

# Paragon Analytics

## Total Extractable Hydrocarbons (Diesel) Case Narrative

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### **Cordilleran Compliance Services, Inc.**

Rulison Area Well monitoring

**Order Number - 0811110**

1. This report consists of 2 water samples. The samples were received cool and intact by Paragon on 11/14/2008.
2. The water samples were extracted by adding hexane to the water sample and shaking the resulting two phase solution according to Paragon Analytics Standard Operating Procedure 603 Revision 10, which was developed at Paragon Analytics. The hydrocarbons partition into the hexane layer, which is then removed for analysis.
3. The extracts were then analyzed using GC with a DB-5.625 capillary column and a flame ionization detector (FID) according to Paragon Analytics Standard Operating Procedure 406 Revision 13 generally based on SW-846 Method 8000B and Method 8015B. The procedures are based on this general method because SW-846 does not have a specific method for total extractable petroleum hydrocarbons (TEPH) or diesel range organics. The only true modification from this method is that TEPH is a multicomponent mixture and is quantitated by summing the entire range, rather than individual peaks. All positive results were quantitated using the responses from the initial calibration curve using the external standard technique. Also, a confirmation column is not used, because the analyte is a multicomponent mixture and the specific carbon range of the peaks detected is specified on the individual sample reporting forms.
4. All initial and continuing calibration criteria were met.
5. The method blank associated with this project was below the MDL for diesel range organics.
6. All laboratory control sample and laboratory control sample duplicate recoveries and RPDs were within the acceptance criteria.

7. Matrix spikes and matrix spike duplicates could not be performed because of insufficient sample. A laboratory control sample and laboratory control sample duplicate were performed instead.
8. All samples were extracted and analyzed within the established holding time.
9. All surrogate recoveries were within the acceptance criteria.
10. The samples were analyzed at a dilution in order to bring the target analyte within the calibration range of the instrument. The reporting limits have been adjusted accordingly.
11. Manual integrations are performed when needed to provide consistent and defensible data following the guidelines in Paragon Analytics Standard Operating Procedure 939 Revision 3. Whenever manual integrations are performed, before and after chromatograms of the peak that was manually integrated are included in the report along with the reason why the re-integration was necessary.

The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, Paragon Analytics certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.

Mindy Norton  
Mindy Norton  
Organics Primary Data Reviewer

12.1.08  
Date

Joe Norton  
Organics Final Data Reviewer

12.2.08  
Date

***Paragon Analytics  
Data Qualifier Flags  
Fuels***

- G:** This flag indicates that a pattern resembling gasoline was detected in this sample.
- D:** This flag indicates that a pattern resembling diesel was detected in this sample.
- M:** This flag indicates that a pattern resembling motor oil was detected in this sample.
- C:** This flag indicates that a pattern resembling crude oil was detected in this sample.
- 4:** This flag indicates that a pattern resembling JP-4 was detected in this sample.
- 5:** This flag indicates that a pattern resembling JP-5 was detected in this sample.
- H:** This flag indicates that the fuel pattern was in the heavier end of the retention time window for the analyte of interest.
- L:** This flag indicates that the fuel pattern was in the lighter end of the retention time window for the analyte of interest.
- Z:** This flag indicates that a significant fraction of the reported result did not resemble the patterns of any of the following petroleum hydrocarbon products:  
gasoline  
JP-8  
diesel  
mineral spirits  
motor oil  
Stoddard solvent  
bunker C

Multiple flags may be used to indicate the presence of more than one product or component.

***Paragon Analytics, Inc.***  
***Data Qualifier Flags***  
***Chromatography and Mass Spectrometry***

- U or ND:** This flag indicates that the compound was analyzed for but not detected.
- J:** This flag indicates an estimated value. This flag is used as follows : (1) when estimating a concentration for tentatively identified compounds (TICs) where a 1:1 response is assumed; (2) when the mass spectral and retention time data indicate the presence of a compound that meets the volatile and semivolatile GC/MS identification criteria, and the result is less than the reporting limit (RL) but greater than the method detection limit (MDL); (3) when the data indicate the presence of a compound that meets the identification criteria, and the result is less than the RL but greater than the MDL; and (4) the reported value is estimated.
- B:** This flag is used when the analyte is detected in the associated method blank as well as in the sample. It indicates probable blank contamination and warns the data user. This flag shall be used for a tentatively identified compound (TIC) as well as for a positively identified target compound.
- E:** This flag identifies compounds whose concentration exceeds the upper level of the calibration range.
- A:** This flag indicates that a tentatively identified compound is a suspected aldol-condensation product.
- X:** This flag indicates that the analyte was diluted below an accurate quantitation level.
- \*:** This flag indicates that a spike recovery is outside the control criteria.
- +:** This flag indicates that the relative percent difference (RPD) exceeds the control criteria.

# Paragon Analytics

## Sample Number(s) Cross-Reference Table

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**Paragon OrderNum:** 0811110

**Client Name:** Cordilleran Compliance Services, Inc.

**Client Project Name:** Rulison Area Well monitoring

**Client Project Number:**

**Client PO Number:**

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Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
A11-15D	0811110-1		WATER	13-Nov-08	8:40
A11-15B	0811110-2		WATER	13-Nov-08	8:30



# Paragon Analytics

A Division of DataChem Laboratories, Inc.

225 Commerce Drive Fort Collins, CO 80524  
800-443-1511 or (970) 490-1511 (970) 490-1522 Fax

Accession Number (LAB ID)

Chain-of-Custody

Date

11/13/08

Page

1

of

1

Originator: Retain pink copy!

Project Name/No.: <u>BULLSON AREA WELL MONITORING</u>				Sampler(s): <u>TPD</u>		Turnaround (circle one) <u>Standard</u> or Rush (Due _____)		Dispose: Date _____ or Return to Client																											
Report To: <u>JAMES HIX</u> Phone: <u>(303) 237 2072</u> Fax: <u>(303) 237 2659</u> E-mail: <u>jameshix@cordcomp.com</u> Company: <u>Cordilleran Compliance Services, Inc.</u> Address: <u>826 2 1/2 Road</u> <u>4690 TABLE MOUNTAIN DR. #200</u> <u>GOLDEN, CO 80403</u>				Circle method (right): provide additional information as needed (comments).																															
Sample ID	Date	Time *	Lab ID	Matrix	Preservative (indicate type... HCl, etc.)	No. of Containers	VOCs	BTEX (only) - MTBE	SVOCs	OC Pesticides	PCBs	Herbicides	Explosives	TCLP Organics SW1311	TCLP Metals SW1311 Hg	Total Metals by ICP Hg	Dissolved Metals by ICP Hg	Total Metals by ICP/MS	Dissolved Metals by ICP/MS	Hexavalent Chromium	Inorganic Anions	Solids:	pH	TPH	Gross Alpha / Beta	Actinides by Paragon SOP	Tritium	Total Alpha-Emitting Radium	Radium 226	Radium 228	Strontium 90 (Total RadioSr)	Gamma Isotopes	Radon 222	PSL	NH <sub>4</sub> <sup>+</sup> , NO <sub>3</sub> <sup>-</sup> , T-Ac <sub>2</sub>
A11-15D	11/13/08	0840	(1)	W	H <sub>2</sub> O	17		X										X		X			X	X		X						X			
A11-15B	11/13/08	0830	(2)	W	"	17		X										X		X			X	X		X						X			
<div style="border: 1px solid black; height: 100px; width: 100%;"></div>																																			
* Time Zone: EST CST MST PST Matrix Key: O = oil, S = soil, NS = non-soil solid, W = water, L = liquid, E = extract, F = filter Comments:							Relinquished By: (1) Signature: <u>[Signature]</u> Printed Name: <u>T.M. DOBRANSKY</u> Date: <u>11/13/08</u> Time: <u>1600</u> Company: <u>Cordilleran Compliance</u>										Relinquished By: (2) Signature: _____ Printed Name: _____ Date: _____ Time: _____ Company: _____																		
Received By: (1) Signature: <u>[Signature]</u> Printed Name: <u>BO-ckane</u> Date: <u>11/14/08</u> Time: <u>0950</u> Company: <u>PA</u>							Received By: (2) Signature: _____ Printed Name: _____ Date: _____ Time: _____ Company: _____																												

## CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: CondillieranWorkorder No: 0811110Project Manager: LSInitials: oo Date: 11-14-08

1. Does this project require any special handling in addition to standard Paragon procedures?		YES	<u>NO</u>
2. Are custody seals on shipping containers intact?	NONE	<u>YES</u>	NO
3. Are Custody seals on sample containers intact?	NONE	<u>YES</u>	NO
4. Is there a COC (Chain-of-Custody) present or other representative documents?		<u>YES</u>	NO
5. Are the COC and bottle labels complete and legible?		<u>YES</u>	NO
6. Is the COC in agreement with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)		<u>YES</u>	NO
7. Were airbills / shipping documents present and/or removable?	DROP OFF	<u>YES</u>	NO
8. Are all aqueous samples requiring preservation preserved correctly? (excluding volatiles)	N/A	<u>YES</u>	<u>NO</u>
9. Are all aqueous non-preserved samples pH 4-9?	N/A	<u>YES</u>	NO
10. Is there sufficient sample for the requested analyses?		<u>YES</u>	NO
11. Were all samples placed in the proper containers for the requested analyses?		<u>YES</u>	NO
12. Are all samples within holding times for the requested analyses?		<u>YES</u>	NO
13. Were all sample containers received intact? (not broken or leaking, etc.)		<u>YES</u>	NO
14. Are all samples requiring no headspace (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u>✓</u> < green pea <u>   </u> > green pea	N/A	YES	<u>NO</u>
15. Do perchlorate LCMS-MS samples have headspace? (at least 1/3 of container required)	<u>N/A</u>	YES	NO
16. Were samples checked for and free from the presence of residual chlorine? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	<u>N/A</u>	YES	NO
17. Were the samples shipped on ice?		<u>YES</u>	NO
18. Were cooler temperatures measured at 0.1-6.0°C?	IR gun used*: <u>#2</u> <u>#4</u>	RAD ONLY <u>YES</u>	NO
Cooler #: <u>1</u>			
Temperature (°C): <u>3.4</u>			
No. of custody seals on cooler: <u>1</u>			
External µR/hr reading: <u>14</u>			
Background µR/hr reading: <u>13</u>			
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> / NO / NA (If no. see Form 008.)			

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16.

Headspace Bottle # (1) 1, 2, 3, 6, 7, 9  
# 2 - 1, 2

Slime layer in -1-15 & -1-16 (Organic?)

If applicable, was the client contacted? YES / NO / NA Contact: J. Hix Date/Time: 11/12/08Project Manager Signature / Date: [Signature] 11/12/08

\*IR Gun #2: Oakton. SN 29922500201-0066

\*IR Gun #4: Oakton. SN 2372220101-0002

CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: Condillivan

Workorder No: 0811110

Project Manager: LS

Initials: as Date: 11-14-08

Additional Information:

Was the laboratory directed to proceed with the analysis of any samples yielding the presence of residual chlorine? YES / NO / NA

NOTE:

No pH adjustments shall be made without prior consent of Project Manager. After pH adjustments, hold metals and radchem samples  $\geq 24$  hrs. before analysis.

Was the pH of any sample adjusted by the laboratory? YES (See Table below) / NO

pH Excursion:

Paragon Sample ID	Client Sample ID	Initial pH	Final pH	Reagent Used	Volume Added (mL)	Lot No. of Reagent	Requested Analysis	Initials / Date / Time
-1-12		7	1.6	conc HNO <sub>3</sub>	1mL			as 11/14/08 1030
-1-15		↓	↓	↓	↓			↓
-1-16		↓	↓	↓	↓			↓
-1-17		↓	↓	↓	↓			↓
-2-12		↓	↓	↓	↓			↓
-2-15		↓	↓	↓	↓			↓
-2-16		↓	↓	↓	↓			↓
-2-17		↓	↓	↓	↓			↓

If applicable, was the client contacted? YES / NO / NA Contact: \_\_\_\_\_ Date/Time: \_\_\_\_\_

Project Manager Signature / Date: LS 11/17/08



ORIGIN ID: GJTA (970) 270-2986  
TIM DOBRANSKY  
CORDILLERAN COMPLIANCE SERVICES, IN  
826 21 1/2 ROAD

GRAND JUNCTION, CO 81505  
UNITED STATES US

Ship Date: 13NOV08  
ActWgt: 20.0 LB MAN  
System#: 390082/CAFE2358  
Account: S 235727234

TO

PARAGON ANALYTICS  
225 COMMERCE DRIVE

FORT COLLINS, CO 80524

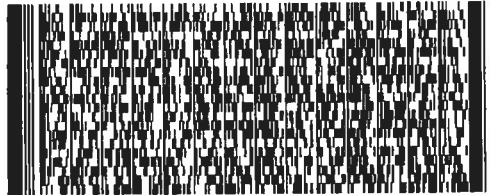
(800) 443-1511

**FedEx**  
Express



CL5050107/22/23

Ref: 8360



Delivery Address  
Barcode

BILL SENDER

PRIORITY OVERNIGHT

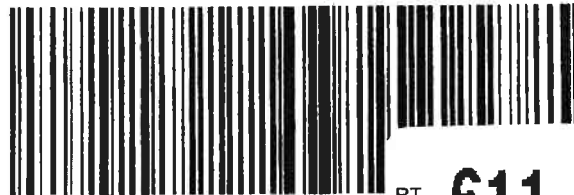
TRK# 9660 0451 2332 Form 0201

**FRI**  
Deliver By:  
14NOV08

**DEN** AA

80524 -CO-US

**72 FTCA**



RT

**611 A**

FZ

2332  
11.14

## Analytical Results

# Diesel Range Organics

Method SW8015MB

Method Blank

Lab Name: Paragon Analytics

Work Order Number: 0811110

Client Name: Cordilleran Compliance Services, Inc.

ClientProject ID: Rulison Area Well monitoring

Lab ID: EX081118-4MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 18-Nov-08

Date Analyzed: 20-Nov-08

Prep Method: METHOD

Prep Batch: EX081118-4

QCBatchID: EX081118-4-1

Run ID: HCD081120-3A

Cleanup: NONE

Basis: N/A

File Name: F3F32996

Sample Aliquot: 160ml

Final Volume: 4ml

Result Units: MG/L

Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
68334-30-5	Diesel Range Organics	1	0.5	0.5	U	

## Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
84-15-1	O-TERPHENYL	1.09		1.25	87	57 - 132

Data Package ID: HCD0811110-1

Date Printed: Monday, December 01, 2008

Paragon Analytics

LIMS Version: 6.212A

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# Diesel Range Organics

Method SW8015MB

## Sample Results

Lab Name: Paragon Analytics

Work Order Number: 0811110

Client Name: Cordilleran Compliance Services, Inc.

ClientProject ID: Rulison Area Well monitoring

<b>Field ID:</b> A11-15D	<b>Sample Matrix:</b> WATER	<b>Prep Batch:</b> EX081118-4	<b>Sample Aliquot:</b> 160 ml
<b>Lab ID:</b> 0811110-1	<b>% Moisture:</b> N/A	<b>QCBatchID:</b> EX081118-4-1	<b>Final Volume:</b> 4 ml
	<b>Date Collected:</b> 13-Nov-08	<b>Run ID:</b> HCD081120-3A	<b>Result Units:</b> MG/L
	<b>Date Extracted:</b> 18-Nov-08	<b>Cleanup:</b> NONE	<b>Clean DF:</b> 1
	<b>Date Analyzed:</b> 20-Nov-08	<b>Basis:</b> As Received	
	<b>Prep Method:</b> METHOD	<b>File Name:</b> F3F32999	

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
68334-30-5	Diesel Range Organics	2	67	1	L	

## Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
84-15-1	O-TERPHENYL	1.06		1.25	84	57 - 132

The chromatogram for Diesel Range Organics indicates the presence of hydrocarbons in the range of C7-C18.

Data Package ID: HCD0811110-1

Date Printed: Monday, December 01, 2008

Paragon Analytics

LIMS Version: 6.212A

Page 1 of 2

# Diesel Range Organics

Method SW8015MB

## Sample Results

Lab Name: Paragon Analytics

Work Order Number: 0811110

Client Name: Cordilleran Compliance Services, Inc.

ClientProject ID: Rulison Area Well monitoring

<b>Field ID:</b> A11-15B	<b>Sample Matrix:</b> WATER	<b>Prep Batch:</b> EX081118-4	<b>Sample Aliquot:</b> 160 ml
<b>Lab ID:</b> 0811110-2	<b>% Moisture:</b> N/A	<b>QCBatchID:</b> EX081118-4-1	<b>Final Volume:</b> 4 ml
	<b>Date Collected:</b> 13-Nov-08	<b>Run ID:</b> HCD081120-3A	<b>Result Units:</b> MG/L
	<b>Date Extracted:</b> 18-Nov-08	<b>Cleanup:</b> NONE	<b>Clean DF:</b> 1
	<b>Date Analyzed:</b> 20-Nov-08	<b>Basis:</b> As Received	
	<b>Prep Method:</b> METHOD	<b>File Name:</b> F3F33000	

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
68334-30-5	Diesel Range Organics	2	100	1	L	

## Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
84-15-1	O-TERPHENYL	1.09		1.25	87	57 - 132

The chromatogram for Diesel Range Organics indicates the presence of hydrocarbons in the range of C7-C28.

Data Package ID: HCD0811110-1

Date Printed: Monday, December 01, 2008

Paragon Analytics

LIMS Version: 6.212A

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## Supporting QA/QC Data

# Surrogate Summary for Diesel Range Organics

Method SW8015MB

Lab Name: Paragon Analytics

Work Order Number: 0811110

Client Name: Cordilleran Compliance Services, Inc.

ClientProject ID: Rulison Area Well monitoring

PrepBatchID: EX081118-4

QC Batch ID: EX081118-4-1

Date Extracted: 11/18/2008

Surrogate Compound	Control Limits	
	Lower	Upper
o-terphenyl	57	132

Lab ID	Client Sample ID	Date Collected	Date Received	% Recovery
EX081118-4MB	XXXXXXX	11/18/2008	11/14/2008	87
EX081118-4LCS	XXXXXXX	11/18/2008	11/14/2008	88
EX081118-4LCSD	XXXXXXX	11/18/2008	11/14/2008	87
0811110-1	A11-15D	11/13/2008	11/14/2008	84
0811110-2	A11-15B	11/13/2008	11/14/2008	87

Data Package ID: HCD0811110-1

Date Printed: Monday, December 01, 2008

Paragon Analytics

LIMS Version: 6.212A

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# Diesel Range Organics

## Method SW8015MB

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0811110

Client Name: Cordilleran Compliance Services, Inc.

ClientProject ID: Rulison Area Well monitoring

Lab ID: EX081118-4LCS

Sample Matrix: WATER  
% Moisture: N/A  
Date Collected: N/A  
Date Extracted: 11/18/2008  
Date Analyzed: 11/20/2008  
Prep Method: METHOD

Prep Batch: EX081118-4  
QCBatchID: EX081118-4-1  
Run ID: HCD081120-3A  
Cleanup: NONE  
Basis: N/A  
File Name: F3F32997

Sample Aliquot: 160ml  
Final Volume: 4 ml  
Result Units: MG/L  
Clean DF: 1

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
68334-30-5	Diesel Range Organics	5	4.89	0.5		98	36 - 150%

Lab ID: EX081118-4LCSD

Sample Matrix: WATER  
% Moisture: N/A  
Date Collected: N/A  
Date Extracted: 11/18/2008  
Date Analyzed: 11/20/2008  
Prep Method: METHOD

Prep Batch: EX081118-4  
QCBatchID: EX081118-4-1  
Run ID: HCD081120-3A  
Cleanup: NONE  
Basis: N/A  
File Name: F3F32998

Sample Aliquot: 160ml  
Final Volume: 4 ml  
Result Units: MG/L  
Clean DF: 1

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
68334-30-5	Diesel Range Organics	5	4.91	0.5		98	20	0

### Surrogate Recovery LCS/LCSD

CASNO	Target Analyte	Spike Added	LCS % Rec.	LCS Flag	LCSD % Rec.	LCSD Flag	Control Limits
84-15-1	O-TERPHENYL	1.25	88		87		57 - 132

Data Package ID: HCD0811110-1

Date Printed: Monday, December 01, 2008

Paragon Analytics  
LIMS Version: 6.212A

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## Prep Batch ID: EX081118-4

Start Date: 11/18/08

End Date: 11/18/08

Concentration Method: NONE

Batch Created By: sjc

Start Time: 14:20

End Time: 14:30

Extract Method: METHOD

Date Created: 11/18/08

Prep Analyst: Sean Collins

Initial Volume Units: ml

Time Created: 15:30

Final Volume Units: ml

Validated By: sjc

Date Validated: 11/18/08

Time Validated: 15:01

Comments:

QC Batch ID: EX081118-4-1

Lab ID	QC Type	Field ID	Matrix	Date Collected	Initial Wt/Vol	Final Wt/Vol	Cleanup Method	Cleanup DF	Order Number
EX081118-4	MB	XXXXXX	WATER	XXXXXX	160	4	NONE	1	0811110
EX081118-4	LCS	XXXXXX	WATER	XXXXXX	160	4	NONE	1	0811110
EX081118-4	LCSD	XXXXXX	WATER	XXXXXX	160	4	NONE	1	0811110
0811110-1	SMP	A11-15D	WATER	11/13/2008	160	4	NONE	1	0811110
0811110-2	SMP	A11-15B	WATER	11/13/2008	160	4	NONE	1	0811110

In generating this benchsheet, prep analyst states that all aspects of sample preparation as set forth in the appropriate PAR SOP's (including Kuderna-Danish temperatures, proper flow settings on the N-evap, and final volumes) were properly adhered to (unless otherwise noted herein).

QC Types

CAR	Carrier reference sample	DUP	Laboratory Duplicate
LCS	Laboratory Control Sample	LCSD	Laboratory Control Sample Duplicat
MB	Method Blank	MS	Laboratory Matrix Spike
MSD	Laboratory Matrix Spike Duplicate	REP	Sample replicate
SMP	Field Sample	SYS	Sample Yield Spike

## Initial Calibration Summary

Paragon Analytics

Inst. ID: FUELS3

C:\HPCHEM\5\METHODS\D111208.M

Calibration Date:11/12/08

1 = F3F32866.D    2 = F3F32867.D    3 = F3F32868.D    4 = F3F32869.D  
 5 = F3F32870.D    6 = F3F32871.D    7 =    8 =

Compounds	Calibration Factors							
	1	2	3	4	5	6	7	8
TEPH	4621.57	4684.28	4844.43	4818.68	5769.21	9243.50		
o-terphenyl	5026.92	5090.10	5185.69	4930.10	5018.60	5157.50		

Compounds	Ave. CF	%RSD	Corr. r2	Curve fit type	Linear curve info			
					slope	y-int	x*x	
TEPH	5663.6	31.8	1.000	Linear	4599.013	141417.51	0.00	
o-terphenyl	5068.2	1.9	0.019	Average	5068.152	0.00	0.00	

Compounds	Calibration Level (µg/ml)							
	1	2	3	4	5	6	7	8
TEPH	5000	2000	1000	500	100	20		
o-terphenyl	500	200	100	50	10	2		

# Calibration Verification Summary

Paragon Analytics

Sample: 500ug/mL DRO ICV

Data File #1: C:\HPCHEM\5\DATA\11122008\F3F32872.D Column #1: DB-5.625

COMPOUND	Column #1			Nom Conc µg/ml	Col. #1	Col. #1
	Exp. RT (min)	Found RT (min)	Dev (min)		Conc µg/ml	%D
TEPH	15.000	15.000	0.000	500	475	5

*Est  
11/17/08*

# Calibration Verification Summary

Paragon Analytics

Sample: 1000ug/mL DRO CCV

Data File #1: C:\HPCHEM\5\DATA\11202008\F3F32995.D Column #1: DB-5.625

COMPOUND	Column #1			Nom Conc µg/ml	Col. #1	Col. #1
	Exp. RT (min)	Found RT (min)	Dev (min)		Conc µg/ml	%D
o-terphenyl	15.660	15.660	0.000	100	89	11
TEPH	15.000	15.000	0.000	1000	984	2

# Calibration Verification Summary

Paragon Analytics

Sample: 1000ug/mL DRO CCV2

Data File #1: C:\HPCHEM\5\DATA\11202008\F3F33007.D Column #1: DB-5.625

COMPOUND	Column #1			Nom Conc µg/ml	Col. #1	Col. #1
	Exp. RT (min)	Found RT (min)	Dev (min)		Conc µg/ml	%D
o-terphenyl	15.660	15.660	0.000	100	96	4
TEPH	15.000	15.000	0.000	1000	1024	2

EB  
11/21/08

## Supporting Raw Data

Instrument Name Fuels 3

## Paragon Analytics

Logbook No./Page 369308Sequence File: C:\HPCHEM\5\SEQUENCE\ 11122008.5Date Analyzed 11/12 - 11/15/08Operator SKSGC Method TEPH12Data Path: C:\HPCHEM\5\DATA\ 11122008Analytical Method 8015m PROSOP 406 Rev. 13Hexane Lot 080013 Dichloromethane Lot NA

Reviewed by / date \_\_\_\_\_

Form 531r2.xls (9/14/2001)

Vial	Data File	Method	Sample Name	Comments	It?
1	F3F32865	D111208	HYDROCARBON MIX	25µL ST071213-5 + 475µL Hexane	NO
2	F3F32866	D111208	5000ug/mL DRO ICAL	ST080516-1	
3	F3F32867	D111208	2000ug/mL DRO ICAL	400µL ST080516-1 + 600µL Hexane	
4	F3F32868	D111208	1000ug/mL DRO ICAL	200µL ST080516-1 + 800µL Hexane	
5	F3F32869	D111208	500ug/mL DRO ICAL	100µL ST080516-1 + 900µL Hexane	
6	F3F32870	D111208	100ug/mL DRO ICAL	20µL ST080516-1 + 980µL Hexane	
7	F3F32871	D111208	20ug/mL DRO ICAL	4µL ST080516-1 + 996µL Hexane	
8	F3F32872	D111208	500ug/mL DRO ICV	Pass 100µL ST080825-7 + 900µL Hexane	
9	F3F32873	D111208	EX081030-8MB		
10	F3F32874	D111208	EX081030-8LCS		
11	F3F32875	D111208	EX081030-8LCSD		
12	F3F32876	D111208	EX081029-6MB		
13	F3F32877	D111208	EX081029-6LCS		
14	F3F32878	D111208	EX081029-6LCSD		
15	F3F32879	D111208	0810235-12		
16	F3F32880	D111208	Hexane		
17	F3F32881	D111208	Hexane		
18	F3F32882	D111208	Hexane		
19	F3F32883	D111208	1000ug/mL DRO CCV1	Pass 200µL ST080516-1 + 800µL Hexane	
20	F3F32884	D111208	0810235-2		
21	F3F32885	D111208	0810235-3		
22	F3F32886	D111208	0810235-5		
23	F3F32887	D111208	0810235-6		
24	F3F32888	D111208	0810235-7		
25	F3F32889	D111208	0810235-7MS		
26	F3F32890	D111208	0810235-7MSD		
27	F3F32891	D111208	0810235-10		
28	F3F32892	D111208	0810235-11		
29	F3F32893	D111208	0810235-18		
30	F3F32894	D111208	1000ug/mL DRO CCV2	Pass 200µL ST080516-1 + 800µL Hexane	
31	F3F32895	D111208	0810235-16 2X		

CONTINUED ON NEXT PAGE

Instrument Name Fuels 3

## Paragon Analytics

Logbook No./Page 369315

Sequence File: C:\HPCHEM\5\SEQUENCE\11202008.AS

Date Analyzed 11/20/08Operator SEGC Method TECH 12

Data Path: C:\HPCHEM\5\DATA\11202008

Analytical Method 8015M DROSOP 406 Rev. 13Hexane Lot 080013 Dichloromethane Lot N/A

Reviewed by / date \_\_\_\_\_

Form 531r2.xls (9/14/2001)

Vial	Data File	Method	Sample Name	Comments	rr?
100	F3F32994	D111208	BLANK		ND
1	F3F32995	D111208	1000ug/mL DRO CCV	200µL ST081119-3 + 800µL Hexane	
2	F3F32996	D111208	EX081118-4MB		
3	F3F32997	D111208	EX081118-4LCS		
4	F3F32998	D111208	EX081118-4LCSD		
5	F3F32999	D111208	0811110-1 2X	150µL Sample + 150µL Hexane	
6	F3F33000	D111208	0811110-2 2X	150µL Sample + 150µL Hexane	
7	F3F33001	D111208	HEXANE		
8	F3F33002	D111208	EX081119-1MB		
9	F3F33003	D111208	EX081119-1LCS		
10	F3F33004	D111208	EX081119-1LCSD		
11	F3F33005	D111208	0811116-2		
12	F3F33006	D111208	0811116-1		
13	F3F33007	D111208	1000ug/mL DRO CCV2	200µL ST081119-3 + 800µL Hexane	
14	F3F33008	D111208	0811117-2		
15	F3F33009	D111208	0811117-1		
16	F3F33010	D111208	HEXANE		
17	F3F33011	D111208	1000ug/mL DRO CCV3	200µL ST081119-3 + 800µL Hexane	

80/100/110  
11/21/08



## Calibration Raw Data

# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32865.D  
 Acq On : 12 Nov 08 01:42 PM  
 Sample : HYDROCARBON MIX  
 Misc : 25uL ST071213-5 + 475uL HEXANE  
 Quant Time: Nov 12 14:24 19108

Vial: 1  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 12:55:51 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	0.00	0	N.D. µg/ml
	Recovery	=	0.00%
Target Compounds			
1) H TEPH	15.00	1493987	340.97 µg/ml

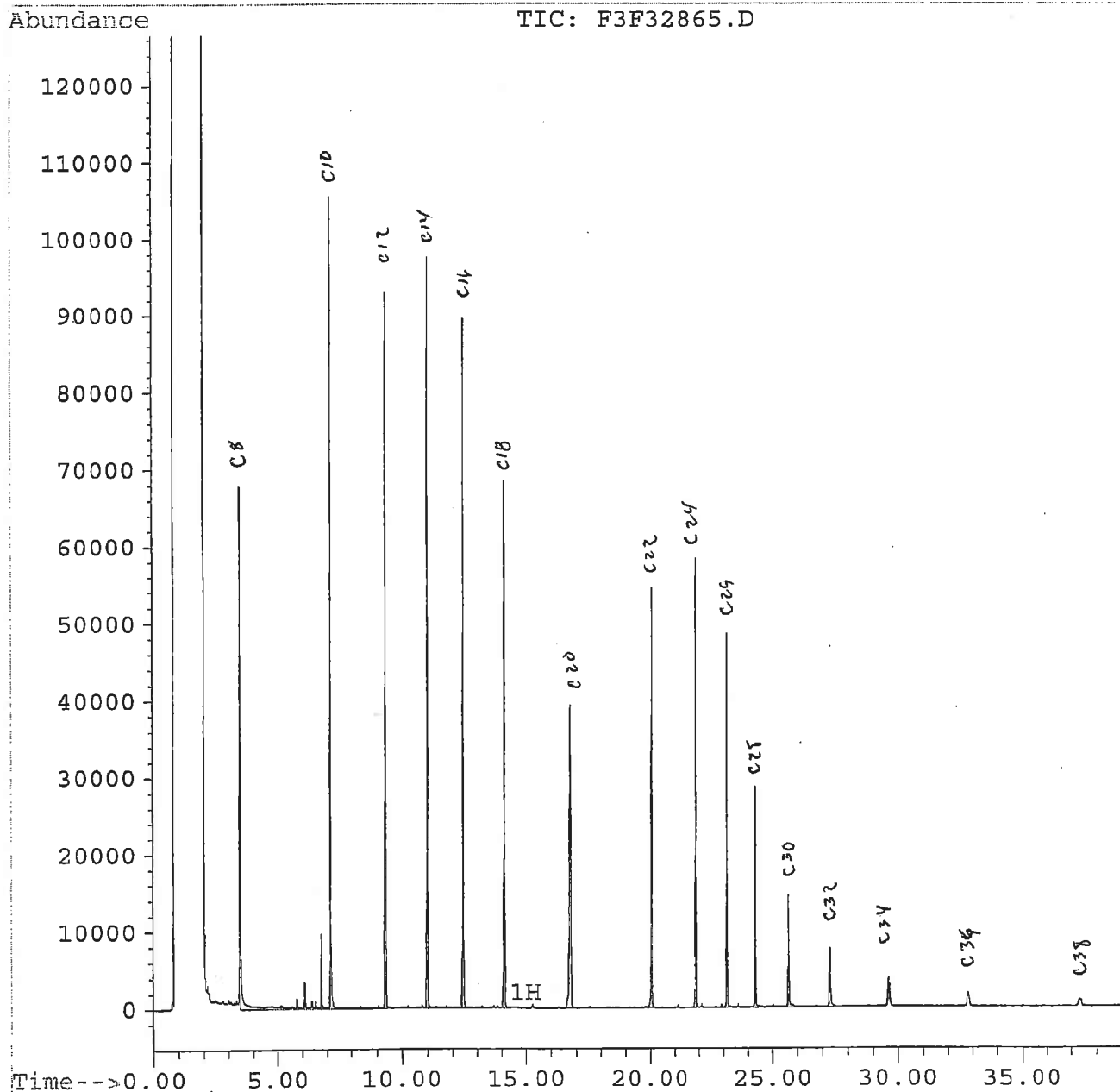
# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32865.D  
 Acq On : 12 Nov 08 01:42 PM  
 Sample : HYDROCARBON MIX  
 Misc : 25uL ST071213-5 + 475uL HEXANE  
 Quant Time: Nov 12 14:24 19108

Vial: 1  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 12:55:51 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5μm  
 Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32866.D  
 Acq On : 12 Nov 08 02:29 PM  
 Sample : 5000ug/mL DRO ICAL  
 Misc : ST080516-1  
 Quant Time: Nov 12 15:13 19108

Vial: 2  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 15:13:37 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5μm  
 Signal Info. : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.86f	2513462	500.00 µg/ml
	Recovery	= 1000.00%	
Target Compounds			
1) H TEPH	15.00	23107873	5000.00 µg/ml

*Ed 11/12/08*

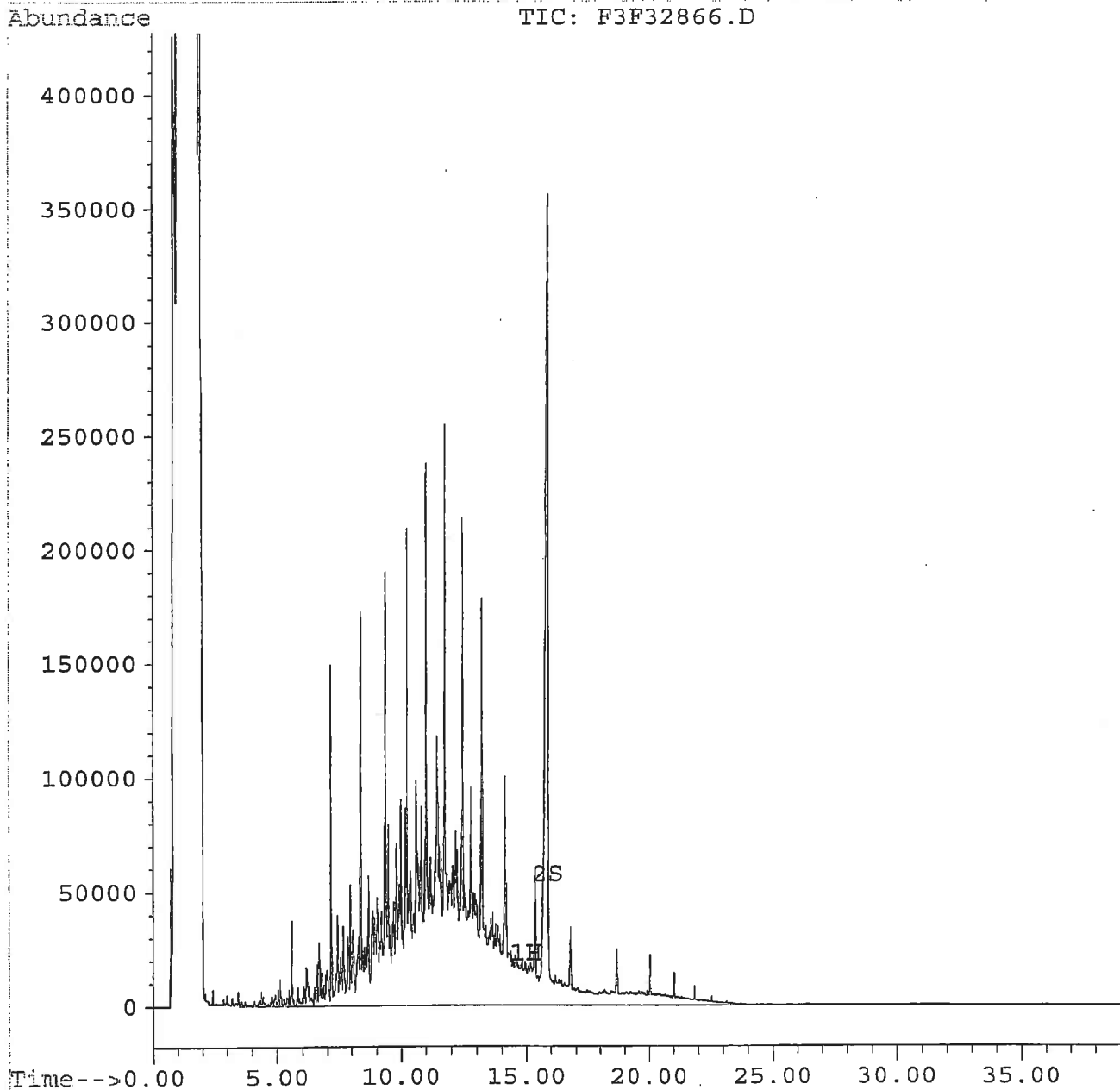
Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32866.D  
Acq On : 12 Nov 08 02:29 PM  
Sample : 5000ug/mL DRO ICAL  
Misc : ST080516-1  
Quant Time: Nov 12 15:13 19108

Vial: 2  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Wed Nov 12 15:13:37 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32867.D  
 Acq On : 12 Nov 08 03:15 PM  
 Sample : 2000ug/mL DRO ICAL  
 Misc : 400uL ST080516-1 + 600uL HEXANE  
 Quant Time: Nov 12 16:01 19108

Vial: 3  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 15:13:37 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5μm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
<b>System Monitoring Compounds</b>			
2) S o-terphenyl	15.79	1018019	202.51 μg/ml
	Recovery	=	405.02%
<b>Target Compounds</b>			
1) H TEPH	15.00	9368563	2027.14 μg/ml

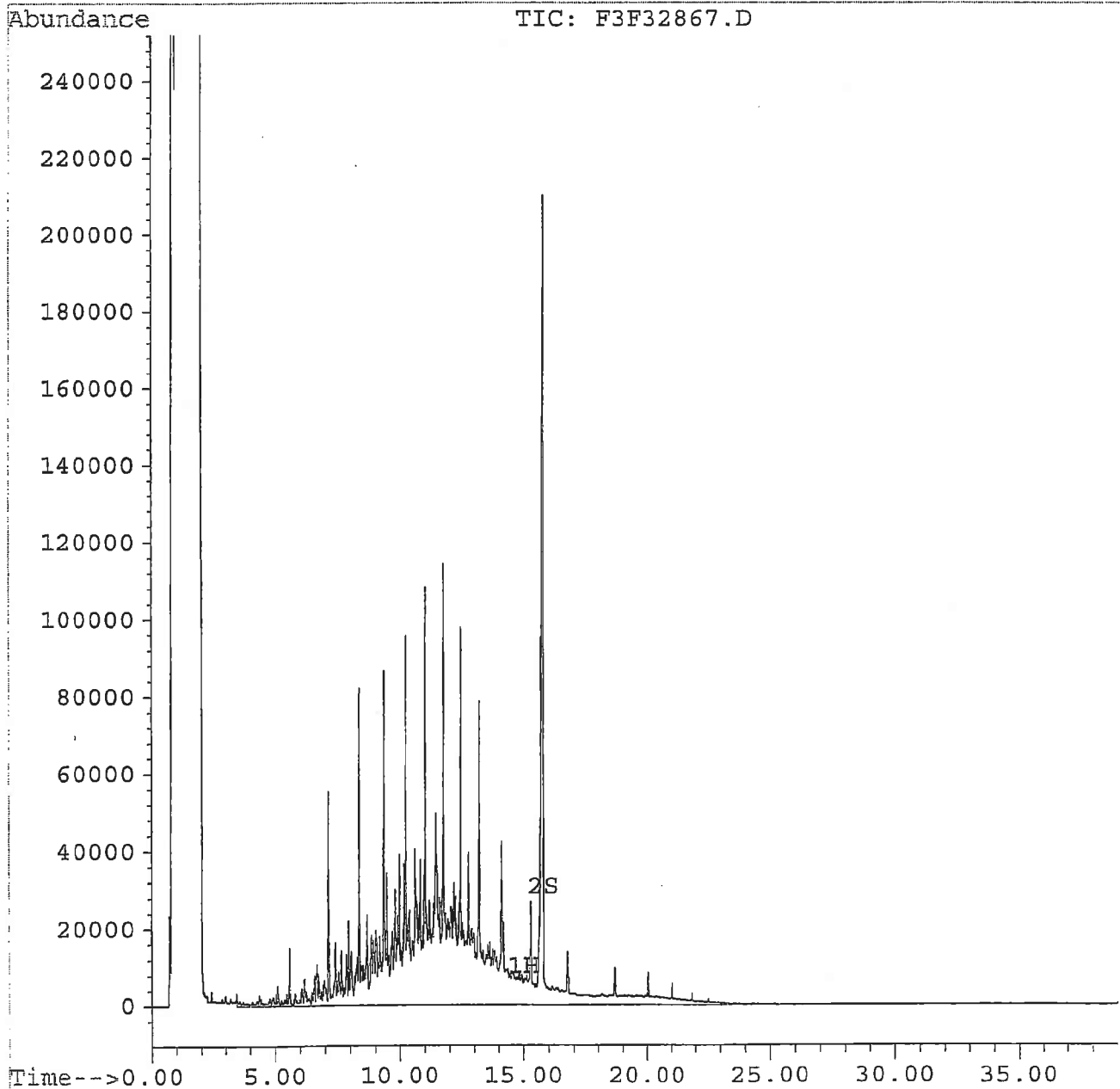
# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32867.D  
Acq On : 12 Nov 08 03:15 PM  
Sample : 2000ug/mL DRO ICAL  
Misc : 400uL ST080516-1 + 600uL HEXANE  
Quant Time: Nov 12 16:01 19108

Vial: 3  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Wed Nov 12 15:13:37 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32868.D  
 Acq On : 12 Nov 08 04:01 PM  
 Sample : 1000ug/mL DRO ICAL  
 Misc : 200uL ST080516-1 + 800uL HEXANE  
 Quant Time: Nov 12 16:43 19108

Vial: 4  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 16:01:31 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.75f	518569	102.51 µg/ml
	Recovery	=	205.02%
Target Compounds			
1) H TEPH	15.00	4844428	1041.16 µg/ml

(f)=RT Delta > 1/2 Window

F3F32868.D D111208.M

Wed Nov 12 16:43:16 2008

(m)=manual int.

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*edb*  
*11/12/08*



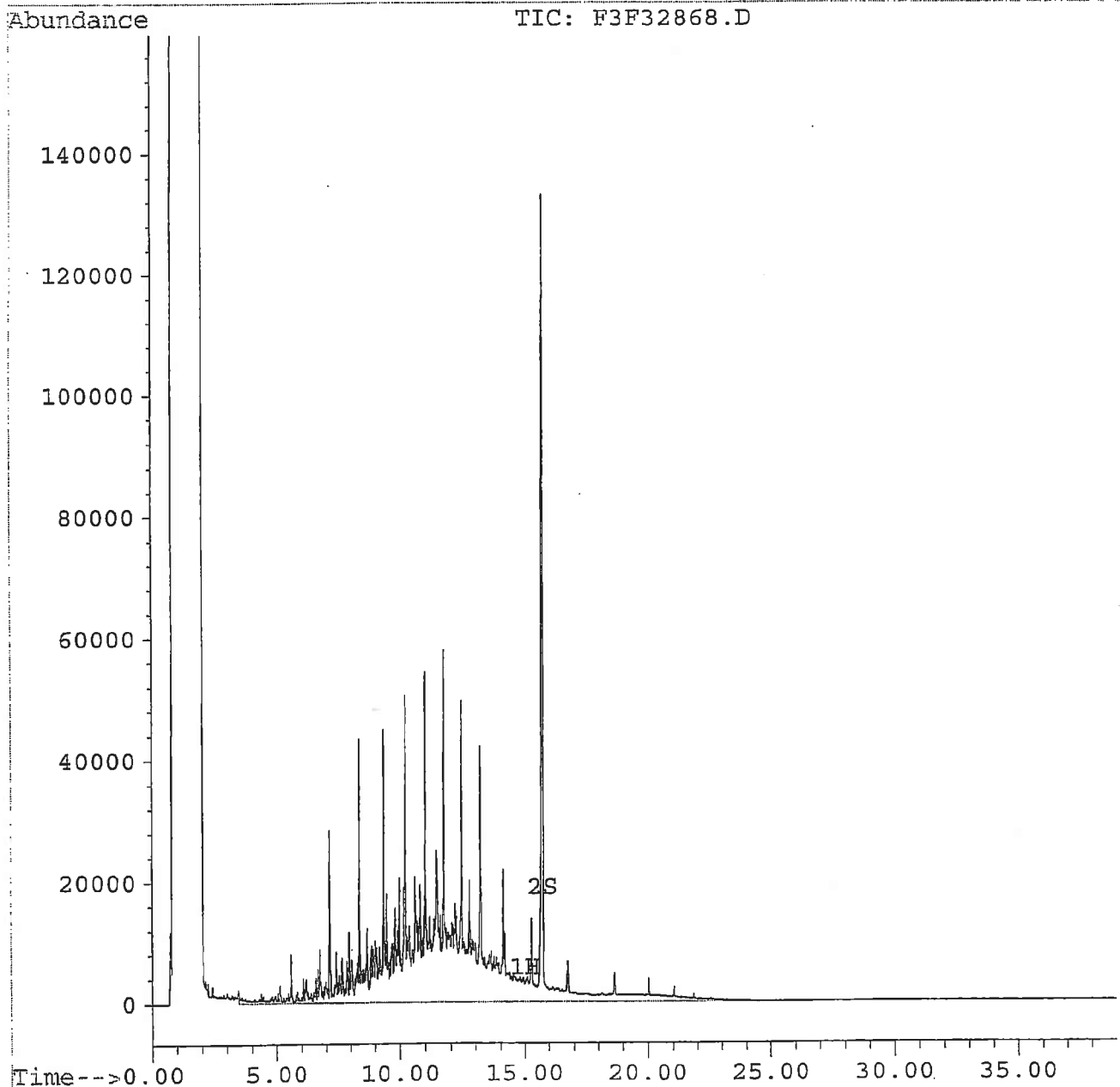
# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32868.D  
Acq On : 12 Nov 08 04:01 PM  
Sample : 1000ug/mL DRO ICAL  
Misc : 200uL ST080516-1 + 800uL HEXANE  
Quant Time: Nov 12 16:43 19108

Vial: 4  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Wed Nov 12 16:01:31 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32869.D  
 Acq On : 12 Nov 08 04:48 PM  
 Sample : 500ug/mL DRO ICAL  
 Misc : 100uL ST080516-1 + 900uL HEXANE  
 Quant Time: Nov 12 17:28 19108

Vial: 5  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 16:43:24 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5μm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
<b>System Monitoring Compounds</b>			
2) S o-terphenyl	15.73f	246505	48.33 μg/ml
	Recovery	=	96.66%
<b>Target Compounds</b>			
1) H TEPH	15.00	2409338	510.80 μg/ml

(f)=RT Delta > 1/2 Window

(m)=manual int.

F3F32869.D D111208.M

Wed Nov 12 17:28:52 2008

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edb  
11/17/08

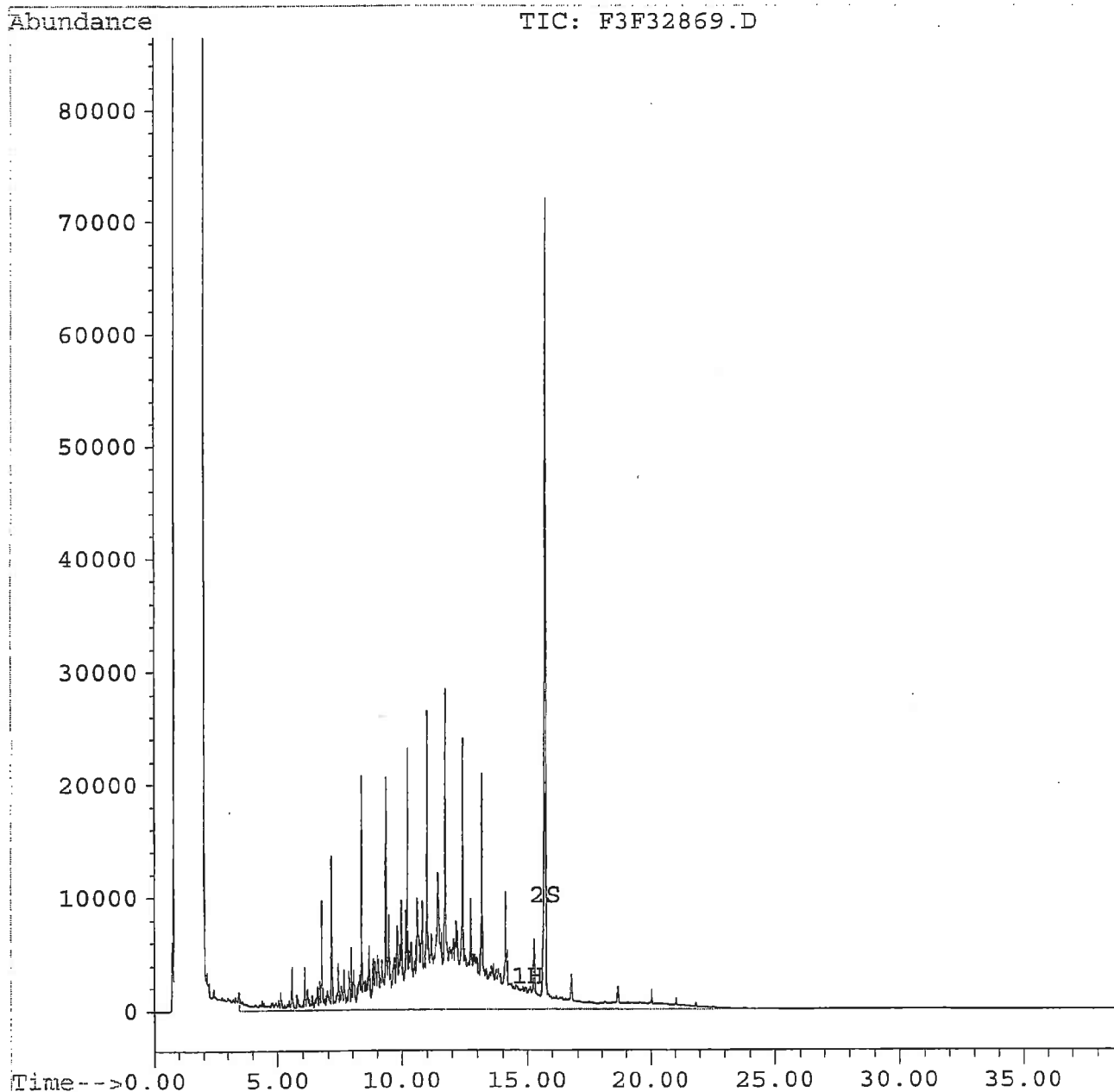
# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32869.D  
Acq On : 12 Nov 08 04:48 PM  
Sample : 500ug/mL DRO ICAL  
Misc : 100uL ST080516-1 + 900uL HEXANE  
Quant Time: Nov 12 17:28 19108

Vial: 5  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Wed Nov 12 16:43:24 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32870.D  
 Acq On : 12 Nov 08 05:34 PM  
 Sample : 100ug/mL DRO ICAL  
 Misc : 20uL ST080516-1 + 980uL HEXANE  
 Quant Time: Nov 12 18:27 19108

Vial: 6  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 17:29:44 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
<b>System Monitoring Compounds</b>			
2) S o-terphenyl	15.70f	50186	9.92 µg/ml
	Recovery	=	19.84%
<b>Target Compounds</b>			
1) H TEPH	15.00	576921	121.66 µg/ml

*Handwritten:* 8/17/05

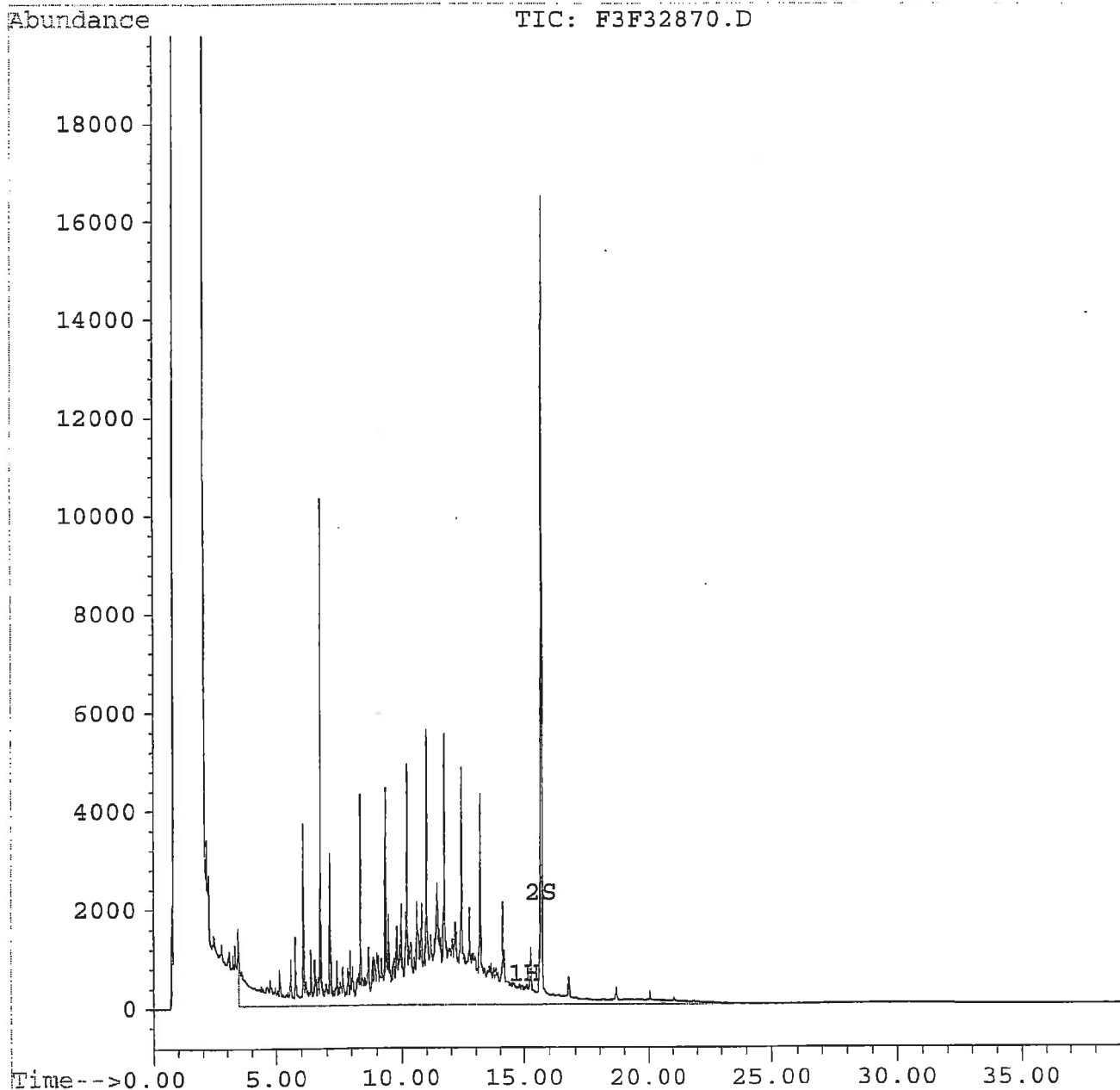
# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32870.D  
Acq On : 12 Nov 08 05:34 PM  
Sample : 100ug/mL DRO ICAL  
Misc : 20uL ST080516-1 + 980uL HEXANE  
Quant Time: Nov 12 18:27 19108

Vial: 6  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Wed Nov 12 17:29:44 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32871.D  
 Acq On : 12 Nov 08 06:20 PM  
 Sample : 20ug/mL DRO ICAL  
 Misc : 4uL ST080516-1 + 996uL HEXANE  
 Quant Time: Nov 12 19:09 19108

Vial: 7  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 19:09:08 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.70f	10315	2.04 µg/ml
	Recovery	=	4.08%
Target Compounds			
1) H TEPH	15.00	184870	17.37 µg/ml

(f)=RT Delta > 1/2 Window

(m)=manual int.

F3F32871.D D111208.M

Wed Nov 12 19:09:39 2008

8/17/08 Page 1  
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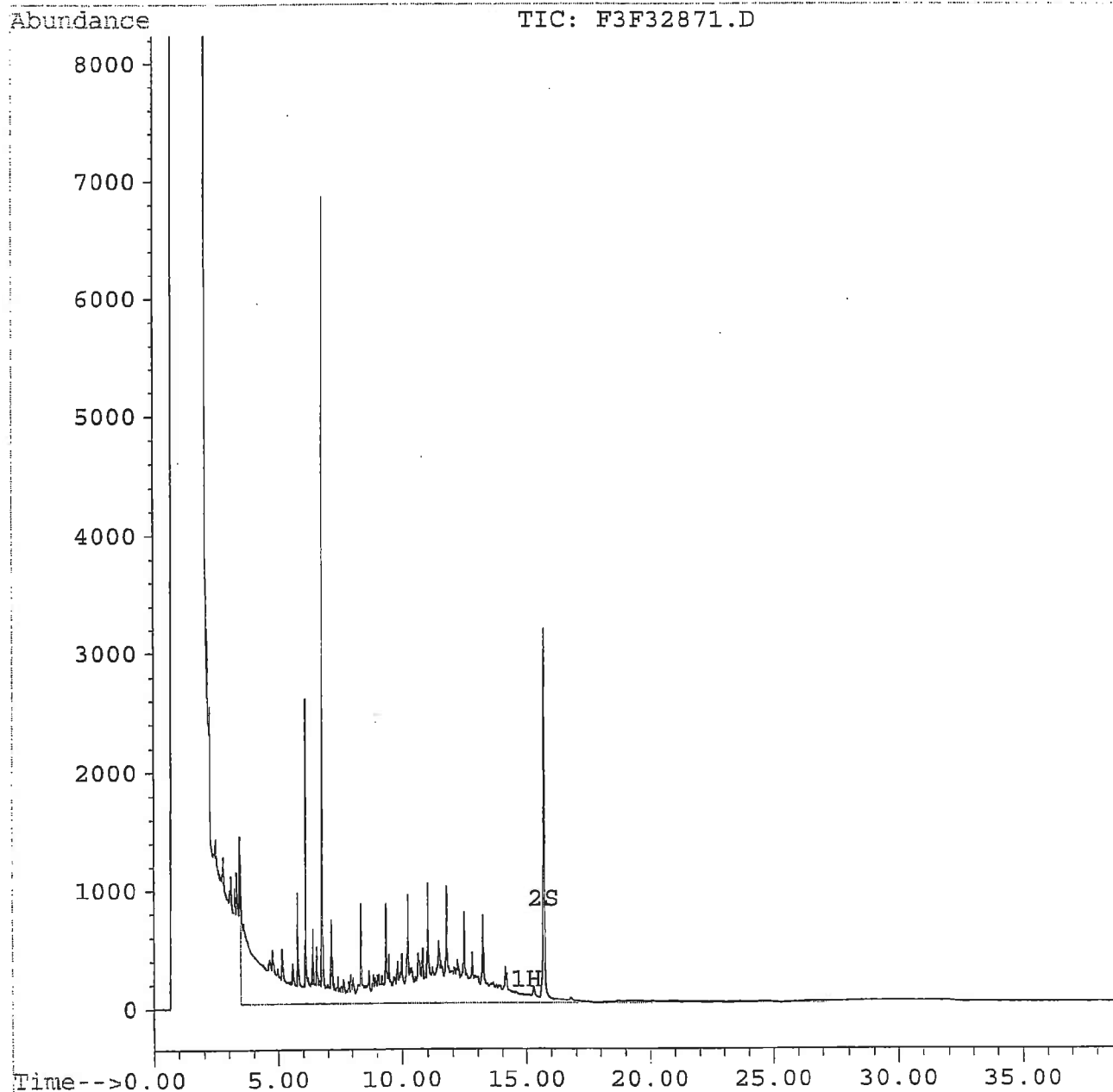
Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32871.D  
Acq On : 12 Nov 08 06:20 PM  
Sample : 20ug/mL DRO ICAL  
Misc : 4uL ST080516-1 + 996uL HEXANE  
Quant Time: Nov 12 19:09 19108

Vial: 7  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Wed Nov 12 19:09:08 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11122008\F3F32872.D  
 Acq On : 12 Nov 08 07:06 PM  
 Sample : 500ug/mL DRO ICV  
 Misc : 100uL ST080825-7 + 900uL HEXANE  
 Quant Time: Nov 13 9:05 19108

Vial: 8  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Wed Nov 12 20:36:13 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5μm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	0.00	0	N.D. μg/ml
	Recovery	=	0.00%
Target Compounds			
1) H TEPH	15.00	2326967	475.22 μg/ml



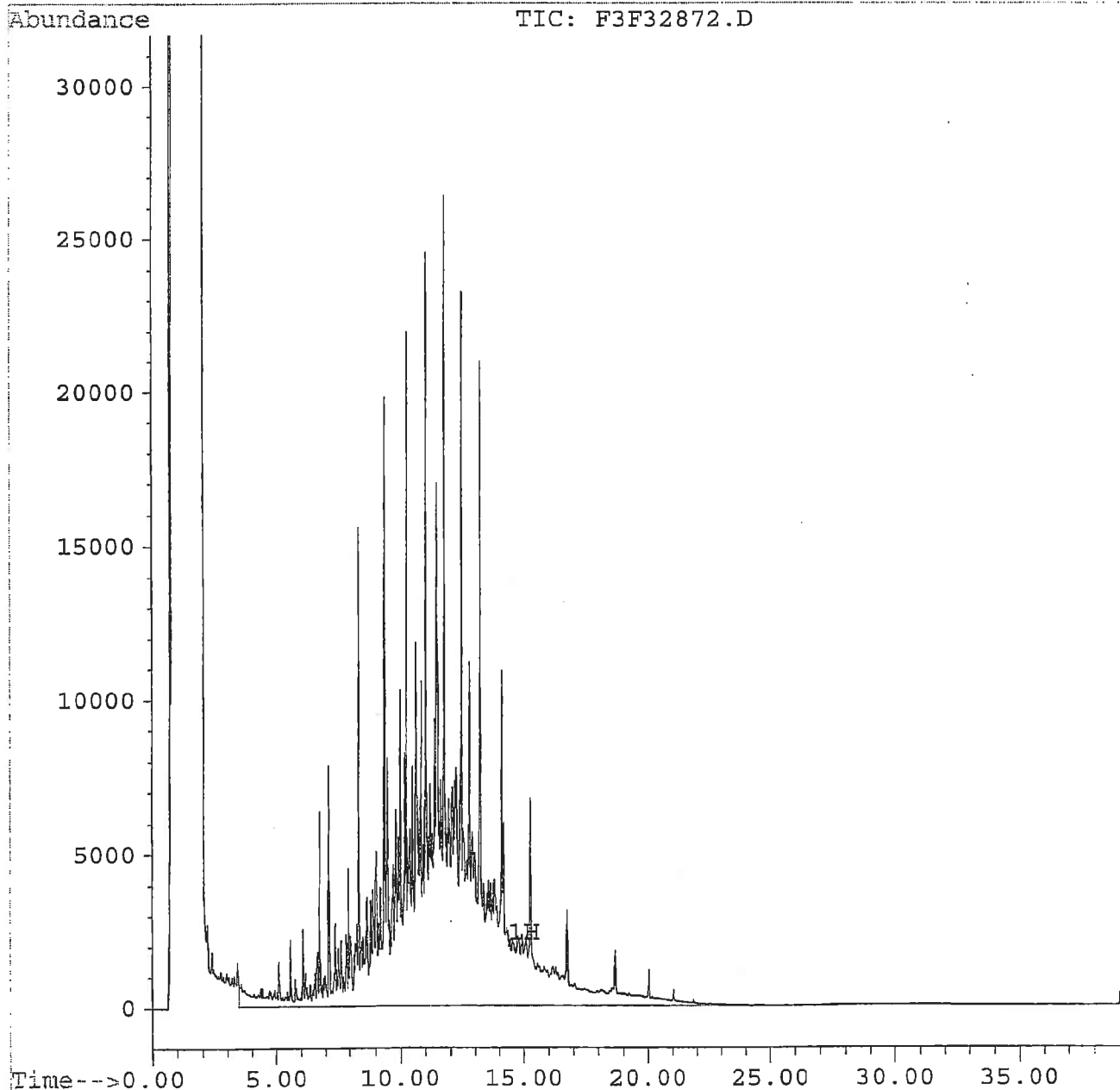
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Data File : C:\HPCHEM\5\DATA\11122008\F3F32872.D  
Acq On : 12 Nov 08 07:06 PM  
Sample : 500ug/mL DRO ICV  
Misc : 100uL ST080825-7 + 900uL HEXANE  
Quant Time: Nov 13 9:05 19108

Vial: 8  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Wed Nov 12 20:36:13 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



## Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32995.D  
Acq On : 20 Nov 08 03:57 PM  
Sample : 1000ug/mL DRO CCV  
Misc : 200uL ST081119-3 + 800uL HEXANE  
Quant Time: Nov 21 11:19 19108

Vial: 1  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.66	453547	89.49 µg/ml
	Recovery	=	178.98%
Target Compounds			
1) H TEPH	15.00	4668113	984.28 µg/ml

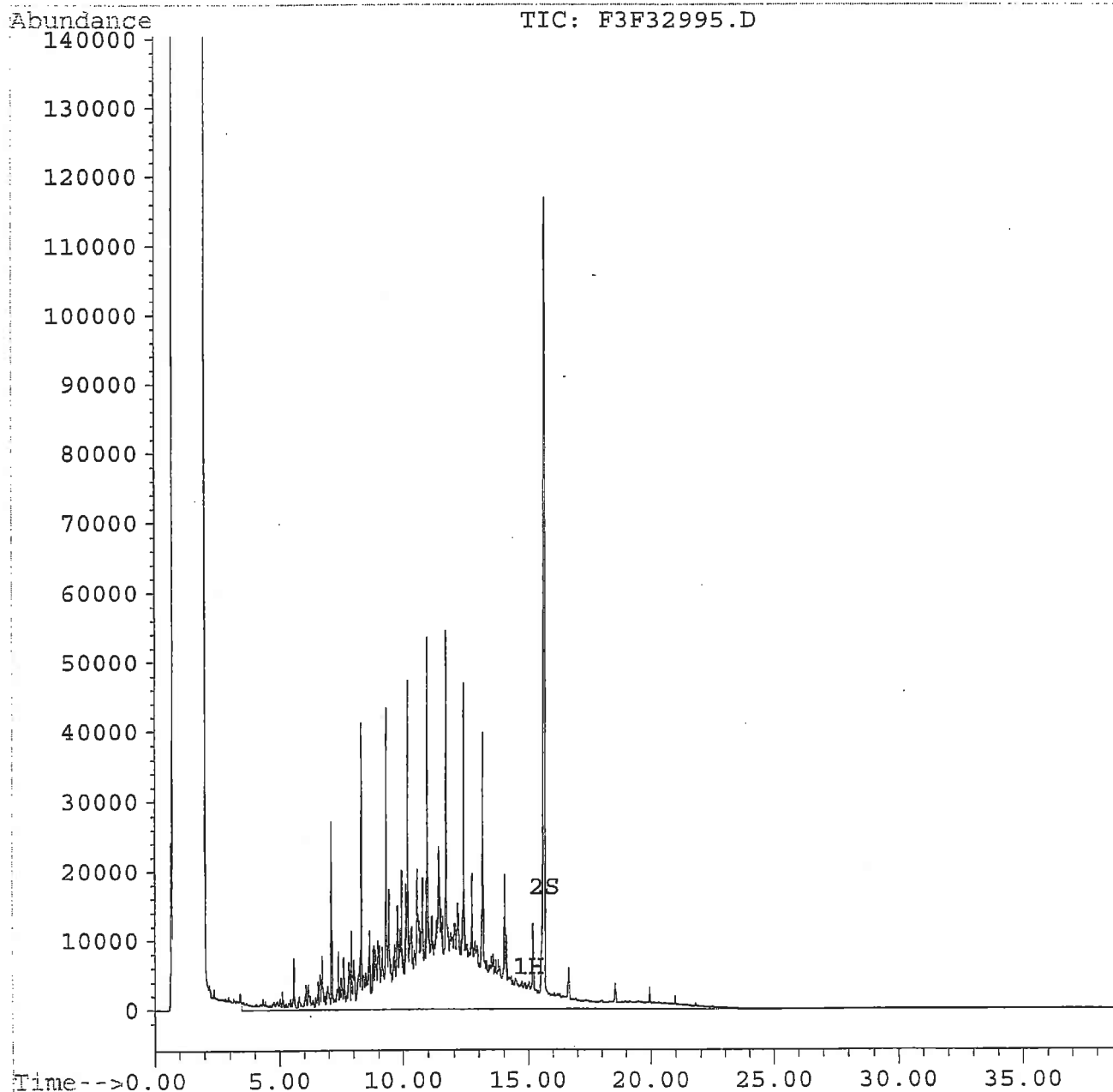
# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32995.D  
Acq On : 20 Nov 08 03:57 PM  
Sample : 1000ug/mL DRO CCV  
Misc : 200uL ST081119-3 + 800uL HEXANE  
Quant Time: Nov 21 11:19 19108

Vial: 1  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F33007.D  
 Acq On : 21 Nov 08 01:09 AM  
 Sample : 1000ug/mL DRO CCV2  
 Misc : 200uL ST081119-3 + 800uL HEXANE  
 Quant Time: Nov 21 12:11 19108

Vial: 13  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Thu Nov 20 12:03:46 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.66	486818	96.05 µg/ml
	Recovery	=	192.10%
Target Compounds			
1) H TEPH	15.00	4850273	1023.88 µg/ml

(f)=RT Delta > 1/2 Window

(m)=manual int.

F3F33007.D D111208.M

Fri Nov 21 12:11:27 2008

8/11/21/06

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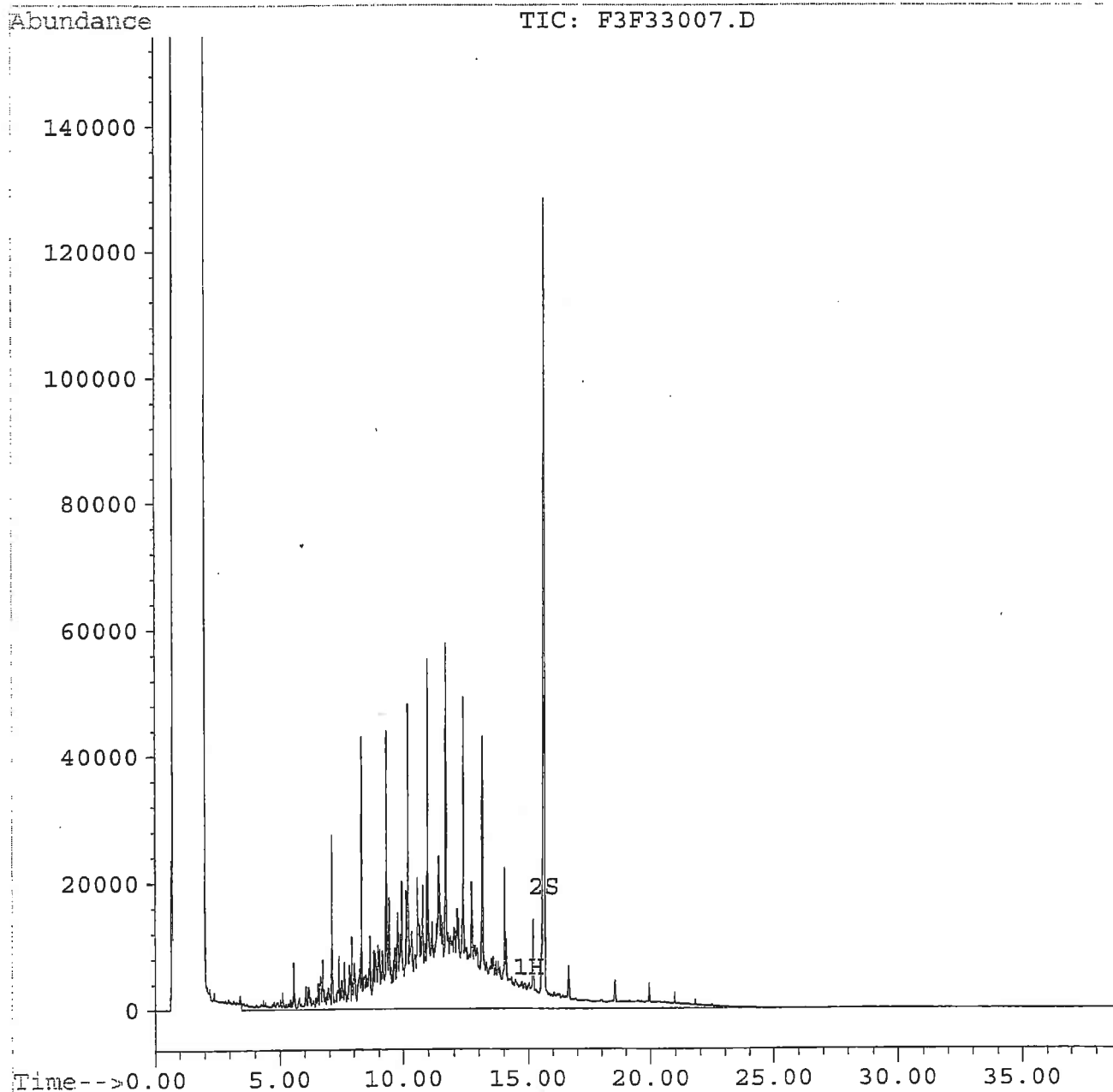
# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F33007.D  
Acq On : 21 Nov 08 01:09 AM  
Sample : 1000ug/mL DRO CCV2  
Misc : 200uL ST081119-3 + 800uL HEXANE  
Quant Time: Nov 21 12:11 19108

Vial: 13  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



## Sample Raw Data

# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32996.D  
 Acq On : 20 Nov 08 04:43 PM  
 Sample : EX081118-4MB  
 Misc : EX081118-4  
 Quant Time: Nov 21 12:14 19108

Vial: 2  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Thu Nov 20 12:03:46 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.63	221015	43.61 µg/ml
	Recovery	=	87.22%
Target Compounds			
1) H TEPH	15.00	139133	N.D. µg/ml

(f)=RT Delta > 1/2 Window

(m)=manual int.

F3F32996.D D111208.M

Fri Nov 21 12:14:12 2008

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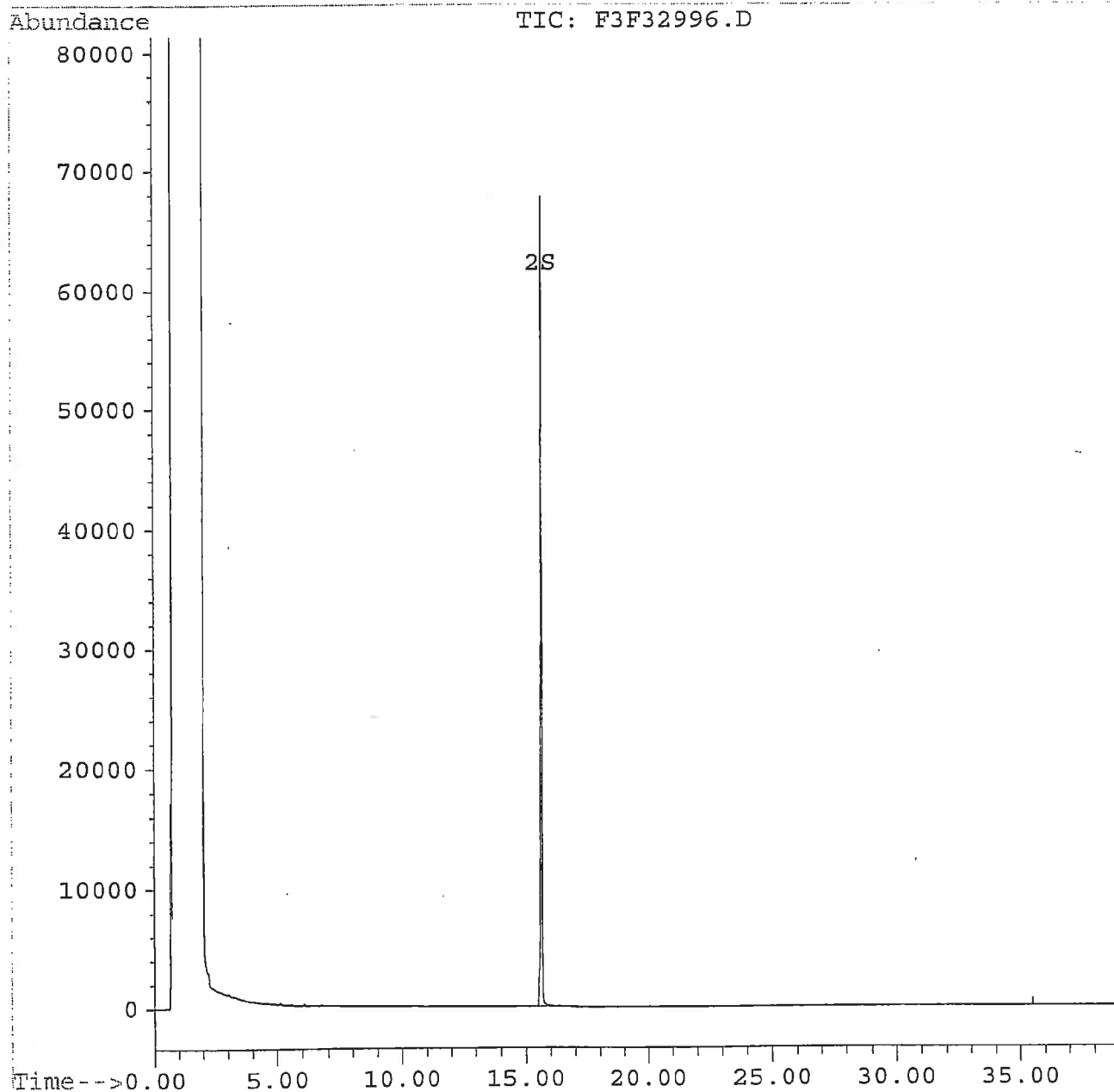
Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32996.D  
Acq On : 20 Nov 08 04:43 PM  
Sample : EX081118-4MB  
Misc : EX081118-4  
Quant Time: Nov 21 12:14 19108

Vial: 2  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
Signal Info : FID





# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32999.D Vial: 5  
 Acq On : 20 Nov 08 07:01 PM Operator: edb  
 Sample : 0811110-1 2X Inst : FUELS3  
 Misc : EX081118-4; 150uL Sample + 150uL Hexane Multiplr: 1.00  
 Quant Time: Nov 21 12:15 19108

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Thu Nov 20 12:03:46 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.62	107044	21.12 µg/ml <sup>(2x)</sup> 84%
	Recovery	=	42.24%
Target Compounds			
1) H TEPH	15.00	6341534	1348.14 µg/ml L
		TEPH	c7-c18

(f)=RT Delta > 1/2 Window

F3F32999.D D111208.M

Fri Nov 21 12:15:14 2008

(m)=manual int.

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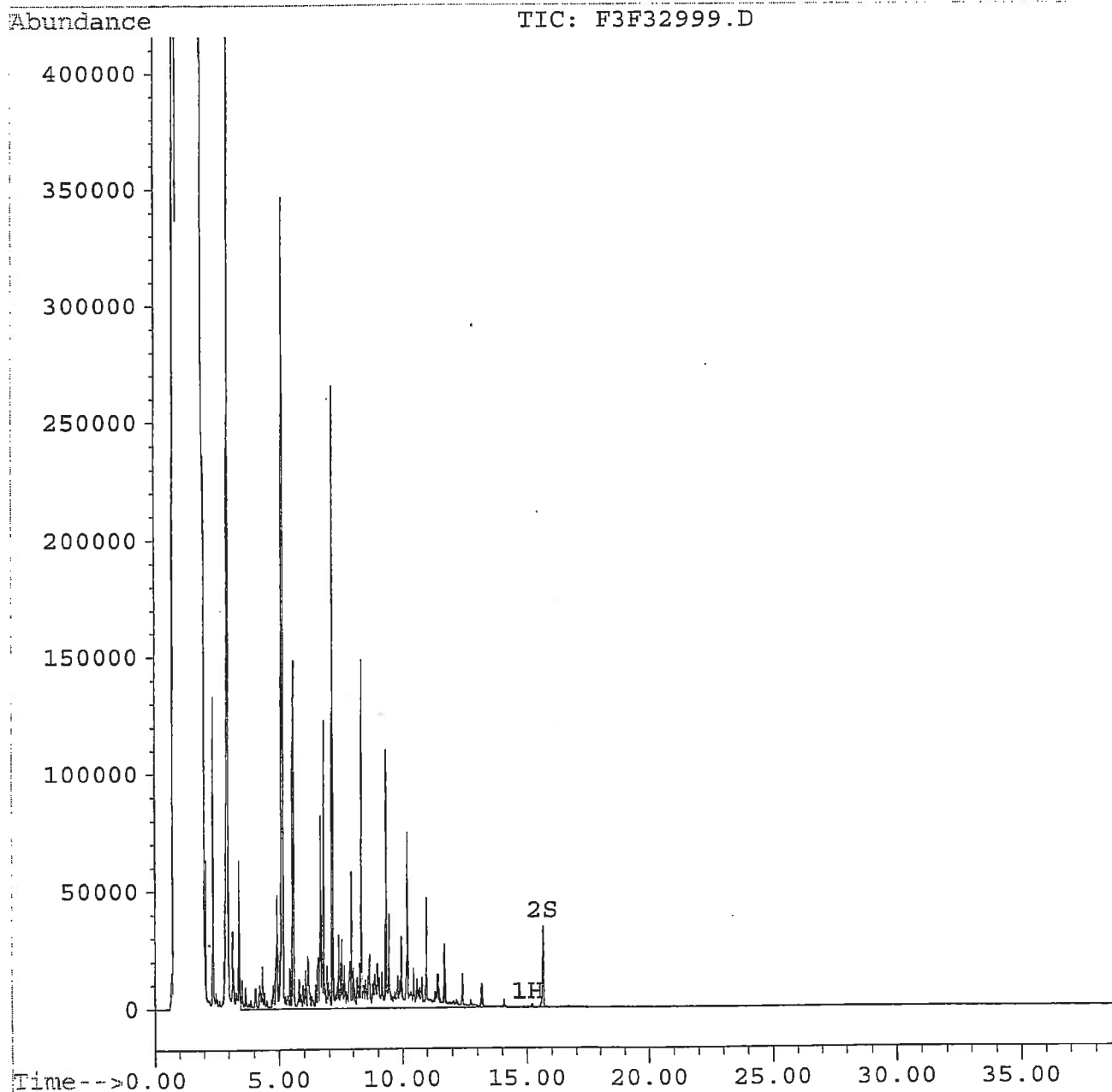
# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32999.D  
Acq On : 20 Nov 08 07:01 PM  
Sample : 0811110-1 2X  
Misc : EX081118-4; 150uL Sample + 150uL Hexane  
Quant Time: Nov 21 12:15 19108

Vial: 5  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F33000.D Vial: 6  
 Acq On : 20 Nov 08 07:47 PM Operator: edb  
 Sample : 0811110-2 2X Inst : FUELS3  
 Misc : EX081118-4; 150uL Sample + 150uL Hexane Multiplr: 1.00  
 Quant Time: Nov 21 12:15 19108

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Thu Nov 20 12:03:46 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.62	110241	(21.75) µg/ml (2%)
	Recovery	=	43.50% 872
Target Compounds			
1) H TEPH	15.00	9631923	(2063.60) µg/ml L
			TEPH C7-C26

(f)=RT Delta > 1/2 Window

F3F33000.D D111208.M

Fri Nov 21 12:15:18 2008

(m)=manual int.

SK  
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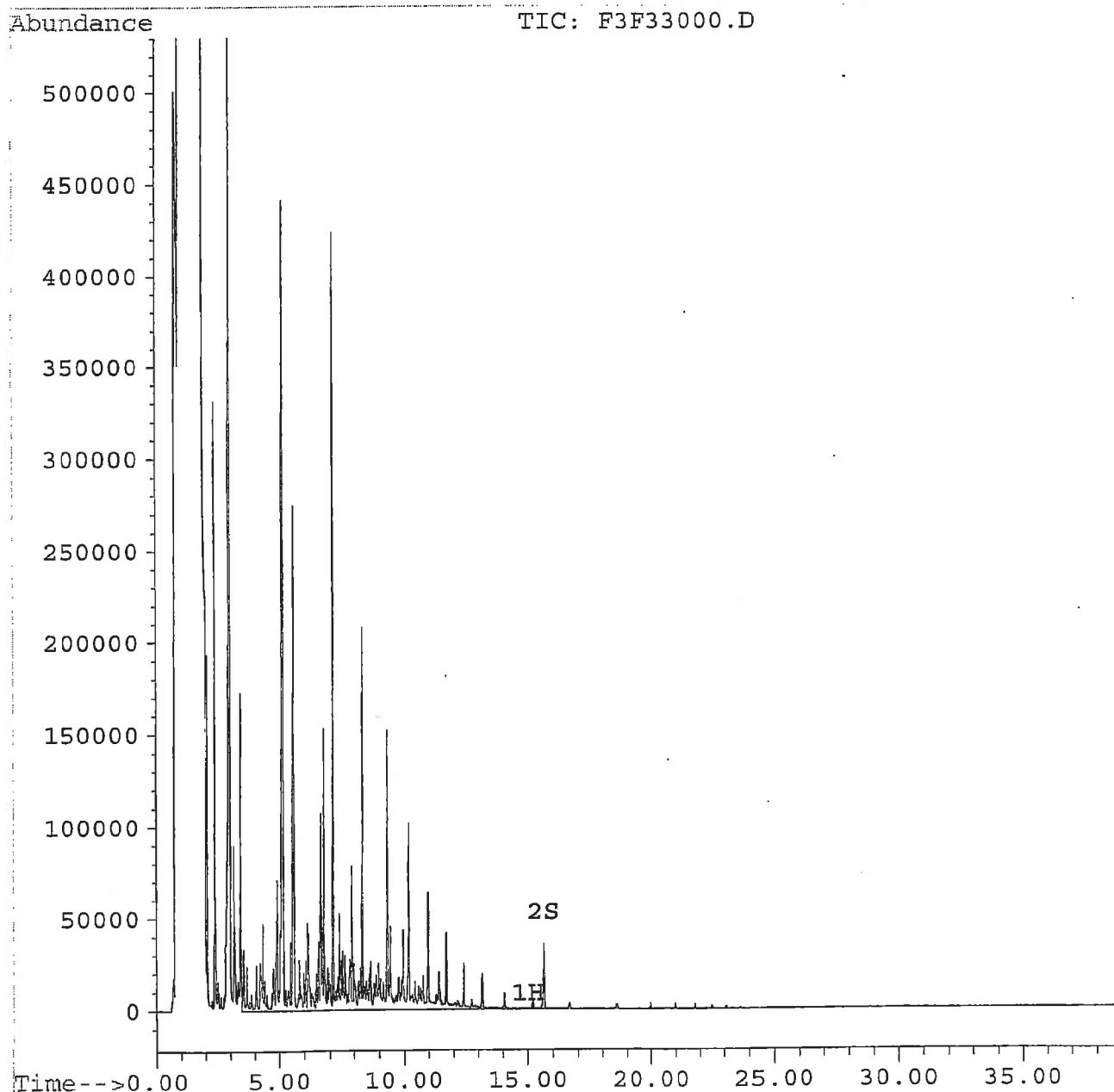
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# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F33000.D Vial: 6  
Acq On : 20 Nov 08 07:47 PM Operator: edb  
Sample : 0811110-2 2X Inst : FUELS3  
Misc : EX081118-4; 150uL Sample + 150uL Hexane Multiplr: 1.00  
Quant Time: Nov 21 12:15 19108

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



## Raw Data Quality Control Samples

# Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32997.D  
 Acq On : 20 Nov 08 05:29 PM  
 Sample : EX081118-4LCS  
 Misc : EX081118-4  
 Quant Time: Nov 21 12:15 19108

Vial: 3  
 Operator: edb  
 Inst : FUELS3  
 Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
 Title : 8015Bmod, CALuft  
 Last Update : Thu Nov 20 12:03:46 2008  
 Response via : Multiple Level Calibration

Volume Inj. : 1uL  
 Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
 Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.63	221793	<u>43.78</u> µg/ml
	Recovery	=	87.52%
Target Compounds			
1) H TEPH	15.00	1041145	<u>195.63</u> µg/ml 98%

(f)=RT Delta > 1/2 Window

F3F32997.D D111208.M

Fri Nov 21 12:15:06 2008

(m)=manual int.

edb  
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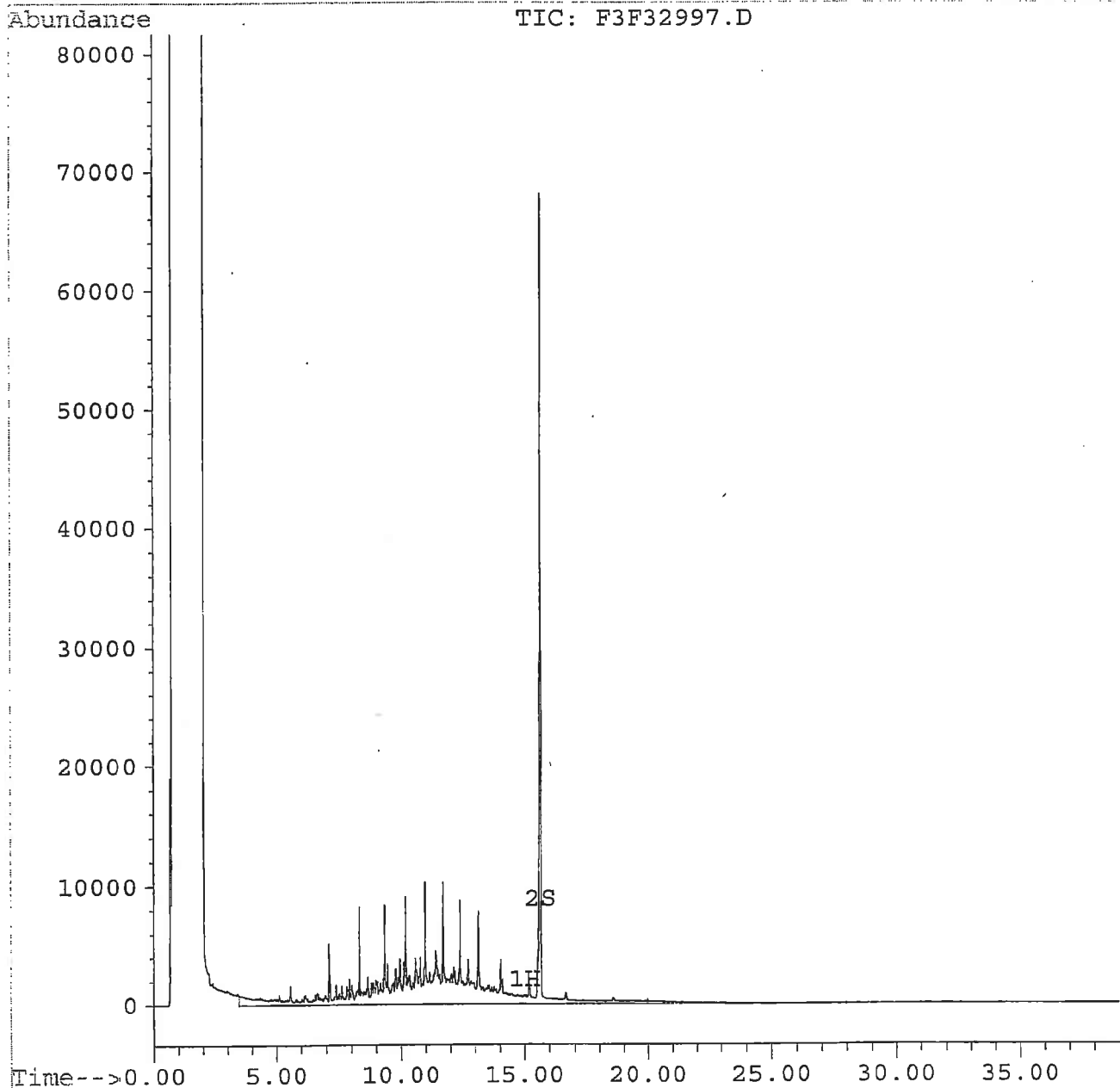
Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32997.D  
Acq On : 20 Nov 08 05:29 PM  
Sample : EX081118-4LCS  
Misc : EX081118-4  
Quant Time: Nov 21 12:15 19108

Vial: 3  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



## Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32998.D  
Acq On : 20 Nov 08 06:15 PM  
Sample : EX081118-4LCSD  
Misc : EX081118-4  
Quant Time: Nov 21 12:15 19108

Vial: 4  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod, CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5µm  
Signal Info : FID

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
2) S o-terphenyl	15.64	219861	<u>43.38</u> µg/ml
	Recovery	=	86.76%
Target Compounds			
1) H TEPH	15.00	1044022	<u>196.26</u> µg/ml 98%

(f)=RT Delta > 1/2 Window

F3F32998.D D111208.M

Fri Nov 21 12:15:10 2008

(m)=manual int.

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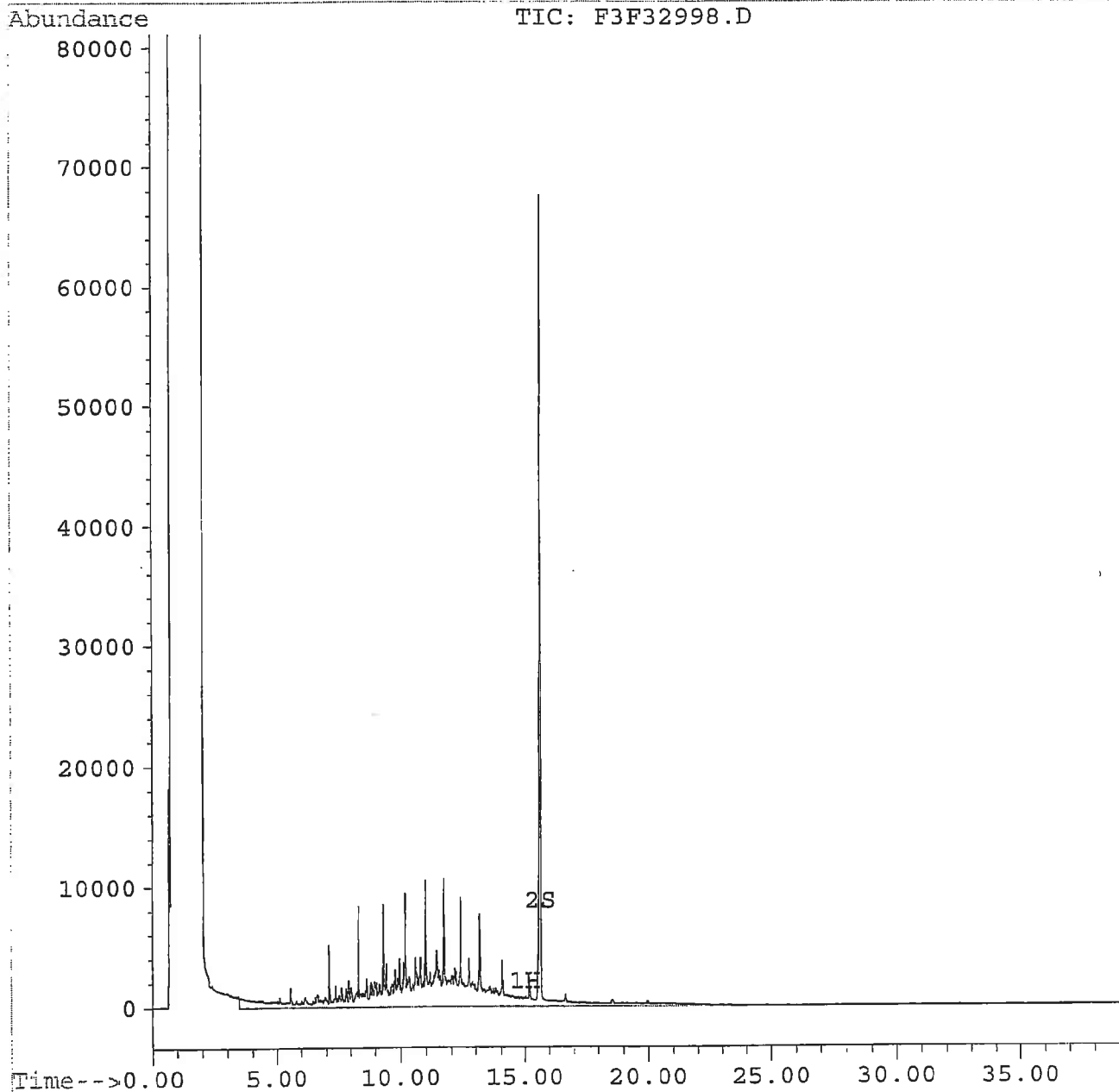
Quantitation Report

Data File : C:\HPCHEM\5\DATA\11202008\F3F32998.D  
Acq On : 20 Nov 08 06:15 PM  
Sample : EX081118-4LCSD  
Misc : EX081118-4  
Quant Time: Nov 21 12:15 19108

Vial: 4  
Operator: edb  
Inst : FUELS3  
Multiplr: 1.00

Method : C:\HPCHEM\5\METHODS\D111208.M  
Title : 8015Bmod; CALuft  
Last Update : Thu Nov 20 12:03:46 2008  
Response via : Multiple Level Calibration

Volume Inj. : 1uL  
Signal Phase : DB-5.625, 30m, 0.25mm 0.5 $\mu$ m  
Signal Info : FID



## Miscellaneous

369701

WO #s 081110 Matrix: AQ SO N/A Batch # EX08118-4 Surrogate Code ST080619-1 M Spike Code ST081107-9 Initials ES Balance ID N/A  
Extr. Start Date/Time 11/18/08/1420 Stop Date/Time 11/18/08/1430 Ext Code 8015M Cal Luft Method 8015M Cal Luft SOP 603 Rev 10 Reviewed by / Date ESK 11/18/08

[illegible]

Reagents: MeOH N/A DCM N/A 80% MeOH / 20% H<sub>2</sub>O Prep ID/Date N/A Hexane 080013 Na<sub>2</sub>SO<sub>4</sub> 083053

Form 602r10.doc (7/20/05)