-80	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	80
F-90	12	9	Approximately 5% silver sheen coverage, 1 stringer up to 2.5 cm wide and 10 cm long	Evidence of oil not observed in bucket	75	90
F-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	100
F-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	110
F-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	120
F-130	12	9	2 particles of oil up to 1 mm in diameter, silver sheen up to 2.5 cm in diameter, less than 5% sheen coverage	1 particle of oil less than 1 mm in diameter, silver sheens stringers up to 2.5 cm long and 2mm wide, less than 5% silver sheen coverage	75	130
F-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	140
F-150	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	150
F-160	12	9	Approximately 0.5" of water in test pit, evidence of oil not observed on surface of water or on the sediment in test pit	Evidence of oil not observed in bucket	75	160
F-170	12	9	Approximately 0.5" of water in test pit, evidence of oil not observed on surface of water or on the sediment in test pit	Evidence of oil not observed in bucket	75	170
F-190	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	75	190
F-200	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with less than 5% coverage	Evidence of oil not observed in bucket	75	200

Notes:

- 1. Test pits were excavated on July 10, 2007.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed
Sheen
1 to 5 particles of oil
6 to 10 particles of oil
11 to 20 particles of oil
21 to 100 particles of oil
Greater than 100 particles of oil

Transect G

Test Pit	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
G-10	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	10
G-20	12	9	Evidence of oil not observed in test pit, surface runoff flowing over test pit during evaluation		90	20
G-30	12	9	Evidence of oil not observed in test pit, surface runoff flowing over test pit during evaluation	Evidence of oil not observed in bucket	90	30
G-40	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	40
G-50	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	50
G-60	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	60
G-70	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	70
G-80	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	80
G-90	12	9	Evidence of oil not observed in test pit	1 particle of oil 1 mm in diameter with silver sheen stringer radiating out approximately 5 cm long, less than 5% sheen coverage	90	90
G-100	12	9	Evidence of oil not observed in test pit	1 silver sheen approximately 3 mm in diameter, less than 5% sheen coverage	90	100
G-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	110
G-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	120
G-130	12	9	Silver sheen stringers up to 2.5 cm long and up to 3 mm wide, less than 5% sheen coverage.	Evidence of oil not observed in bucket	90	130
G-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	140
G-150	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	150
G-160	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	160
G-170	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	90	170
G-190	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with less than 5% coverage	90	190
G-200	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 75% coverage	Evidence of oil not observed in bucket	90	200

Notes:

- 1. Test pits were excavated on July 10, 2007.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed

Sheen
1 to 5 particles of oil
6 to 10 particles of oil
11 to 20 particles of oil

21 to 100 particles of oil

Greater than 100 particles of oil

Transect H

Test Pit	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
H-50	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	50
H-60	12	9	Evidence of oil not observed in test pit;	Evidence of oil not observed in bucket;	105	60
			organic "scum" observed on surface of	organic "scum" observed on surface of		
			test pit with approximately 50% coverage	water covering spoils with approximately		
				50% coverage		
H-70	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	70
H-80	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	80
H-90	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	90
H-100	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	100
H-110	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	110
H-120	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	120
H-130	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	130
H-140	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	140
H-150	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	150
H-160	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 80% coverage	Evidence of oil not observe din bucket	105	160
H-170	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	170
H-190	12	9	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 80% coverage	Evidence of oil not observed in bucket; organic "scum" observed on surface of water covering spoils with approximately 70% coverage	105	190
H-200	12	9	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	105	200

Notes:

- 1. Test pits were excavated on July 10, 2007.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed
Sheen
1 to 5 particles of oil
6 to 10 particles of oil
11 to 20 particles of oil
21 to 100 particles of oil

Greater than 100 particles of oil

Vertical Distribution

	Diameter	Depth	Field Team Observation - Test Pit	Field Team Observation - Spoils in	X	Y
ID	(inches)	(inches)		Bucket	Coordinate	Coordinate
A-70	12	2	Large particles of oil interconnected in an area 7.5 cm in diameter and in an area 5 cm in diameter, an oiled rock was located adjacent to the test pit	approximately 10% sheen coverage	0	70
A-70	12	4		More than 200 particles of oil up to 5 mm in diameter with rainbow sheen radiating out	0	70
A-70	12	6	Approximately 35 particles of oil, 3 up to 2 cm in diameter, most of them approximately 1 mm in diameter with rainbow sheen radiating out	Approximately 10 particles of oil up to 1 mm in diameter, approximately 20% rainbow sheen coverage	0	70
D-60	12	2	Evidence of oil not observed in test pi	Evidence of oil not observed in bucket	45	60
D-60	12	4	Evidence of oil not observed in test pi	Evidence of oil not observed in bucket	45	60
D-60	12	6	2 particles of oil less than 1 mm in diameter with rainbow sheen stringer radiating out approximately 5 mm wide and 5 cm long	7 particles of oil up to 1 mm in diameter with rainbow sheen stringers radiating out up to 18 cm long and 5 mm wide	45	60
E-60	12	2	Evidence of oil not observed in test pit; organic "scum" observed on surface of test pit with approximately 50% coverage	Evidence of oil not observed in bucket	60	60
E-60	12	4	Evidence of oil not observed in test pit	Evidence of oil not observed in bucket	60	60
E-60	12	6	Evidence of oil not observed in test pit	3 sheens up to 3 mm in diameter, less than 5% sheen coverage	60	60

Notes:

- 1. Test pits were excavated on July 10, 2007.
- 2. Depths are approximate values.
- 3. Coordinates are reported in feet.
- 4. Trenches were excavated at 10-foot intervals along the transect.
- 5. mm = millimeters
- 6. cm = centimeters
- 7. Observations:

Evidence of oil not observed

Sheen
1 to 5 particles of oil
6 to 10 particles of oil
11 to 20 particles of oil
21 to 100 particles of oil
Greater than 100 particles of oil



APPENDIX E HEALTH AND SAFETY PLAN

SITE-SPECIFIC HEALTH & SAFETY PLAN

for

Project Name:

Bouchard B120 Oil Spill Buzzards Bay, Massachusetts

Project Location:

Brandt Island West (Leisure Shores and Howard's Beach) Mattapoisett, Massachusetts

Project Number: 3871-002

Prepared by:

GeoInsight, Inc. 5 Lan Drive, Suite 200 Westford, MA 01886

Reviewed by:		
Health & Safety Officer:		Date:
•	Christene A. Binger	
Project Manager:	Kevin D. Trainer	Date:
Field Team Leader/ On-Site Safety Officer:		Date:
	Kristin E. Zeman	
Field Personnel:	(1)	Date:
	(2)	Date:
	(3)	Date:
	(4)	Date:
	(5)	Date:

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FIGURES

FIGURE 1 Site Locus FIGURE 2 Site Map

APPENDICES

APPENDIX A Daily Safety Meeting and Employee Injury Report Forms

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SITE-SPECIFIC HEALTH & SAFETY PLAN

1.0 INTRODUCTION

The purpose of this Site Health & Safety Plan (HASP) is to define personal protection and monitoring protocols to be followed during remedial activities at the Leisure Shores and Howard's Beach portions of Brandt Island West in Mattapoisett, Massachusetts (the site). This HASP was prepared according to the requirements of 29 CFR 1910.120. These requirements and protocols are applicable to GeoInsight, Inc. (GeoInsight) employees and visitors. GeoInsight personnel, contractors, subcontractors, and visitors will be informed of the site emergency procedures and potential safety hazards involved with the anticipated activities. Subcontractors will be solely responsible for the health and safety of their personnel and will prepare and enforce their own HASP, which will be, at a minimum, consistent with the provisions of this HASP. This HASP summarizes those hazards and identifies personal protective measures required for this site. This plan must be reviewed by personnel prior to entering the site, and an agreement to comply with the requirements specified in the plan must be signed.

GeoInsight does not guarantee the health and safety of any person entering this site. Due to the hazardous nature of this site, and the activity occurring thereon, it is not possible to discover, evaluate, and provide protection for all possible hazards that may be encountered. Strict adherence to the health and safety guidelines set forth herein will reduce, but not entirely eliminate, the potential for injury at this site. The health and safety guidelines in this plan were prepared specifically for GeoInsight and this site, and should not be used on any other site without prior review and evaluation by trained health and safety personnel.

2.0 BACKGROUND INFORMATION

2.1 GENERAL SITE INFORMATION

The site is located at the Leisure Shores and Howard's Beach portions of Brandt Island West in Mattapoisett, Massachusetts. Refer to Figure 1 for the location of the site. Brandt Island West consists of Howard's Beach and Leisure Shores that are located on the western portion of the segment. An approximately 1,500 foot causeway connects to Brandt Island. This segment consists of the western portion of the causeway and western portion of Brandt Island. Nasketucket bay is located to the southwest, an unnamed pond is located to the north, and Buzzards Bay is located to the south and southeast of the segment. The primary shoreline consists of rip rap seawalls, bulkheads, piers, docks, and pilings. Residents and visitors use the beaches located at Leisure Shores which is primarily a mixed sand and gravel with cobble shoreline. In general, people use these shorelines primarily for seasonal recreational activities, including sunbathing, swimming, fishing, walking, and boating. Figure 2 shows the proposed work area, which is located in the intertidal zone in areas of the Leisure Shores and Howard's Beach portions of Brandt Island West.

2.2 HISTORICAL INFORMATION

Portions of the intertidal zone at Leisure Shores and Howard's Beach were impacted by Number 6 (No. 6) fuel oil released from Bouchard Barge B120 on April 23, 2003. Cleanup operations removed most of the No. 6 fuel oil stranded in the shoreline, but small amounts of No. 6 fuel oil remain present at the site. The residual oil consists primarily of "flecks" of oil in the sediment, sheen, isolated tarballs, and limited areas of buried oily sediment. In general, the oil is soft and tacky to the touch. Small areas of sheen are often produced when oiled sediment is disturbed (i.e., excavating test pits/trenches, removing oiled sediment/cobbles).

2.3 ANTICIPATED ACTIVITIES

The proposed field activities consist of digging test pits and trenches and conducting visual inspections to further characterize the extent or residual oil, and to identify whether additional cleanup activities are warranted. If warranted, cleanup activities will include using hand tools and/or light equipment to remove residual oil. Post-cleanup visual inspections will also be conducted to evaluate the cleanup effectiveness. During the cleanup activities, residual pavement will be removed from oil-impacted portions of the intertidal zone using hand tools (e.g., gardening trowels) and potentially light equipment (i.e., small excavator). Rocks with residual oil splatter will be either cleaned in-place using hand tools (e.g., wire brushes) or removed for disposal. It is anticipated that the characterization and cleanup activities will be initiated in September 2007.

A subcontractor may assist GeoInsight during cleanup activities. A representative from ENTRIX, Inc., the environmental consulting firm assisting with the Natural Resource Damage Assessment process, will be present during the cleanup and post-cleanup inspection activities.

Work will be conducted during the low tide "window" (i.e., approximately three hours before and after low tide).

3.0 POTENTIAL ON-SITE HAZARDS

3.1 0	VERALL HAZARD		
() Hi	igh () Moderate Low () Unknown		
3.2 H	AZARD TYPES		
X X X	sical Hazards (check all that apply): Slips, Trips, Falls Electrical Equipment Traffic Open Excavations Heavy Equipment Sharps Cold/Heat Stress Flammable Liquids Corrosive Liquids Manual Lifting	X X	Direct Sunlight Poisonous Plants Insects Poisonous Animals Noise Exposures Underground Utilities Overhead Electric Lines Other: Other:
Potent release aroma at this derma	emical Hazards ial chemical hazards identified based upon histori e of No. 6 fuel oil. Specific compounds of concer tic hydrocarbons (PAH) that were likely derived f site is weathered and not volatile, the expected ex l exposure.	n in media rom this s	a at the site are polynuclear ource. Because the residual oil
3) D 10	logical Hazards (check all that apply):		
	Raw Sewage/Septic Wastes		Viruses
	_ Medical Waste _ Bloodborne Pathogens	X	Bacteria Biting/Stinging Insects
3.3 W	ASTE TYPES		

Petroleum Hydrocarbons

<u>Health Hazard:</u> Causes eye irritation, skin irritation including redness, and a burning sensation. Prolonged or repeated contact can cause drying and cracking of the skin leading to dermatitis (inflammation). Harmful effects from skin adsorption are not expected.

First Aid:

Eye – if irritation or redness develops, immediately flush eyes with clean water for 5 minutes. Skin – wipe material from skin, remove contaminated clothing, wash affected area with mild soap and water.

For all other issues or if symptoms persist, seek medical attention.

Polynuclear Aromatic Hydrocarbons (PAHs)

<u>Health Hazard:</u> Causes eye, skin, and respiratory tract irritation, other symptoms include headache, nausea, confusion, and excitement. Chronic ingestion may result in gastrointestinal and kidney pain. Some PAHs are carcinogenic.

First Aid:

Eye – if irritation or redness develops, immediately flush eyes with clean water for 5 minutes. Skin – wipe material from skin, remove contaminated clothing, wash affected area with mild soap and water.

For all other issues or if symptoms persist, seek medical attention.

4.0 MONITORING PROTOCOL/INSTRUMENTATION

Because residual oil is present in small isolated areas and is typically weathered and often hardened on the surface, this oil does not contain sufficient volatiles to adversely affect ambient air, field monitoring of VOCs in ambient air will not be conducted.

Ambient air temperature will be monitored periodically and work periods will be adjusted to provide adequate rest and cool down periods for personnel. Personnel will be checked periodically for symptoms of heat-related problems. If high temperatures cause personnel to exhibit signs of heat stroke, heat exhaustion or dehydration, then field activities will be immediately halted and personnel will seek shelter and consumable liquids.

5.0 PERSONAL PROTECTIVE EQUIPMENT (PPE)

5.1 LEVELS OF PROTECTION

Level D

Tasks: Cleanup oversight and visual inspections

Modified Level D (based upon potential for dermal contact)

Tasks: Cleanup activities.

Level C

Tasks: None

5.2 PPE FOR SPECIFIED PROTECTION LEVELS

LEVEL D:

Work clothes
Steel toe workboots
Safety glasses with side shields
:

MODIFIED LEVEL D:

Level D PPE and; Tyvek outer suit (spun olefin) Chemical resistant boots or disposable, protective boot covers Nitrile gloves

6.0 PERSONAL DECONTAMINATION

On-site personnel will employ some measure of decontamination when leaving the work zone. PPE to be decontaminated may include boots, chemical resistant gloves, and other reusable equipment or materials that may have contacted affected media. Actual decontamination procedures will ultimately depend upon the level of protection, field screening results, and the results of ambient air monitoring. Typical decontamination procedures for the anticipated work levels are summarized below.

Level D:

Wash hands prior to eating and remove or clean work boots before returning to office.

Modified Level D:

Segregated equipment drop, boot and glove wash/rinse, boot removal, Tyvek suit removal, outer glove removal, field wash. Disposable items that become contaminated will be segregated and disposed.

Cleaning Solution:

Mixture of "Alconox" and potable water.

Rinse: potable water.

Field personnel can use alcohol-based or detergent "hand wipes" to remove oil adhered to fingers.

7.0 WORK ZONE DELINEATION

7.1 GENERAL

Work areas at the facility will vary with the type of work or task being conducted. The Site Health and Safety Officer (SHSO) will delineate the work zones prior to the start of work and document them in the project field logbook. The SHSO will use the following descriptions to determine the work zones.

7.2 DESCRIPTION OF WORK ZONES

Support Zone

Support activities, such as equipment deliveries, preparation for site activities, and meetings with personnel unauthorized to enter work zones will be performed within the designated Support Zone. The Support Zone is located a safe distance away from work activities at the site. The exact location of this zone will vary depending on where the work is being performed at the site, but should remain at least 20 feet away from the Work Zone.

Decontamination Zone

The Decontamination Zone will be adjacent to, but separate from, the Support Zone. Final decontamination of personnel and equipment will take place within the Decontamination Zone, as necessary, to minimize relocation or spreading of impacted materials. Certain initial decontamination activities, such as removal of large quantities of debris from equipment, will be performed within exclusion zones to the extent practicable to minimize the amount of material brought into the Decontamination Zone.

Work Zone

The Work Zone (exclusion area) will depend on the type of work being performed. Localized Work Zones will include the immediate vicinity of drilling and excavation activities, remote monitoring wells, and surface water sampling locations. Persons not directly involved in the site activities will be required to remain at least 20 feet away from the perimeter of an established Work Zone.

8.0 ON-SITE ORGANIZATION AND CONTROL

8.1 ORGANIZATION

Project Manager (Kevin Trainer):

Responsible for allocation of resources for the implementation of the HASP; assignment of personnel who meet the medical and training requirements of the HASP; and allocation of resources to resolve health and safety issues identified during the performance of project tasks.

GeoInsight Corporate Health and Safety Officer (CHSO: Michael Redding):

The CHSO is responsible for the overall coordination of the GeoInsight Corporate Health and Safety Program. The CHSO should be informed of any exceedence of a PEL, injuries, near misses, and general health and safety concerns.

GeoInsight Office Health and Safety Officer (OHSO: Christene Binger):

The OHSO is responsible to the CHSO in matters related to health and safety, including investigation of health and safety related incidents at a site.

Site Health and Safety Officer (Kristin Zeman):

Responsible to the GeoInsight CHSO in matters related to health and safety the on site, including development and implementation of the site-specific HASP; conducting site safety meetings and site-specific training of site personnel; investigation of health and safety related incidents at the site; accompanying authorized visitors on site tours; and updating and modifying this HASP, as necessary, if site or environmental conditions change. GeoInsight is not responsible for public present at the site, municipal employees, or contractors.

8.2 SITE SECURITY

The field team leader will control access to the site during GeoInsight site activities. GeoInsight personnel present at the site will be recorded in the field log of daily activities. The site HSO or field team leader will be responsible for enforcing adherence to work zone delineations described above.

8.3 COMMUNICATION

Because of the relatively small size of the site and the level of protection typically used, voice and hand signals will likely be sufficient for the anticipated work activities.

8.4 MEDICAL MONITORING REQUIREMENTS

GeoInsight personnel who perform on-site activities where there is potential for exposure to hazardous substances must have completed a medical monitoring examination no earlier than 12 months prior to commencing these site activities. The examination must comply with requirements specified by 29 CFR 1910.120 (f). A certification, signed by a medical doctor, must indicate work limitations, if any, placed on the individual. The certification must also

specify that the individual is capable of working while wearing respiratory protection equipment. The certification must be in the corporate health and safety file before the individual may begin on-site activities.

8.5 TRAINING REQUIREMENTS

Personnel who perform activities where there is potential for exposure to hazardous substances must have completed an initial Hazardous Waste Operations and Emergency Response (HAZWOPER) course or an annual refresher course of the initial training, within 12 months prior to the beginning of site activities. The HAZWOPER training must comply with requirements outlined in 29 CFR 1910.120 (e). A certificate indicating successful completion of this training must be in GeoInsight's project file for GeoInsight personnel.

Subcontractors performing excavation and trenching work are required to have a "competent person," as defined by OSHA and requiring OSHA-specified training, on-site while this type of work is in progress.

The site HSO will hold daily meetings with field personnel before work commences to discuss safety issues. During the meeting, personnel working on-site will be provided access to this HASP. The HASP will be reviewed and discussed and questions answered. Personnel who will work on-site will sign this HASP (or, in the case of subcontractor personnel, the subcontractor's HASP) to indicate that they have reviewed and understand site conditions and agree to comply with HASP requirements. The site HSO will record the daily meetings on the form titled "Daily Safety Meeting Form," a copy of which is attached in Appendix A.

9.0 GENERAL HEALTH AND SAFETY REQUIREMENTS

9.1 STANDING ORDERS

The following standing orders apply to the activities anticipated to be performed at the site during the project:

- 1. No eating, drinking, chewing tobacco or toothpicks, application of cosmetics, storing food or food containers or open flames permitted within the Work Zone.
- 2. No smoking within the perimeter of the site.
- 3. Wear the appropriate level of protection as defined in this HASP.
- 4. Wear latex or nitrile surgical gloves and use a physical barrier when providing emergency first aid or CPR.
- 5. Work must be restricted to daylight hours only.
- 6. Maintain close contact with your work partner while in the Exclusion Zone.
- 7. Persons with beards and mustaches that interfere with respirator fit and seal will not be allowed to work at activities requiring Level C or Level B protection.
- 8. Report any unusual conditions to the field team leader immediately.

9.2 INCIDENT REPORTING

Any incident or accident involving field personnel must be documented. Situations covered by this requirement include, but are not limited to, fires, explosions, exposures above the Permissible Exposure Limit (PEL) or Short Term Exposure Limit (STEL), illness and injuries, however minor. The site HSO or field team leader must be notified immediately so that first aid requirements can be assessed and transportation to the nearest medical treatment facility provided, if required. Reports of the incident must be provided to the GeoInsight Corporate HSO within 24 hours of the incident and includes completion of the accident reporting form in Appendix B.

10.0 APPENDED INFORMATION

The following documents (attached in Appendix B) provide additional information regarding issues that are considered applicable to anticipated site activities:

- 1. PPE Checklist
- 2. Heat Stress and Heat Stroke Prevention Guidelines
- 3. Emergency First Aid





SOURCE:

MASSACHUSETTS GEOGRAPHIC INFORMATION SYSTEMS (MASSGIS)

2001 AERIAL PHOTOGRAPH



ROJECT:				
	B120	OIL	SPII	I

LOCATION:

WIF-02 MATTAPOISETT, MASSACHUSETTS

TITLE:

APPROXIMATE LOCATIONS OF PROPOSE PHASE IV FIELD ACTIVITIES

DESIGNED:	DRAWN:	CHECKED:	APPROVED:
KEZ	KEZ	KDT	MJW
SCALE:	DATE:	FILE NO.:	PROJECT NO.:
AS SHOWN	08/03/07	387I-OILING	3871-002



Environmental Strategy & Engineering <u>Practical in Nature</u>

FIGURE NO.:

4

11.0 EMERGENCY INFORMATION/REFERENCES

EMERGENCY PHONE NUMBERS (for the Town of Fairhaven)

AMBULANCE: 911 POLICE: 911 FIRE: 911

POISON CONTROL CENTER: 1-800-562-8236

HOSPITAL St Luke's Hospital

101 Page Street New Bedford, MA 508-997-1515

HOSPITAL ROUTE: (Refer to attached map and directions)

OTHER CONTACTS PHONE NUMBERS

GeoInsight Inc. 978-692-1114 Christene Binger – OHSO (cell) 617-803-8108 Kevin Trainer – Project Manager (cell) 978-790-5294



CLICK HERE or go to lasikplus.com

Start: 190 Brandt Island Rd

Mattapoisett, MA 02739-1789,

US

End:

101 Page St

New Bedford, MA 02740-3464, US

Notes:

Only text visible within note field will print.





Direction	ons	Distance
Total E	Est. Time: 21 minutes Total Est. Distance: 8.62 miles	
START	1: Start out going NORTHWEST on BRANDT ISLAND RD toward DAVID	O ST. 2.2 miles
WEST 6	2: Turn LEFT onto FAIRHAVEN RD / US-6. Continue to follow US-6 W.	4.9 miles
\Leftrightarrow	3: Turn LEFT onto COUNTY ST.	0.7 miles
\Leftrightarrow	4: Turn RIGHT onto ALLEN ST.	0.4 miles
\Leftrightarrow	5: Turn RIGHT onto PAGE ST.	<0.1 miles
END	6: End at 101 Page St New Bedford, MA 02740-3464, US	
Total E	st. Time: 21 minutes Total Est. Distance: 8.62 miles	



Start: 190 Brandt Island Rd Mattapoisett, MA 02739-1789, US

MAPQUEST B

O

ISSUE | 1600 m

4800 ft

Mattapoisett
Harbor

Nasketucket
Bay

© 2007 MapQuest Inc. West

© 2007 NAVTEQ

End: 101 Page St New Bedford, MA 02740-3464, US



All rights reserved. Use Subject to License/Copyright

These directions are informational only. No representation is made or warranty given as to their content, road conditions or route usability or expeditiousness. User assumes all risk of use. MapQuest and its suppliers assume no responsibility for any loss or delay resulting from such use.

APPENDIX A DAILY SAFETY MEETING AND EMPLOYEE INJURY REPORT FORMS



Daily Safety Meeting Form

Weather Conditions:	Site Location:		
Site Conditions:	GeoInsight Proj. #:		
	Date:		
TOPICS DISCUSSED (Please Check Ones That App	oly)		
 Health & Safety Emergency Numbers Hospital Locations Work Areas Posted Designated Smoking Areas Confined Space Entry Slip, Trip, & Fall Manual Lifting Utility Locations Mechanical Hazards PERSONAL PROTECTIVE EQUIPMENT	Bonding & Grounding Heavy Equipment Traffic Hazards Heat or Cold Stress Air Compressor Lock Out/Tag Out Excavation Hazards Venting/Inerting		
 Eye Protection Hearing Protection Gloves Respiratory Protection Engineering Controls 	Hard HatProtective ClothingRetrieval SystemBackup System		
ADDITIONAL COMMENTS:			
MEETING ATTENDED BY THE FOLLOWING:			
This meeting was conducted by:	on .		



EMPLOYEE INJURY REPORT

Please answer all questions completely. This report must be forwarded to the Corporate Heath and Safety Officer within 24 hours of the injury.

Injured's Name			Sex	_ S.S.#	Birth date		
Home Address							
City	State	Zip		Phone			
Job Title			Hire Date		Hourly wage		
Number of days wo	orked per week:						
Date of Incident		Time_		Time reported	To whom?_		
Project Name			Project #_		Time work be	egan	
Has employee retur	ned to work? Yes _	No W	hen?	Did the e	mployee miss scheduled	l work?	
At what position?_							
Medical attention:	None First a	aid on-site	Docto	or's office	_Hospital ERHosp	oitalized	
Doctor/Hospital na	me		Ao	ddress			
Witness name(s)					Statement attached? _	Yes No	
Nature of injury or	Illness		Exa	ct body part			
Job assignment at ti	me of incident						
Was this his/her reg	gular occupation? If not,	state regula	ar occupatio	n			
Describe incident_							
	ral condition or unsafe action has been taken to pre	vent recurr	rence?		Date		
Site Safety Officer	Print	Signature		Date	Date		
	Print		Signature		Date	Date	
Comments on incid Manager's name	ent and corrective action						
Print Sig			nature		Date		
Concur with action	taken? Yes I	No Rema	arks				
Reviewed by Corpo	orate H&S Officer						
Print	Signature				Date		
Completed by	Title		Phone_		_ Date		
J:\Health and Safety\HASP\HAS	SPappendix\INJURY REPORT.doc						

APPENDIX B

GUIDELINES AND ADDITIONAL INFORMATION FOR SITE-SPECIFIC ACTIVITIES

PPE INSPECTION CHECKLIST

Before using personal protective clothing (PPE), inspect each article for defects, according to the checklist below.

Determine that appropriate clothing material is compatible with the anticipated chemical and exposure conditions of the project and its required tasks.

Visually inspect for:

- X imperfect seams;
- X non-uniform coatings;
- X tears; and
- X defective zippers and other closures.

Hold up to the light and check for pinholes.

- X Flex the product and inspect it for:
- X cracks; and
- X any defects which indicate that the product=s shelf life has been exceeded.
- X If the product has been used previously, inspect it inside and out for signs of chemical attack.

For example:

- X discoloration;
- X swelling; or
- X stiffness.

Before using gloves, pressurize each one with air to make sure that it has no holes. Also, visually inspect the gloves for the following defects:

- X imperfect seams;
- X tears; and
- X non-uniform coating.

While performing work, be alert for evidence of PPE failure, breakthrough, or excessive wearand-tear, as described below.

- X Evidence of chemical attack such as discoloration, swelling, stiffening and softening. Keep in mind, however, that chemical permeation can occur without any visible effects.
- X Closure failure.
- X Tears.
- X Punctures.
- X Seam discontinuities.

HEAT STRESS AND HEAT STROKE PREVENTION GUIDELINES

Because protective outerwear decreases ventilation, some fieldwork in hot weather increases the potential for heat-related casualties. Thus, field personnel must be alert to the symptoms of heat stress and must respond to them promptly and effectively.

1.0 PREVENTION OF HEAT STRESS AND HEAT STROKE

One of the major causes of heat stress and stroke is the depletion of body fluids. On-site there should be enough potable fluids available to prevent this. Personnel should replace water and salts lost in perspiration. Salts can be replaced either by consuming a 0.1% salt solution, more heavily salted foods, or commercial products such as Gatorade. Commercial products may be the preferred choice for personnel on low sodium diets.

A work schedule should be established so that the majority of work will be performed early in the day, before ambient air temperatures peak.

No work should be permitted at any site under Level B protection without prior approval of the Corporate HSO. A work/rest guideline will be implemented for personnel who wear Level B protection, as follows:

	<u> Maximum Work Time</u>		
Ambient Temperatures	Between 15 min. Rest Periods		
Above 90°F	1/2 hour		
80°F to 90°F	1 hour		
70°F to 80°F	2 hours		
60°F to 70°F	3 hours		
<60°F	4 hours		

Adequate time will be allowed for personnel to cool down. This may require shifts of workers during field activities.

2.0 HEAT STRESS MONITORING

One of both of the following techniques should be used to determine each site worker's ability to recuperate from working under hot conditions. Personnel wearing protective clothing should be monitored first when the ambient temperature is 70°F or above. They should be monitored with increasing frequency as the ambient temperature increases or if they are slow to recover after a rest work period. If temperatures exceed 80°F, workers must be monitored for heat stress after each work period.

- X Heart Rate Heart rate (HR) should be measured by the radial pulse for 30 seconds, as early as possible during the resting period. The HR at the beginning of the rest period should not exceed 110 beats per minute. If the HR is higher, the next work period should be shortened by 10 minutes (or 33%), while the length of the rest period stays the same. If the pulse rate is 110 beats per minute at the beginning of the next rest period, the following work period should be shortened by 33%.
- Body Temperature Body temperature should be measured orally with a clinical thermometer as early as possible during the resting period. Oral Temperature (OT) at the beginning of the rest period should not exceed 99°F. If it does, the next work period should be shortened by 10 minutes (or 33%), while the length of the rest period stays the same. If, however, the OT exceeds 99°F at the beginning of the next work period, the following work period should be further shortened by 33%. OT should be measured again at the end of the rest period to make sure that it has dropped below 99°F. It the OT exceeds 100.6°F, the individual should be removed from their chemical protective clothing.

Good Hygienic standards must be maintained by frequent change of clothing and daily showering. Clothing should be permitted to dry during rest periods. Personnel who notice skin problems should immediately consult medical processionals.

3.0 HEAT EXHAUSTION - IDENTIFICATION AND TREATMENT

3.1 Symptoms

Heat Exhaustion usually begins with muscular weakness, dizziness, nausea, and a staggering gait. Vomiting is frequent. The bowels may move involuntarily. The victim may be very pale, his/her skin may be clammy, may perspire profusely and may faint unless he/she lies down. This may pass, but sometimes it remains as heat stress progresses to heat stroke, a life-threatening condition.

3.2 First Aid

Immediately remove the victim to the Contamination Reduction Zone, preferably in a shady or cool area with good air circulation. Remove all protective outer wear. Call a physician. Treat the victim for shock:

- X Make the victim lie down;
- X Raise the victim's feet 6 to 12 inches above his/her head; and
- X Keep the victim warm but loosen all clothing.

If the victim is conscious, it may be helpful to administer sips of a salt water solution: one teaspoon of salt to ton 8-ounce glass of water. Transport the victim to a medical facility as soon as possible.

4.0 HEAT STROKE - IDENTIFICATION AND TREATMENT

4.1 Symptoms

This is the most serious of heat casualties because the body overheats to potentially life-threatening levels. Body temperatures of heat stroke victims may rise rapidly to between 107 to 110°F. First, the victim may experience headache, dizziness, and nausea. Almost always the victim's skin will be dry, red, and hot. Unconsciousness follows quickly and death is imminent if exposure continues. Onset of heat stroke usually is sudden.

4.2 First Aid

Immediately evacuate the victim to a cool, shady area in the Contamination Reduction Zone. Remove all protective outerwear and all personal clothing. Lay the victim on the back with the head and shoulder slightly elevated. It is imperative that the body temperature be lowered immediately. Apply cold, wet towels, ice bags, etc. to the head, sponge off the bare skin with cool water or rubbing alcohol, if available, or even place the victim in a tub of cool water. The main objective is to cool the victim without chilling him/her. Give no stimulants. Transport the victim to a medical facility as soon as possible.

EMERGENCY FIRST AID PROCEDURES

Red Cross first aid procedures should be used to treat personnel who are injured while working in both contaminated and uncontaminated areas. If the injured person can be moved, he/she should be taken outside of the work area or the Contamination Reduction Zone, as appropriate. Any contaminated clothing should be removed, if possible, and fist aid should be administered. Depending on the nature and severity of the injury, fist aid should continue during transport to a medical facility and until treatment is obtained.

For most injuries, the earliest measures, described below, will be among the most important to effectively administer first aid.

First, survey the scene. Is the injured person in imminent danger or further injury if left in place until his/her injuries can be assessed? Are others in the area in danger of injury?

Perform a primary survey of injuries to determine if emergency first aid measures are needed (e.g., as in the case of severe bleeding, etc.) before moving the victim outside the work area or to the Contamination Reduction Zone. Determine if it is safe for those other than emergency medical personnel to move the victim and, if so, which methods are appropriate to avoid compounding his/her injuries. Request emergency medical services (EMS) if the primary survey indicates that this is necessary because of the nature or severity of the injuries.

Perform a secondary survey of the victim's injuries. Determine if there are signs and symptoms of internal bleeding, imminent shock, or other potentially life-threatening conditions. If such symptoms are suspected, request EMS if this has not already been done, then immediately administer appropriate first aid until medical treatment is obtained.

Control any external bleeding using direct pressure. Elevate injured or bleeding areas, unless a fracture is suspected. Monitor the victim's airway, breathing, and level of consciousness frequently. Reassure the victim and keep him/her warm and as comfortable as possible. Even if symptoms of shock are not apparent, keep in mind that onset of shock may be sudden. Take measures to avoid shock immediately should it occur. If the victim vomits, place him/her on his/her side, and clear the airway if it becomes obstructed.

To reduce the risk of being infected by the victim when attempting to control bleeding, the caregiver should use some sort of barrier (e.g., several dressings, latex gloves, or a piece of plastic wrap). Hands should always be washed thoroughly after first aid is given. If there is more than one victim, a caregiver should always change gloves or wash his/her hands after touching one victim and before touching another (or any other individual) to prevent crossing-contamination.

GeoInsight, Inc.



APPENDIX F NOTICE OF DOCUMENT AVAILABILITY



August 2, 2007

GeoInsight Project 3871-002 delivered by Certified Mail

Chief Municipal Officer Town Hall 16 Main Street Mattapoisett, Massachusetts 02739

RE: Notice of Document Availability

Phase IV Remedy Implementation Plan

Barge B120 Spill

Buzzards Bay, Massachusetts

Release Tracking Number (RTN) 4-17786

Dear Chief Municipal Officer:

In accordance with Public Notification requirements of the Massachusetts Contingency Plan (MCP; 310 CMR 40.1403), please accept this letter as notification that a Phase IV Remedy Implementation Plan to remediate a small amount of residual oil located at portions of Leisure Shores and Howard's Beach will be filed with and available for your review after August 3, 2007 at:

> Massachusetts Department of Environmental Protection Southeast Regional Office, Bureau of Waste Site Cleanup 20 Riverside Drive Lakeville, Massachusetts 02347 Service Center: 508-946-2718 Fax: 508-946-2865

http://www.mass.gov/dep/about/region/serofile.htm

File reviews are conducted Tuesdays and Wednesdays from 9:00 a.m. to 11:30 a.m. and 2:00 p.m. to 4:30 p.m. (except state holidays). An electronic copy of this report will also be posted at www.buzzardsbay.org. Field activities are scheduled to be conducted in September 2007.

Fax (203) 271-8038



We trust this information is sufficient for your files. Please contact me at (978) 692-1114 if you have questions regarding the Phase IV Remedy Implementation Plan.

Sincerely,

GEOINSIGHT, INC.

Kevin D. Trainer, C.P.G., P.G., LS.P.

Senior Project Geologist

cc: MADEP, SERO - Lakeville, Massachusetts

Mattapoisett Board of Health - Daniel C. Lee, Jr.

Richard J. Wozmak, P.E., P.H., L.S.P. – EnviroLogic LLC.