TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

December 15, 2010

TestAmerica Project Number: G0K300434
PO/Contract: 2027.07

Ted Splitter
Tronox LLC / AIU Henderson, NV
PO Box 268859
Oklahoma City, OK 73126-8859

Dear Mr. Splitter,

This report contains the analytical results for the samples received under chain of custody by TestAmerica on November 30, 2010. These samples are associated with your Tronox Henderson Air Monitoring project.

The test results in this report meet all NELAC requirements for parameters that accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The case narrative is an integral part of this report.

If you have any questions, please feel free to call me at (916) 374-4383.

Sincerely,

DAVID R. ALLTUCKER
Project Manager

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TestAmerica West Sacramento Project Number G0K300434

Case Narrative

TestAmerica West Sacramento Project Number G0K300434

AIR, TO-13, Semivolatile Organics
Sample(s): 1, 4
The method blank associated with extraction batch 0334389 has d4-1,2-Dichlorobenzene pre-spike surrogate low outside stated control limits (58% vs. 60% lower limit). Re-injection confirms the low surrogate recovery. The field samples are in control for all surrogates. As air samples are unique, there is no possibility for re-extraction.

There were no other anomalies associated with this project.

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

ACCREDITED IN ACCORDANCE WITH nelac

TestAmerica Laboratories West Sacramento Certifications/Accreditations

Certifying State	Certificate #	Certifying State	Certificate #
Alaska	UST-055	New York*	11666
Arizona	AZ0708	Oregon*	CA 200005
Arkansas	88-0691	Pennsylvania	68-1272
California*	01119CA	South Carolina	87014
Colorado	NA	Texas	T104704399-08-TX
Connecticut	PH-0691	Utah*	QUAN1
Florida*	E87570	Virginia	00178
Georgia	960	Washington	C1281
Hawaii	NA	West Virginia	9930C, 334
Illinois	200060	Wisconsin	998204680
Kansas*	E-10375	NFESC	NA
Louisiana*	30612	USACE	NA
Michigan	9947	USDA Foreign Plant	37-82605
Nevada	CA44	USDA Foreign Soil	P330-09-00055
New Jersey*	CA005	US Fish & Wildlife	LE148388-0
New Mexico	NA	Guam	09-014r

*NELAP accredited. A more detailed parameter list is available upon request. Updated 3/25/2009

QC Parameter Definitions

QC Batch: The QC batch consists of a set of up to 20 field samples that behave similarly (i.e., same matrix) and are processed using the same procedures, reagents, and standards at the same time.

Method Blank: An analytical control consisting of all reagents, which may include internal standards and surrogates, and is carried through the entire analytical procedure. The method blank is used to define the level of laboratory background contamination.

Laboratory Control Sample and Laboratory Control Sample Duplicate (LCS/LCSD): An aliquot of blank matrix spiked with known amounts of representative target analytes. The LCS (and LCSD as required) is carried through the entire analytical process and is used to monitor the accuracy of the analytical process independent of potential matrix effects. If an LCSD is performed, it may also be used to evaluate the precision of the process.

Duplicate Sample (DU): Different aliquots of the same sample are analyzed to evaluate the precision of an analysis.

Surrogates: Organic compounds not expected to be detected in field samples, which behave similarly to target analytes. These are added to every sample within a batch at a known concentration to determine the efficiency of the sample preparation and analytical process.

Matrix Spike and Matrix Spike Duplicate (MS/MSD): An MS is an aliquot of a matrix fortified with known quantities of specific compounds and subjected to an entire analytical procedure in order to indicate the appropriateness of the method for a particular matrix. The percent recovery for the respective compound(s) is then calculated. The MSD is a second aliquot of the same matrix as the matrix spike, also spiked, in order to determine the precision of the method.

Isotope Dilution: For isotope dilution methods, isotopically labeled analogs (internal standards) of the native target analytes are spiked into the sample at time of extraction. These internal standards are used for quantitation, and monitor and correct for matrix effects. Since matrix effects on method performance can be judged by the recovery of these analogs, there is little added benefit of performing MS/MSD for these methods. MS/MSD are only performed for client or QAPP requirements.

Control Limits: The reported control limits are either based on laboratory historical data, method requirements, or project data quality objectives. The control limits represent the estimated uncertainty of the test results.

Sample Summary

TestAmerica West Sacramento Project Number G0K300434

WO#	Sample #	Client Sample ID	Sampling Date	Received Date
MAMLW	1	UW-11242010B	11/24/2010 02:42 PM	11/30/2010 09:10 AM
MAML1	2	UW-11242010B	11/24/2010 02:48 PM	11/30/2010 09:10 AM
MAML6	3	DW-11242010B	11/24/2010 03:04 PM	11/30/2010 09:10 AM
MAML8	4	DW-11242010B	11/24/2010 03:02 PM	11/30/2010 09:10 AM

Notes(s):

- The analytical results of the samples listed above are presented on the following pages.
- All calculations are performed before rounding to avoid round-off errors in calculated results.
- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity, pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight.

northgate
environmental management, inc.
300 Frank H. Ogawa Plaza, Ste 510
Oakland, CA 94612 (510) 839-0688

CHAIN-OF-CUSTODY / Analytical Request Document
The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed and accurate.

COC # 2027.07.0018

Total # of Samples: 4 Event Complete?

Regular / Rush / 5 day / Mark One

Required Ship to Lab

Lab Name	Test America Laboratories Inc
Address	880 Riverside Parkway
	West Sacramento, CA 95605
Lab PM	David Alltucker
Phone/FaH	(916) 373-5600
Lab PM email	David.Alltucker@testamericainc.com
Applicable Lab Quote #:	

Required Project Information:

Site ID #:102	TRONOX LLC. HENDERSON
Project #	2027.07
Site Address	560 W Lake Mead Pkwy
City	Henderson
State, Zip	NV, 89015
Site PM Name	Ted Splitter
Phone/FaH:	(610) 435-4609
Site PM Email:	Ted Splitter@ngem.com

Required Invoice Information:

Send Invoice to	Susan Crowley Tronox LLC.
Address	PO Box 55
City/State	Henderson, NV 89009
Phone #:	(949) 260-9293
PO #	
Send EDD to	Frank.Hagar@ngem.com
CC Hardcopy report to	PDF Electronic Version Only – FTP Upload
CC Hardcopy report to	See Additional Comments Below

ITEM #	SAMPLE ID Samples IDs MUST BE UNIQUE	SAMPLE LOCATION	MATRIX CODE	G=GRAB C=COMP	SAMPLE TYPE	SAMPLE DATE	SAMPLE TIME	#OF CONTAINERS	Comments/Lab Sample I.D. Volume (m^3)	Filtered	Preservative	Analysis: TO-8A/Dioxins, Furans	TO-13A/8270C/HCB	T&P	6020/As/Mu/CPMS
	UW-11242010B		AA			11/24/2010	2.42 PM	1	474.16				X		
	UW-11242010B		AA			11/24/2010	2:48 PM	1	698.9					X	X
	DW-11242010B		AA			11/24/2010	3:04 PM	1	667.57					X	X
	DW-11242010B		AA			11/24/2010	3:02 PM	1	450.17				X		

Additional Comments/Special Instructions:
3-5 DAY TURN AROUND

RELEASED BY / AFFILIATION	DATE	TIME	RECEIVED BY / AFFILIATION	DATE	TIME
Ronda S. Bailey	11/29/10	1500		11/30/10	0920

Sample Receipt Conditions

Temp in °C	Samples on ice?	Sample intact?	Trip Blank?
4.2	Y/N	Y/N	Y/N
	Y/N	Y/N	Y/N
	Y/N	Y/N	Y/N
	Y/N	Y/N	Y/N

SHIPPING INFO

SAMPLER NAME AND SIGNATURE

PRINT Name of SAMPLER: Ronda Bailey

SIGNATURE of SAMPLER: DATE Signed: 11/29/10 1500 Time

TestAmerica

411111 THE LEADER IN ENVIRONMENTAL TESTING

LOT RECEIPT CHECKLIST
TestAmerica West Sacramento

CLIENT Northgate PM DA LOG # 68332

LOT# (QUANTIMS ID) G0K300434 QUOTE# 84087 LOCATION W14D AC

Checked (✓)

DATE RECEIVED 11/30/10 TIME RECEIVED 0910 ☑

DELIVERED BY ☒ FEDEX ☐ ON TRAC ☐ CLIENT

☐ GOLDENSTATE ☐ UPS ☐ GO-GETTERS ☐ OTHER

☐ TAL COURIER ☐ TAL SF ☐ VALLEY LOGISTICS ☑

CUSTODY SEAL STATUS ☒ INTACT ☐ BROKEN ☐ N/A ☑

CUSTODY SEAL #(S) NA

SHIPPPING CONTAINER(S) ☐ TAL ☒ CLIENT ☐ N/A ☑

COC #(S) 2027.07.0018 ☑

TEMPERATURE BLANK Observed: NA Corrected:

SAMPLE TEMPERATURE - (TEMPERATURES ARE IN °C)

Observed: 4 Average 4 Corrected Average 4

LABORATORY THERMOMETER ID:

IR UNIT: #4 ☒ #5 ☐ ☐ OTHER ☑

CV 11/30/10
Initials Date

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pH MEASURED ☐ YES ☐ ANOMALY ☒ N/A ☑

LABELED BY ☑

LABELS CHECKED BY ☑

PEER REVIEW ☒ NA

SHORT HOLD TEST NOTIFICATION

SAMPLE RECEIVING ☑

WETCHEM ☒ N/A ☑

VOA-ENCORES ☒ N/A ☑

☐ METALS NOTIFIED OF FILTER/PRESERVE VIA VERBAL & EMAIL ☒ N/A ☑

☒ COMPLETE SHIPMENT RECEIVED IN GOOD CONDITION WITH APPROPRIATE TEMPERATURES, CONTAINERS, PRESERVATIVES ☐ N/A ☑

☐ CLOUSEAU ☐ TEMPERATURE EXCEEDED (2 °C – 6 °C)*1 ☒ N/A

☐ WET ICE ☐ BLUE ICE ☐ GEL PACK ☐ NO COOLING AGENTS USED ☐ PM NOTIFIED

CV 11/30/10
Initials Date

Notes

*1 Acceptable temperature range for State of Wisconsin samples is ≤4°C.

THE LEADER IN ENVIRONMENTAL TESTING

Lot ID: G0K300434

	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20
VOA*																				
VOAh*																				
AGB																				
AGBs																				
250AGB																				
250AGBs																				
250AGBn																				
500AGB																				
____AGJ																				
500AGJ																				
250AGJ																				
125AGJ																				
____CGJ																				
500CGJ																				
250CGJ																				
125CGJ																				
PJ																				
PJn																				
500PJ																				
500PJn																				
500PJna																				
500PJzn/na																				
250PJ																				
250PJn																				
250PJna																				
250PJzn/na																				
Acetate Tube																				
_____"CT																				
Encore																				
Folder/filter		1	1																	
PUF	1			1																
Petri/Filter																				
XAD Trap																				
Ziploc																				
	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20

h = hydrochloric acid **s** = sulfuric acid **na** = sodium hydroxide **n** = nitric acid **zn** = zinc acetate

Number of VOAs with air bubbles present / total number of VOA's

AIR, TO-13, Semivolatile Organics

Northgate Environmental Management, Inc.

Sample ID: UW-11242010B

Trace Level Compounds

Lot - Sample #....:	G0K300434 - 001	**Work Order #....:**	MAMLW1AA	**Matrix....:**	AA
Date Sampled....:	11/24/10	**Date Received....:**	11/30/10	**Dilution Factor....:**	1
Prep Date....:	11/30/10	**Analysis Date....:**	12/04/10	**Volume....:**	474.16
Prep Batch # ....:	0334389	**Instrument ID....:**	5MH	**Method....:**	EPA-2 TO-13
Initial Wgt/Vol....:	1 Sample	**Analyst ID....:**	Mark Onishi		

PARAMETER	RESULT	REPORTING LIMIT	DETECTION LIMIT	UNITS
Hexachlorobenzene	ND	0.021	0.0027	ug/m3

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
1,2-Dichlorobenzene-d4	71	60 - 120
2-Fluorobiphenyl	85	58 - 105
2-Fluorophenol	73	41 - 105
Nitrobenzene-d5	81	46 - 118
Phenol-d5	85	43 - 122
Terphenyl-d14	92	69 - 110
2,4,6-Tribromophenol	106	61 - 118

QUALIFIERS

Northgate Environmental Management, Inc.

Sample ID: DW-11242010B

Trace Level Compounds

Lot - Sample #....:	G0K300434 - 004	**Work Order #....:**	MAML81AA	**Matrix....:**	AA
Date Sampled....:	11/24/10	**Date Received....:**	11/30/10	**Dilution Factor....:**	1
Prep Date....:	11/30/10	**Analysis Date....:**	12/04/10	**Volume....:**	450.17
Prep Batch # ....:	0334389	**Instrument ID....:**	5MH	**Method....:**	EPA-2 TO-13
Initial Wgt/Vol....:	1 Sample	**Analyst ID....:**	Mark Onishi		

PARAMETER	RESULT	REPORTING LIMIT	DETECTION LIMIT	UNITS
Hexachlorobenzene	ND	0.022	0.0029	ug/m3

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
1,2-Dichlorobenzene-d4	68	60 - 120
2-Fluorobiphenyl	78	58 - 105
2-Fluorophenol	64	41 - 105
Nitrobenzene-d5	73	46 - 118
Phenol-d5	80	43 - 122
Terphenyl-d14	93	69 - 110
2,4,6-Tribromophenol	106	61 - 118

QUALIFIERS

QC DATA ASSOCIATION SUMMARY

G0K300434

Sample Preparation and Analysis Control Numbers

SAMPLE#	MATRIX	ANALYTICAL METHOD	LEACH BATCH #	PREP BATCH #	MS RUN#
001	AA	EPA-2 TO-13		0334389	
002	AA	SW846 6020		0340010	
003	AA	SW846 6020		0340010	
004	AA	EPA-2 TO-13		0334389	

Method Blank Report

Trace Level Compounds

Lot - Sample #....:	G0K300000 - 389B	**Work Order #....:**	MANG51AA	**Matrix....:**	AIR
Date Sampled....:	11/24/10	**Date Received....:**	11/30/10	**Dilution Factor....:**	1
Prep Date....:	11/30/10	**Analysis Date....:**	12/04/10	**Volume....:**	0
Prep Batch # ....:	0334389	**Instrument ID....:**	5MH	**Method....:**	EPA-2 TO-13
Initial Wgt/Vol....:	1 Sample	**Analyst ID....:**	Mark Onishi		

PARAMETER	RESULT	REPORTING LIMIT	DETECTION LIMIT	UNITS
Hexachlorobenzene	ND	10.0	1.3	ug

SURROGATE	PERCENT RECOVERY		RECOVERY LIMITS
1,2-Dichlorobenzene-d4	58	*	60 - 120
2-Fluorobiphenyl	73		58 - 105
2-Fluorophenol	63		41 - 105
Nitrobenzene-d5	69		46 - 118
Phenol-d5	70		43 - 122
Terphenyl-d14	93		69 - 110
2,4,6-Tribromophenol	101		61 - 118

QUALIFIERS

* Surrogate recovery is outside stated control limits.

LABORATORY CONTROL SAMPLE DATA REPORT

Trace Level Compounds

Client Lot # ...:	G0K300434	**Work Order # ...:**	MANG51AC-LCS MANG51AD-LCSD	**Matrix .....:**	AIR
LCS Lot-Sample# :	G0K300000 - 389				
Prep Date .....:	11/30/10	**Analysis Date ..:**	12/04/10		
Prep Batch # ...:	0334389				
Dilution Factor :	1				
Analyst ID.....:	Mark Onishi	**Instrument ID..:**	5MH	**Method.....:**	EPA-2 TO-13
Initial Wgt/Vol:	1 Sample				

PARAMETER	SPIKE AMOUNT	MEASURED AMOUNT	UNITS	PERCENT RECOVERY	RECOVERY LIMITS	RPD	RPD LIMITS
Hexachlorobenzene	**100**	**97.9**	**ug**	**98**	**(70 - 110)**		
	100	**97.0**	**ug**	**97**	**(70 - 110)**	**0.86**	**(0 - 30)**

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
2-Fluorobiphenyl	90	(58 - 105)
	91	(58 - 105)
2-Fluorophenol	77	(41 - 105)
	74	(41 - 105)
Nitrobenzene-d5	85	(46 - 118)
	84	(46 - 118)
Phenol-d5	86	(43 - 122)
	79	(43 - 122)
Terphenyl-d14	91	(69 - 110)
	90	(69 - 110)
2,4,6-Tribromophenol	109	(61 - 118)
	112	(61 - 118)

Notes:

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

AIR, Metals by ICPMS (As and Mn)

Northgate Environmental Management, Inc.

Sample ID: UW-11242010B

Trace Level Compounds

Lot - Sample #....:	G0K300434 - 002	**Work Order #....:**	MAML11AC	**Matrix....:**	AA
Date Sampled....:	11/24/10	**Date Received....:**	11/30/10	**Dilution Factor....:**	1
Prep Date....:	12/06/10	**Analysis Date....:**	12/07/10	**Volume....:**	698.9
Prep Batch # ....:	0340010	**Instrument ID....:**	M02	**Method....:**	SW846 6020
Initial Wgt/Vol....:	0.08333 L	**Analyst ID....:**	Sabine Hargrave		

PARAMETER	RESULT		REPORTING LIMIT	DETECTION LIMIT	UNITS
Arsenic	0.0016	B J	0.0034	0.00070	ug/m3
Manganese	3.02	J	0.00172	0.000243	ug/m3

QUALIFIERS

B Estimated result. Result is less than RL and greater than or equal to the IDL.

J Estimated Result.

Northgate Environmental Management, Inc.

Sample ID: DW-11242010B

Trace Level Compounds

Lot - Sample #....:	G0K300434 - 003	**Work Order #....:**	MAML61AC	**Matrix....:**	AA
Date Sampled....:	11/24/10	**Date Received....:**	11/30/10	**Dilution Factor....:**	1
Prep Date....:	12/06/10	**Analysis Date....:**	12/07/10	**Volume....:**	667.57
Prep Batch # ....:	0340010	**Instrument ID....:**	M02	**Method....:**	SW846 6020
Initial Wgt/Vol....:	0.08333 L	**Analyst ID....:**	Sabine Hargrave		

PARAMETER	RESULT		REPORTING LIMIT	DETECTION LIMIT	UNITS
Arsenic	0.0021	B J	0.0036	0.00073	ug/m3
Manganese	0.199	J	0.00180	0.000255	ug/m3

QUALIFIERS

B Estimated result Result is less than RL and greater than or equal to the IDL

J Estimated Result.

QC DATA ASSOCIATION SUMMARY

G0K300434

Sample Preparation and Analysis Control Numbers

SAMPLE#	MATRIX	ANALYTICAL METHOD	LEACH BATCH #	PREP BATCH #	MS RUN#
002	AA	SW846 6020		0340010	
003	AA	SW846 6020		0340010	

Method Blank Report

Trace Level Compounds

Lot - Sample #....:	G0L060000 - 010B	**Work Order #....:**	MAWLK1AA	**Matrix.....:**	AIR
Date Sampled....:	11/24/10	**Date Received....:**	11/30/10	**Dilution Factor....:**	1
Prep Date....:	12/06/10	**Analysis Date....:**	12/07/10	**Volume....:**	0
Prep Batch # ....:	0340010	**Instrument ID....:**	M02	**Method....:**	SW846 6020
Initial Wgt/Vol....:	0.08333 L	**Analyst ID....:**	Sabine Hargrave		

PARAMETER	RESULT		REPORTING LIMIT	DETECTION LIMIT	UNITS
Arsenic	**0.77**	**B**	**2.4**	**0.49**	**ug**
Manganese	**0.21**	**B**	**1.2**	**0.17**	**ug**

QUALIFIERS

B Estimated result Result is less than RL and greater than or equal to the IDL.

LABORATORY CONTROL SAMPLE DATA REPORT

Trace Level Compounds

Client Lot # ...:	G0K300434	**Work Order # ...:**	MAWLK1AD-LCS MAWLK1AE-LCSD	**Matrix .....:**	AIR
LCS Lot-Sample# :	G0L060000 - 010				
Prep Date .....:	12/06/10	**Analysis Date ..:**	12/07/10		
Prep Batch # ...:	0340010				
Dilution Factor :	1				
Analyst ID.....:	Sabine Hargrave	**Instrument ID..:**	M02	**Method.....:**	SW846 6020
Initial Wgt/Vol:	0.08333 L				

PARAMETER	SPIKE AMOUNT	MEASURED AMOUNT	UNITS	PERCENT RECOVERY	RECOVERY LIMITS	RPD	RPD LIMITS
Arsenic	**240**	**221**	**ug**	**92**	**(86 - 110)**		
	240	**224**	**ug**	**93**	**(86 - 110)**	**1.6**	**(0 - 15)**
Manganese	**240**	**223**	**ug**	**93**	**(88 - 110)**		
	240	**227**	**ug**	**95**	**(88 - 110)**	**1.9**	**(0 - 15)**

Notes:

Calculations are performed before rounding to avoid round-off errors in calculated results.

Bold print denotes control parameters

AIR, TSP-
Total Suspended
Particulates

Northgate Environmental Management, Inc.

Sample ID: UW-11242010B

Trace Level Compounds

Lot - Sample #....:	G0K300434 - 002	**Work Order #....:**	MAML11AA	**Matrix....:**	AA
Date Sampled....:	11/24/10	**Date Received....:**	11/30/10	**Dilution Factor....:**	1
Prep Date....:	11/30/10	**Analysis Date....:**	12/07/10	**Volume....:**	698.9
Prep Batch # ....:	0341292	**Instrument ID....:**	QA-045	**Method....:**	CFR50B APDX B
Initial Wgt/Vol....:	0	**Analyst ID....:**	Thep Phomsopha		

PARAMETER	RESULT	REPORTING LIMIT	DETECTION LIMIT	UNITS
Total Suspended Particulates	0.0000306	0.000000715	--	g/m3

QUALIFIERS

Northgate Environmental Management, Inc.

Sample ID: DW-11242010B

Trace Level Compounds

Lot - Sample #....:	G0K300434 - 003	**Work Order #....:**	MAML61AA	**Matrix....:**	AA
Date Sampled....:	11/24/10	**Date Received....:**	11/30/10	**Dilution Factor....:**	1
Prep Date....:	11/30/10	**Analysis Date....:**	12/07/10	**Volume....:**	667.57
Prep Batch # ....:	0341292	**Instrument ID....:**	QA-045	**Method....:**	CFR50B APDX B
Initial Wgt/Vol....:		**Analyst ID....:**	Thep Phomsopha		

PARAMETER	RESULT	REPORTING LIMIT	DETECTION LIMIT	UNITS
Total Suspended Particulates	**0.000144**	**0.000000749**	--	**g/m3**

QUALIFIERS

QC DATA ASSOCIATION SUMMARY

G0K300434

Sample Preparation and Analysis Control Numbers

SAMPLE#	MATRIX	ANALYTICAL METHOD	LEACH BATCH #	PREP BATCH #	MS RUN#
002	AA	CFR50B APDX B		0341292	
003	AA	CFR50B APDX B		0341292	

AIR, TO-13, Semivolatile Organics

Raw Data Package

Run/Batch Data

Includes (as applicable):

runlogs

continuing calibration standards

interference/performance check standards

continuing calibration blanks

method blanks

lcs

ms/sd

sample raw data

ms tune data

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

West Sacramento

8270C CCV Checklist

Instrument: SV5 _____

ICAL Date: 10/02/10 _____ DFTPP ID: DFT1204

Initiator/Date: KT-12/06/10 _____ Standard ID: HSL1204

Reviewer/Date: [signature] 12/6/10 NCM #: _____

I: 8270C Criteria

	Initiated	Reviewed
Log Book page included.	☒	☑
CCV compared to correct ICAL.	☒	☑
Tune documentation is present and meets criteria.	☒	☑
Manual re-integrations are checked, initialed and hardcopies included.	☒	☑
Retention time correct for Isomers and all other analytes.	☒	☑
CCV Internal Standards are within 50-200% of ICAL mid-point.	☒	☑
Samples analyzed within 12 hours of Tune time.	☒	☑
Tailing and degradation criteria are met.	☒	☑
Spot check manual integrations in Target. Analyte checked:_____	NA	☑
Non-CCC ≤ 50% D	☒	☑

II: 8270C SPCC Check SPCC RRFs must be greater than 0.050

	Initiated	Reviewed		Initiated	Reviewed
N-nitroso-di-n-propylamine	☒	☑	2,4-Dinitrophenol	☒	☑
Hexachlorocyclopentadiene	☒	☑	4-Nitrophenol	☒	☑

III: 8270C CCC Check CCC must be ≤ 20%D (If CCC are not targets, all analytes must be <20%D.)

	Initiated	Reviewed		Initiated	Reviewed
Phenol	☒	☑	Acenaphthene	☒	☑
1,4-Dichlorobenzene	☒	☑	N-nitrosodiphenylamine	☒	☑
2-Nitrophenol	☒	☑	Pentachlorophenol	☒	☑
2,4-Dinitrophenol	☒	☑	Flouranthene	☒	☑
Hexachlorobutadiene	☒	☑	Di-n-octyl phthalate	☒	☑
4-Chloro-3-methylphenol	☒	☑	Benzo(a)pyrene	☒	☑
2,4,6-Trichlorophenol	☒	☑			

IV: AFCEE 3.1 and 4.0 QAPP Criteria

	Initiated	Reviewed
All analytes in CCV +/- 20%D compared to ICAL. See Note	☒	See ☐ Note
CCV and Sample Internal Standards are within 50-200% of ICAL mid-point.	☒	☑
Are the compounds which required manual integrations documented in the MI spreadsheet?	☒	☑

Benzo(B) Fluoranthene = +22% high and ND ok.

KT 12/6/10

V: DOD QSM V3 Criteria

	Initiated	Reviewed
For 8270, CCCs must be ≤ 20% D.	☒	☑
RRFs for SPCCs must meet minimum response factor criteria	☒	☑
CCV and sample Internal Standards are within 50-200% of ICAL mid-point.	☒	☑
SIM: All analytes must be ≤ 20%	☐NA	☑NA
Are the compounds which required manual integrations documented in the MI spreadsheet?	☒	☑

TestAmerica West Sacramento Page# _____

GC/MS INSTRUMENT LOG
SEMI-VOLATILES

Method Key (MTH Column)
QL = EPA 8270C (WS-MS-0005)
JZ = EPA TO-13A (WS-MS-0005)
VX = EPA 8270C-SIM (mod) CWM (WS-MS-0003)
QI = EPA 8270C-SIM (WS-MS-0008)
FX = PAH-SIM Isotope Dilution (WS-MS-0006)
F9 = EPA 8270C-SIM (mod) 1,4-Dioxane (WS-MS-0011)

Inst ID : sv5.i
Batch ID : 120410.B
ICAL Date: See Calib Report
See raw data for standard IDs

Date	Time	USER	Sample ID	File ID	Vol or Wt	Extract Vol	Diln	MTH	Comments
04-DEC-2010	09:53	KT	PRIMER	QC001A.D	NA	NA	NA		
04-DEC-2010	10:15	KT	DFTPP 50ug/ml	DFT1204.D	NA	NA	NA		
04-DEC-2010	10:36	KT	HSL_050 ug/ml CS-4	HSL1204.D	NA	NA	NA		
04-DEC-2010	11:00	KT	MALNT1AA G0K290000-259B	S120401.D	1000 Sa	1 mL	1	JZ	
04-DEC-2010	11:25	KT	MALNT1AC G0K290000-259C	S120402.D	1000 Sa	1 mL	1	JZ	
04-DEC-2010	11:49	KT	MALNT1AD G0K290000-259L	S120403.D	1000 Sa	1 mL	1	JZ	High % R
04-DEC-2010	12:14	KT	L99F61DL G0K180595-1S	S120404.D	30.06 g	1 mL	1	QL	
04-DEC-2010	12:38	KT	L99F61DM G0K180595-1D	S120405.D	29.95 g	1 mL	1	QL	
04-DEC-2010	13:03	KT	MAK001AA G0K270427-1	S120406.D	1000 Sa	1 mL	1	JZ	
04-DEC-2010	13:27	KT	MAK011AA G0K270427-2	S120407.D	1000 Sa	1 mL	1	JZ	
04-DEC-2010	13:52	KT	MAK021AA G0K270427-3	S120408.D	1000 Sa	1 mL	1	JZ	Low Surr.
04-DEC-2010	14:16	KT	MAK031AA G0K270427-4	S120409.D	1000 Sa	1 mL	1	JZ	low surr.
04-DEC-2010	14:41	KT	MANG51AA G0K30000-389B	S120410.D	1000 Sa	1 mL	1	JZ	low surr.
04-DEC-2010	15:05	KT	MANG51AC G0K30000-389C	S120411.D	1000 Sa	1 mL	1	JZ	
04-DEC-2010	15:30	KT	MANG51AD G0K30000-389L	S120412.D	1000 Sa	1 mL	1	JZ	
04-DEC-2010	15:54	KT	MAML W1AA G0K300434-1	S120413.D	1000 Sa	1 mL	1	JZ	
04-DEC-2010	16:19	KT	MAML81AA G0K300434-4	S120414.D	1000 Sa	1 mL	1	JZ	
04-DEC-2010	16:43	KT	L99G21AE G0K180595-8	S120415.D	29.99 g	1 mL	1	QL	
04-DEC-2010	17:08	KT	L99G31AE G0K180595-9	S120416.D	30.07 g	1 mL	1	QL	
04-DEC-2010	17:32	KT	L99G41AE G0K180595-10	S120417.D	30.08 g	1 mL	1	QL	
04-DEC-2010	17:57	KT	L99G51AE G0K180595-11	S120418.D	30.05 g	1 mL	1	QL	
04-DEC-2010	18:21	KT	L99G61AE G0K180595-12	S120419.D	29.92 g	1 mL	1	QL	
04-DEC-2010	18:46	KT	L99G91AE G0K180595-13	S120420.D	30.07 g	1 mL	1	QL	
04-DEC-2010	19:10	KT	L99HA1AE G0K180595-14	S120421.D	30.03 g	1 mL	1	QL	
04-DEC-2010	19:35	KT	L99HD1AE G0K180595-15	S120422.D	29.97 g	1 mL	1	QL	
04-DEC-2010	19:59	KT	L99HF1AE G0K180595-16	S120423.D	30.01 g	1 mL	1	QL	
04-DEC-2010	20:24	KT	L99HH1AE G0K180595-17	S120424.D	30.07 g	1 mL	1	QL	

TestAmerica West Sacramento

CONTINUING CALIBRATION COMPOUNDS

Instrument ID: sv5.i Injection Date: 04-DEC-2010 10:36
Lab File ID: HSL1204.D Init. Cal. Date(s): 17-AUG-2010 02-OCT-2010
Analysis Type: Init. Cal. Times: 17:32 15:00
Lab Sample ID: HSL 050 ug/ml CS-4 Quant Type: ISTD
Method: \\SV5\C\chem\sv5.i\120410.B\8270f.m

COMPOUND	RRF / AMOUNT	RF50	CCAL RRF50	MIN RRF	%D / %DRIFT	MAX %D / %DRIFT	CURVE TYPE
$ 7 2-Fluorophenol	1.40992	1.46404	1.46404	0.010	3.83874	50.00000	Averaged
$ 8 Phenol-d5	1.77296	1.83375	1.83375	0.010	3.42888	50.00000	Averaged
$ 9 2-Chlorophenol-d4	1.55698	1.61867	1.61867	0.010	3.96181	50.00000	Averaged
$ 10 1,2-Dichlorobenzene-d4	0.98513	1.01892	1.01892	0.010	3.43026	50.00000	Averaged
$ 11 Nitrobenzene-d5	0.33879	0.34944	0.34944	0.010	3.14130	50.00000	Averaged
$ 12 2-Fluorobiphenyl	1.28852	1.29558	1.29558	0.010	0.54766	50.00000	Averaged
$ 13 2,4,6-Tribromophenol	0.17381	0.18767	0.18767	0.010	7.97121	50.00000	Averaged
$ 14 Terphenyl-d14	0.78789	0.81013	0.81013	0.010	2.82244	50.00000	Averaged
15 N-Nitrosodimethylamine	0.92154	0.90858	0.90858	0.010	-1.40620	50.00000	Averaged
16 Pyridine	1.54111	1.50575	1.50575	0.010	-2.29461	50.00000	Averaged
23 Aniline	2.25673	2.27939	2.27939	0.010	1.00413	50.00000	Averaged
24 Phenol	2.03729	2.08653	2.08653	0.010	2.41708	20.00000	Averaged
26 Bis(2-chloroethyl)ether	1.42859	1.46102	1.46102	0.010	2.26983	50.00000	Averaged
27 2-Chlorophenol	1.56381	1.59024	1.59024	0.010	1.68995	50.00000	Averaged
28 1,3-Dichlorobenzene	1.70337	1.79554	1.79554	0.010	5.41089	50.00000	Averaged
29 1,4-Dichlorobenzene	1.78118	1.80374	1.80374	0.010	1.26683	20.00000	Averaged
30 Benzyl Alcohol	1.05101	1.07490	1.07490	0.010	2.27311	50.00000	Averaged
31 1,2-Dichlorobenzene	1.63746	1.69840	1.69840	0.010	3.72162	50.00000	Averaged
32 2-Methylphenol	1.43012	1.43294	1.43294	0.010	0.19697	50.00000	Averaged
33 2,2'-oxybis(1-Chloropropane	2.27365	2.27329	2.27329	0.010	-0.01588	50.00000	Averaged
34 4-Methylphenol	1.51904	1.51474	1.51474	0.010	-0.28351	50.00000	Averaged
36 Hexachloroethane	0.60636	0.65572	0.65572	0.010	8.13887	50.00000	Averaged
37 N-Nitrosodinpropylamine	1.01180	1.02542	1.02542	0.050	1.34598	50.00000	Averaged
42 Nitrobenzene	0.33116	0.34537	0.34537	0.010	4.28887	50.00000	Averaged
44 Isophorone	0.63679	0.65503	0.65503	0.010	2.86481	50.00000	Averaged
45 2-Nitrophenol	0.19648	0.20808	0.20808	0.010	5.90704	20.00000	Averaged
46 2,4-Dimethyphenol	0.34911	0.36705	0.36705	0.010	5.13622	50.00000	Averaged
47 Bis(2-chloroethoxy)methane	0.38908	0.38995	0.38995	0.010	0.22393	50.00000	Averaged
49 2,4-Dichlorophenol	0.27010	0.27837	0.27837	0.010	3.06220	20.00000	Averaged
50 Benzoic Acid	0.19324	0.20608	0.20608	0.010	6.64448	50.00000	Averaged
51 1,2,4-Trichlorobenzene	0.29246	0.31045	0.31045	0.010	6.15434	50.00000	Averaged
52 Naphthalene	1.10443	1.13982	1.13982	0.010	3.20449	50.00000	Averaged
54 4-Chloroaniline	0.43288	0.44148	0.44148	0.010	1.98835	50.00000	Averaged
57 Hexachlorobutadiene	0.14313	0.15626	0.15626	0.010	9.17619	20.00000	Averaged
60 4-Chloro-3-Methylphenol	0.30164	0.31836	0.31836	0.010	5.54416	20.00000	Averaged
63 2-Methylnaphthalene	0.69378	0.73392	0.73392	0.010	5.78600	50.00000	Averaged
66 Hexachlorocyclopentadiene	0.29846	0.32035	0.32035	0.050	7.33524	50.00000	Averaged
69 2,4,6-Trichlorophenol	0.31913	0.33680	0.33680	0.010	5.53446	20.00000	Averaged
70 2,4,5-Trichlorphenol	0.34380	0.36886	0.36886	0.010	7.28819	50.00000	Averaged
71 2-Chloronaphthalene	1.12571	1.15486	1.15486	0.010	2.58993	50.00000	Averaged
73 2-Nitroaniline	0.34119	0.36528	0.36528	0.010	7.06266	50.00000	Averaged
76 Dimethylphthalate	1.29606	1.31651	1.31651	0.010	1.57746	50.00000	Averaged

Manual calculation for Nitrobenzene:

$\frac{252501}{584888} \times \frac{40}{50} = 0.34537$ BY 12/6/10

12/6/10

TestAmerica West Sacramento

CONTINUING CALIBRATION COMPOUNDS

Instrument ID: sv5.i Injection Date: 04-DEC-2010 10:36
Lab File ID: HSL1204.D Init. Cal. Date(s): 17-AUG-2010 02-OCT-2010
Analysis Type: Init. Cal. Times: 17:32 15:00
Lab Sample ID: HSL_050 ug/ml CS-4 Quant Type: ISTD
Method: \\SV5\C\chem\sv5.i\120410.B\8270f.m

COMPOUND	RRF / AMOUNT	RF50	CCAL RRF50	MIN RRF	%D / %DRIFT	MAX %D / %DRIFT	CURVE TYPE
77 Acenaphthylene	1.96037	1.99664	1.99664	0.010	1.85039	50.00000	Averaged
79 2,6-Dinitrotoluene	0.30197	0.30862	0.30862	0.010	2.20268	50.00000	Averaged
80 3-Nitroaniline	0.37691	0.38778	0.38778	0.010	2.88550	50.00000	Averaged
81 Acenaphthene	1.24787	1.26244	1.26244	0.010	1.16789	20.00000	Averaged
82 2,4-Dinitrophenol	50.00000	52.52788	0.18729	0.050	5.05576	0.000e+000	Quadratic
83 Dibenzofuran	1.65612	1.69133	1.69133	0.010	2.12631	50.00000	Averaged
84 4-Nitrophenol	0.15634	0.17314	0.17314	0.050	10.74786	50.00000	Averaged
86 2,4-Dinitrotoluene	0.39633	0.42188	0.42188	0.010	6.44559	50.00000	Averaged
91 Fluorene	1.37139	1.39249	1.39249	0.010	1.53868	50.00000	Averaged
92 Diethylphthalate	1.32699	1.37656	1.37656	0.010	3.73540	50.00000	Averaged
93 4-Chlorophenyl-phenylether	0.57019	0.57433	0.57433	0.010	0.72546	50.00000	Averaged
94 4-Nitroaniline	0.37361	0.39990	0.39990	0.010	7.03707	50.00000	Averaged
97 4,6-Dinitro-2-methylphenol	50.00000	54.80273	0.15726	0.010	9.60546	0.000e+000	Linear
98 N-Nitrosodiphenylamine	0.60628	0.62806	0.62806	0.010	3.59210	20.00000	Averaged
100 Azobenzene	0.78660	0.80903	0.80903	0.010	2.85135	50.00000	Averaged
101 4-Bromophenyl-phenylether	0.19527	0.21480	0.21480	0.010	10.00406	50.00000	Averaged
108 Hexachlorobenzene	0.21807	0.22934	0.22934	0.010	5.17043	50.00000	Averaged
110 Pentachlorophenol	50.00000	48.80415	0.12675	0.010	-2.39170	0.000e+000	Linear
114 Phenanthrene	1.26074	1.27780	1.27780	0.010	1.35252	50.00000	Averaged
115 Anthracene	1.25955	1.31316	1.31316	0.010	4.25662	50.00000	Averaged
118 Carbazole	1.15061	1.17061	1.17061	0.010	1.73845	50.00000	Averaged
120 Di-n-Butylphthalate	1.38442	1.46886	1.46886	0.010	6.09975	50.00000	Averaged
126 Fluoranthene	1.12969	1.22360	1.22360	0.010	8.31338	20.00000	Averaged
127 Benzidine	0.81067	0.86515	0.86515	0.010	6.71949	50.00000	Averaged
128 Pyrene	1.25025	1.26571	1.26571	0.010	1.23579	50.00000	Averaged
134 3,3'-dimethylbenzidine	0.71564	0.76105	0.76105	0.010	6.34630	50.00000	Averaged
136 Butylbenzylphthalate	0.62663	0.67343	0.67343	0.010	7.46792	50.00000	Averaged
138 Benzo(a)Anthracene	1.06548	1.12285	1.12285	0.010	5.38490	50.00000	Averaged
139 Chrysene	1.08994	1.07955	1.07955	0.010	-0.95314	50.00000	Averaged
140 3,3'-Dichlorobenzidine	0.40189	0.42573	0.42573	0.010	5.93261	50.00000	Averaged
141 bis(2-ethylhexyl)Phthalate	0.86316	0.90793	0.90793	0.010	5.18612	50.00000	Averaged
142 Di-n-octylphthalate	1.37975	1.53772	1.53772	0.010	11.44900	20.00000	Averaged
144 Benzo(b)fluoranthene	0.90549	1.10333	1.10333	0.010	21.84870	50.00000	Averaged
145 Benzo(k)fluoranthene	1.16236	1.07851	1.07851	0.010	-7.21363	50.00000	Averaged
147 Benzo(e)pyrene	0.94425	0.98783	0.98783	0.010	4.61585	50.00000	Averaged
148 Benzo(a)pyrene	1.02655	1.07597	1.07597	0.010	4.81409	20.00000	Averaged
151 Indeno(1,2,3-cd)pyrene	0.83029	0.94746	0.94746	0.010	14.11161	50.00000	Averaged
152 Dibenzo(a,h)anthracene	0.92758	1.02615	1.02615	0.010	10.62663	50.00000	Averaged
153 Benzo(g,h,i)perylene	1.00427	1.09166	1.09166	0.010	8.70201	50.00000	Averaged
M 162 benzo b,k Fluoranthene Tota	2.06785	2.18184	2.18184	0.010	5.51256	50.00000	Averaged

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\120410.B\HSL1204.D
Lab Smp Id: HSL_050 ug/ml CS-4 Client Smp ID: 8270F.M
Inj Date : 04-DEC-2010 10:36
Operator : KT Inst ID: sv5.i
Smp Info : HSL_050 ug/ml CS-4;2;;4;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0310;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\120410.B\8270f.m
Meth Date : 06-Dec-2010 09:21 semivoa Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 97 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	138327	40.0000	
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	584888	40.0000	
* 3	Acenaphthene-d10	164	7.045	7.045	(1.000)	316560	40.0000	
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	498828	40.0000	
* 5	Chrysene-d12	240	13.190	13.190	(1.000)	516170	40.0000	
* 6	Perylene-d12	264	15.543	15.543	(1.000)	510499	40.0000	
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	253146	50.0000	51.92
$ 8	Phenol-d5	99	3.231	3.231	(0.912)	317072	50.0000	51.71
$ 9	2-Chlorophenol-d4	132	3.345	3.345	(0.944)	279882	50.0000	51.98
$ 10	1,2-Dichlorobenzene-d4	152	3.739	3.739	(1.056)	176180	50.0000	51.72
$ 11	Nitrobenzene-d5	82	4.174	4.174	(0.843)	255477	50.0000	51.57
$ 12	2-Fluorobiphenyl	172	6.257	6.257	(0.888)	512661	50.0000	50.27
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	74261	50.0000	53.98
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	522707	50.0000	51.41
15	N-Nitrosodimethylamine	74	1.314	1.314	(0.371)	157102	50.0000	49.30
16	Pyridine	79	1.325	1.325	(0.374)	260357	50.0000	48.85
23	Aniline	93	3.252	3.252	(0.918)	394126	50.0000	50.50
24	Phenol	94	3.252	3.252	(0.918)	360779	50.0000	51.21
26	Bis(2-chloroethyl)ether	93	3.314	3.314	(0.936)	252623	50.0000	51.13
27	2-Chlorophenol	128	3.356	3.356	(0.947)	274966	50.0000	50.84
28	1,3-Dichlorobenzene	146	3.501	3.501	(0.988)	310465	50.0000	52.70
29	1,4-Dichlorobenzene	146	3.553	3.553	(1.003)	311883	50.0000	50.63
30	Benzyl Alcohol	108	3.718	3.718	(1.050)	185860	50.0000	51.14
31	1,2-Dichlorobenzene	146	3.760	3.760	(1.061)	293669	50.0000	51.86
32	2-Methylphenol	108	3.874	3.874	(1.094)	247768	50.0000	50.10
33	2,2'-oxybis(1-Chloropropane)	45	3.895	3.895	(1.099)	393072	50.0000	49.99
34	4-Methylphenol	108	4.040	4.040	(1.140)	261911	50.0000	49.86
36	Hexachloroethane	117	4.081	4.081	(1.152)	113379	50.0000	54.07
37	N-Nitrosodinpropylamine	70	4.040	4.040	(1.140)	177304	50.0000	50.67
42	Nitrobenzene	77	4.185	4.185	(0.845)	252501	50.0000	52.14
44	Isophorone	82	4.454	4.454	(0.900)	478900	50.0000	51.43
45	2-Nitrophenol	139	4.547	4.547	(0.918)	152132	50.0000	52.95
46	2,4-Dimethyphenol	107	4.630	4.630	(0.935)	268351	50.0000	52.57

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	4.734	4.734	(0.956)	285100	50.0000	50.11
49	2,4-Dichlorophenol	162	4.827	4.827	(0.975)	203520	50.0000	51.53
50	Benzoic Acid	122	4.765	4.765	(0.962)	150668	50.0000	53.32
51	1,2,4-Trichlorobenzene	180	4.910	4.910	(0.992)	226976	50.0000	53.08
52	Naphthalene	128	4.972	4.972	(1.004)	833332	50.0000	51.60
54	4-Chloroaniline	127	5.086	5.086	(1.027)	322772	50.0000	50.99
57	Hexachlorobutadiene	225	5.200	5.200	(1.050)	114243	50.0000	54.59
60	4-Chloro-3-Methylphenol	107	5.687	5.687	(1.149)	232755	50.0000	52.77
63	2-Methylnaphthalene	142	5.915	5.915	(1.195)	536574	50.0000	52.89
66	Hexachlorocyclopentadiene	237	6.060	6.060	(0.860)	126762	50.0000	53.67
69	2,4,6-Trichlorophenol	196	6.174	6.174	(0.876)	133270	50.0000	52.77
70	2,4,5-Trichlorphenol	196	6.216	6.216	(0.882)	145957	50.0000	53.64
71	2-Chloronaphthalene	162	6.361	6.361	(0.903)	456980	50.0000	51.29
73	2-Nitroaniline	65	6.537	6.537	(0.928)	144542	50.0000	53.53
76	Dimethylphthalate	163	6.817	6.817	(0.968)	520942	50.0000	50.79
77	Acenaphthylene	152	6.858	6.858	(0.974)	790071	50.0000	50.92
79	2,6-Dinitrotoluene	165	6.889	6.889	(0.978)	122120	50.0000	51.10
80	3-Nitroaniline	138	7.045	7.045	(1.000)	153445	50.0000	51.44
81	Acenaphthene	153	7.076	7.076	(1.004)	499548	50.0000	50.58
82	2,4-Dinitrophenol	184	7.169	7.169	(1.018)	74110	50.0000	52.53
83	Dibenzofuran	168	7.273	7.273	(1.032)	669261	50.0000	51.06
84	4-Nitrophenol	109	7.304	7.304	(1.037)	68513	50.0000	55.37
86	2,4-Dinitrotoluene	165	7.356	7.356	(1.044)	166937	50.0000	53.22
91	Fluorene	166	7.687	7.687	(1.091)	551010	50.0000	50.77
92	Diethylphthalate	149	7.677	7.677	(1.090)	544705	50.0000	51.87
93	4-Chlorophenyl-phenylether	204	7.708	7.708	(1.094)	227261	50.0000	50.36
94	4-Nitroaniline	138	7.781	7.781	(1.104)	158242	50.0000	53.52
97	4,6-Dinitro-2-methylphenol	198	7.843	7.843	(0.881)	98059	50.0000	54.80
98	N-Nitrosodiphenylamine	169	7.874	7.874	(0.885)	458977	58.6000	60.70
100	Azobenzene	77	7.905	7.905	(0.888)	504459	50.0000	51.42
101	4-Bromophenyl-phenylether	248	8.330	8.330	(0.936)	133936	50.0000	55.00
108	Hexachlorobenzene	284	8.496	8.496	(0.955)	143002	50.0000	52.58
110	Pentachlorophenol	266	8.755	8.755	(0.984)	79036	50.0000	48.80
114	Phenanthrene	178	8.931	8.931	(1.003)	796750	50.0000	50.68
115	Anthracene	178	8.993	8.993	(1.010)	818801	50.0000	52.13
118	Carbazole	167	9.263	9.263	(1.041)	729917	50.0000	50.87
120	Di-n-Butylphthalate	149	9.957	9.957	(1.119)	915888	50.0000	53.05
126	Fluoranthene	202	10.745	10.745	(1.207)	762960	50.0000	54.16
127	Benzidine	184	11.035	11.035	(0.837)	558203	50.0000	53.36
128	Pyrene	202	11.087	11.087	(0.841)	816649	50.0000	50.62
134	3,3'-dimethylbenzidine	212	12.309	12.309	(0.933)	491042	50.0000	53.17
136	Butylbenzylphthalate	149	12.434	12.434	(0.943)	434505	50.0000	53.73
138	Benzo(a)Anthracene	228	13.159	13.159	(0.998)	724478	50.0000	52.69
139	Chrysene	228	13.232	13.232	(1.003)	696538	50.0000	49.52
140	3,3'-Dichlorobenzidine	252	13.221	13.221	(1.002)	274689	50.0000	52.97
141	bis(2-ethylhexyl)Phthalate	149	13.563	13.563	(1.028)	585805	50.0000	52.59
142	Di-n-octylphthalate	149	14.610	14.610	(1.108)	992154	50.0000	55.72
144	Benzo(b)fluoranthene	252	14.973	14.973	(0.963)	704060	50.0000	60.92
145	Benzo(k)fluoranthene	252	15.014	15.014	(0.966)	688223	50.0000	46.39
147	Benzo(e)pyrene	252	15.387	15.387	(0.990)	630359	50.0000	52.31
148	Benzo(a)pyrene	252	15.460	15.460	(0.995)	686601	50.0000	52.41
151	Indeno(1,2,3-cd)pyrene	276	17.076	17.076	(1.099)	604594	50.0000	57.06
152	Dibenzo(a,h)anthracene	278	17.118	17.118	(1.101)	654808	50.0000	55.31
153	Benzo(g,h,i)perylene	276	17.439	17.439	(1.122)	696616	50.0000	54.35

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				1392283	50.0000	52.76(A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\120410.B\HSL1204.D
Lab Smp Id: HSL_050 ug/ml CS-4 Client Smp ID: 8270F.M
Inj Date : 04-DEC-2010 10:36
Operator : KT Inst ID: sv5.i
Smp Info : HSL_050 ug/ml CS-4;2;;4;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0310;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\120410.B\8270f.m
Meth Date : 06-Dec-2010 09:21 semivoa Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 97 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	138327	40.0000	
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	584888	40.0000	
* 3	Acenaphthene-d10	164	7.045	7.045	(1.000)	316560	40.0000	
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	498828	40.0000	
* 5	Chrysene-d12	240	13.190	13.190	(1.000)	516170	40.0000	
* 6	Perylene-d12	264	15.543	15.543	(1.000)	510499	40.0000	
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	253146	50.0000	51.92
$ 8	Phenol-d5	99	3.231	3.231	(0.912)	317072	50.0000	51.71
$ 9	2-Chlorophenol-d4	132	3.345	3.345	(0.944)	279882	50.0000	51.98
$ 10	1,2-Dichlorobenzene-d4	152	3.739	3.739	(1.056)	176180	50.0000	51.72
$ 11	Nitrobenzene-d5	82	4.174	4.174	(0.843)	255477	50.0000	51.57
$ 12	2-Fluorobiphenyl	172	6.257	6.257	(0.888)	512661	50.0000	50.27
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	74261	50.0000	53.98
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	522707	50.0000	51.41
15	N-Nitrosodimethylamine	74	1.314	1.314	(0.371)	157102	50.0000	49.30
16	Pyridine	79	1.325	1.325	(0.374)	260357	50.0000	48.85
23	Aniline	93	3.252	3.252	(0.918)	394126	50.0000	50.50
24	Phenol	94	3.252	3.252	(0.918)	360779	50.0000	51.21
26	Bis(2-chloroethyl)ether	93	3.314	3.314	(0.936)	252623	50.0000	51.13
27	2-Chlorophenol	128	3.356	3.356	(0.947)	274966	50.0000	50.84
28	1,3-Dichlorobenzene	146	3.501	3.501	(0.988)	310465	50.0000	52.70
29	1,4-Dichlorobenzene	146	3.553	3.553	(1.003)	311883	50.0000	50.63
30	Benzyl Alcohol	108	3.718	3.718	(1.050)	185860	50.0000	51.14
31	1,2-Dichlorobenzene	146	3.760	3.760	(1.061)	293669	50.0000	51.86
32	2-Methylphenol	108	3.874	3.874	(1.094)	247768	50.0000	50.10
33	2,2'-oxybis(1-Chloropropane)	45	3.895	3.895	(1.099)	393072	50.0000	49.99
34	4-Methylphenol	108	4.040	4.040	(1.140)	261911	50.0000	49.86
36	Hexachloroethane	117	4.081	4.081	(1.152)	113379	50.0000	54.07
37	N-Nitrosodinpropylamine	70	4.040	4.040	(1.140)	177304	50.0000	50.67
42	Nitrobenzene	77	4.185	4.185	(0.845)	252501	50.0000	52.14
44	Isophorone	82	4.454	4.454	(0.900)	478900	50.0000	51.43
45	2-Nitrophenol	139	4.547	4.547	(0.918)	152132	50.0000	52.95
46	2,4-Dimethyphenol	107	4.630	4.630	(0.935)	268351	50.0000	52.57

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	4.734	4.734	(0.956)	285100	50.0000	50.11
49	2,4-Dichlorophenol	162	4.827	4.827	(0.975)	203520	50.0000	51.53
50	Benzoic Acid	122	4.765	4.765	(0.962)	150668	50.0000	53.32
51	1,2,4-Trichlorobenzene	180	4.910	4.910	(0.992)	226976	50.0000	53.08
52	Naphthalene	128	4.972	4.972	(1.004)	833332	50.0000	51.60
54	4-Chloroaniline	127	5.086	5.086	(1.027)	322772	50.0000	50.99
57	Hexachlorobutadiene	225	5.200	5.200	(1.050)	114243	50.0000	54.59
60	4-Chloro-3-Methylphenol	107	5.687	5.687	(1.149)	232755	50.0000	52.77
63	2-Methylnaphthalene	142	5.915	5.915	(1.195)	536574	50.0000	52.89
66	Hexachlorocyclopentadiene	237	6.060	6.060	(0.860)	126762	50.0000	53.67
69	2,4,6-Trichlorophenol	196	6.174	6.174	(0.876)	133270	50.0000	52.77
70	2,4,5-Trichlorphenol	196	6.216	6.216	(0.882)	145957	50.0000	53.64
71	2-Chloronaphthalene	162	6.361	6.361	(0.903)	456980	50.0000	51.29
73	2-Nitroaniline	65	6.537	6.537	(0.928)	144542	50.0000	53.53
76	Dimethylphthalate	163	6.817	6.817	(0.968)	520942	50.0000	50.79
77	Acenaphthylene	152	6.858	6.858	(0.974)	790071	50.0000	50.92
79	2,6-Dinitrotoluene	165	6.889	6.889	(0.978)	122120	50.0000	51.10
80	3-Nitroaniline	138	7.045	7.045	(1.000)	153445	50.0000	51.44
81	Acenaphthene	153	7.076	7.076	(1.004)	499548	50.0000	50.58
82	2,4-Dinitrophenol	184	7.169	7.169	(1.018)	74110	50.0000	52.53
83	Dibenzofuran	168	7.273	7.273	(1.032)	669261	50.0000	51.06
84	4-Nitrophenol	109	7.304	7.304	(1.037)	68513	50.0000	55.37
86	2,4-Dinitrotoluene	165	7.356	7.356	(1.044)	166937	50.0000	53.22
91	Fluorene	166	7.687	7.687	(1.091)	551010	50.0000	50.77
92	Diethylphthalate	149	7.677	7.677	(1.090)	544705	50.0000	51.87
93	4-Chlorophenyl-phenylether	204	7.708	7.708	(1.094)	227261	50.0000	50.36
94	4-Nitroaniline	138	7.781	7.781	(1.104)	158242	50.0000	53.52
97	4,6-Dinitro-2-methylphenol	198	7.843	7.843	(0.881)	98059	50.0000	54.80
98	N-Nitrosodiphenylamine	169	7.874	7.874	(0.885)	458977	58.6000	60.70
100	Azobenzene	77	7.905	7.905	(0.888)	504459	50.0000	51.42
101	4-Bromophenyl-phenylether	248	8.330	8.330	(0.936)	133936	50.0000	55.00
108	Hexachlorobenzene	284	8.496	8.496	(0.955)	143002	50.0000	52.58
110	Pentachlorophenol	266	8.755	8.755	(0.984)	79036	50.0000	48.80
114	Phenanthrene	178	8.931	8.931	(1.003)	796750	50.0000	50.68
115	Anthracene	178	8.993	8.993	(1.010)	818801	50.0000	52.13
118	Carbazole	167	9.263	9.263	(1.041)	729917	50.0000	50.87
120	Di-n-Butylphthalate	149	9.957	9.957	(1.119)	915888	50.0000	53.05
126	Fluoranthene	202	10.745	10.745	(1.207)	762960	50.0000	54.16
127	Benzidine	184	11.035	11.035	(0.837)	558203	50.0000	53.36
128	Pyrene	202	11.087	11.087	(0.841)	816649	50.0000	50.62
134	3,3'-dimethylbenzidine	212	12.309	12.309	(0.933)	491042	50.0000	53.17
136	Butylbenzylphthalate	149	12.434	12.434	(0.943)	434505	50.0000	53.73
138	Benzo(a)Anthracene	228	13.159	13.159	(0.998)	724478	50.0000	52.69
139	Chrysene	228	13.232	13.232	(1.003)	696538	50.0000	49.52
140	3,3'-Dichlorobenzidine	252	13.221	13.221	(1.002)	274689	50.0000	52.97
141	bis(2-ethylhexyl)Phthalate	149	13.563	13.563	(1.028)	585805	50.0000	52.59
142	Di-n-octylphthalate	149	14.610	14.610	(1.108)	992154	50.0000	55.72
144	Benzo(b)fluoranthene	252	14.973	14.973	(0.963)	704060	50.0000	60.92
145	Benzo(k)fluoranthene	252	15.014	15.014	(0.966)	688223	50.0000	46.39
147	Benzo(e)pyrene	252	15.387	15.387	(0.990)	630359	50.0000	52.31
148	Benzo(a)pyrene	252	15.460	15.460	(0.995)	686601	50.0000	52.41
151	Indeno(1,2,3-cd)pyrene	276	17.076	17.076	(1.099)	604594	50.0000	57.06
152	Dibenzo(a,h)anthracene	278	17.118	17.118	(1.101)	654808	50.0000	55.31
153	Benzo(g,h,i)perylene	276	17.439	17.439	(1.122)	696616	50.0000	54.35

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				1392283	50.0000	52.76(A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1204.D
Lab Smp Id: HSL_050 ug/ml CS-4
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\SV5\C\chem\sv5.i\120410.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0310;0;8270F.M

Calibration Date: 04-DEC-2010
Calibration Time: 09:53
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	138327	12.80
2 Naphthalene-d8	530514	265257	1061028	584888	10.25
3 Acenaphthene-d10	282538	141269	565076	316560	12.04
4 Phenanthrene-d10	462722	231361	925444	498828	7.80
5 Chrysene-d12	435850	217925	871700	516170	18.43
6 Perylene-d12	422284	211142	844568	510499	20.89

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.54	3.04	4.04	3.54	0.00
2 Naphthalene-d8	4.95	4.45	5.45	4.95	0.00
3 Acenaphthene-d10	7.05	6.55	7.55	7.05	0.00
4 Phenanthrene-d10	8.90	8.40	9.40	8.90	0.00
5 Chrysene-d12	13.19	12.69	13.69	13.19	0.00
6 Perylene-d12	15.54	15.04	16.04	15.54	0.00

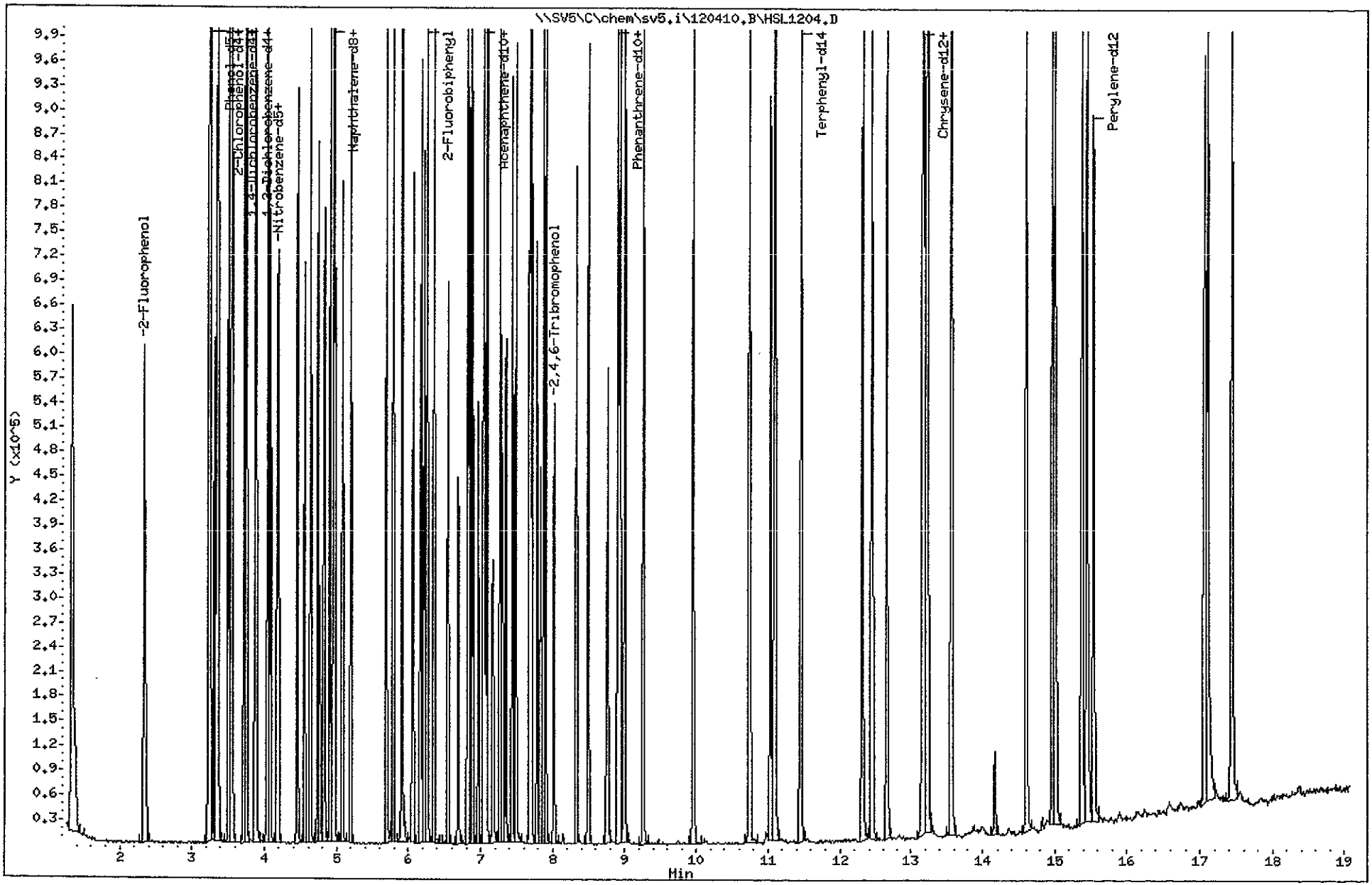
AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV5\C\chem\sv5.i\120410.B\HSL1204.D
Date : 04-DEC-2010 10:36
Client ID: 8270F.M
Sample Info: HSL_050 ug/ml CS-4;2;;4;;;;;4

Instrument: sv5.i

Operator: KT
Column diameter: 2.00

Column phase:



TAILING FACTOR/DEGRADATION SUMMARY RESULTS

TAILING ANALYSIS SUMMARY

Compound	Tail Factor	Max Allowed	Test
Pentachlorophenol	0.9091184	5.000	PASS
Benzidine	0.3577715	3.000	PASS

DDT DEGRADATION BREAKDOWN ANALYSIS SUMMARY

Compound	Response	%Breakdown	Max Allowed	Test
4,4-DDD + DDE	197771	9.6	20.5	PASS

****
*** PASSED ***
****

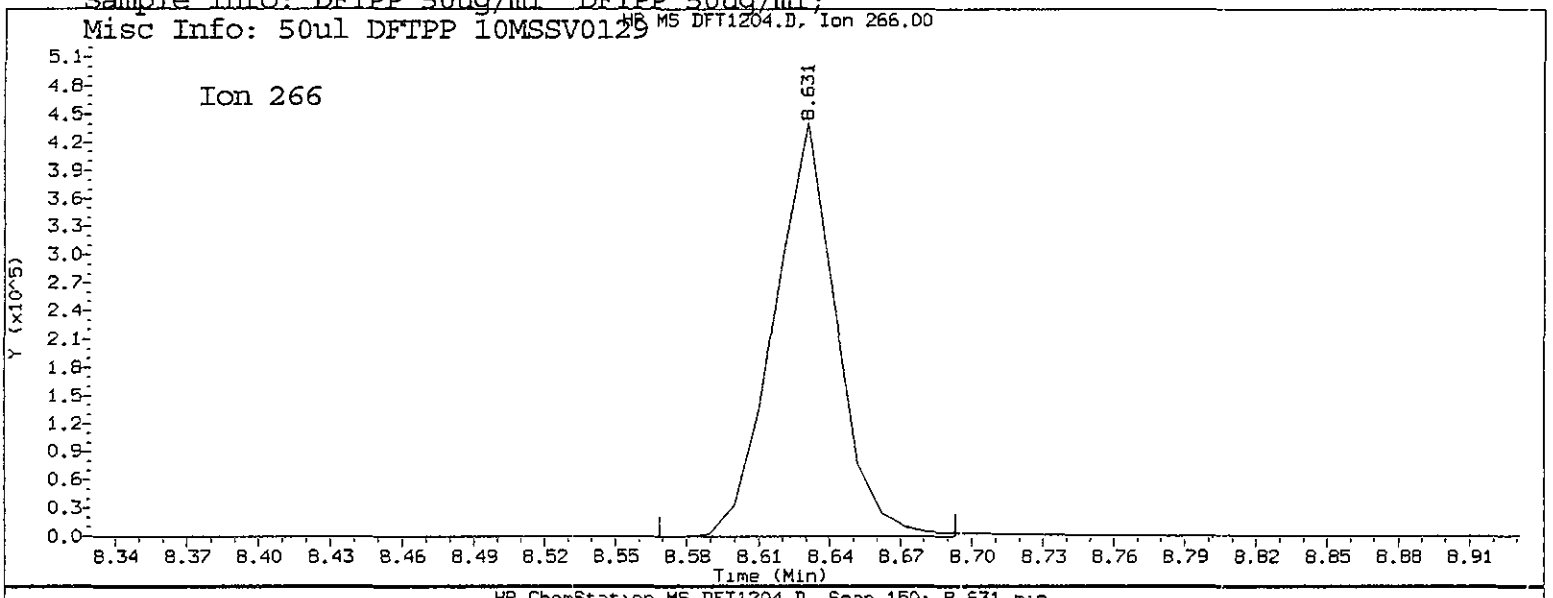
Sample //SV5/C/chem/sv5.i/120410.B/DFT1204.D/DFT1204.D

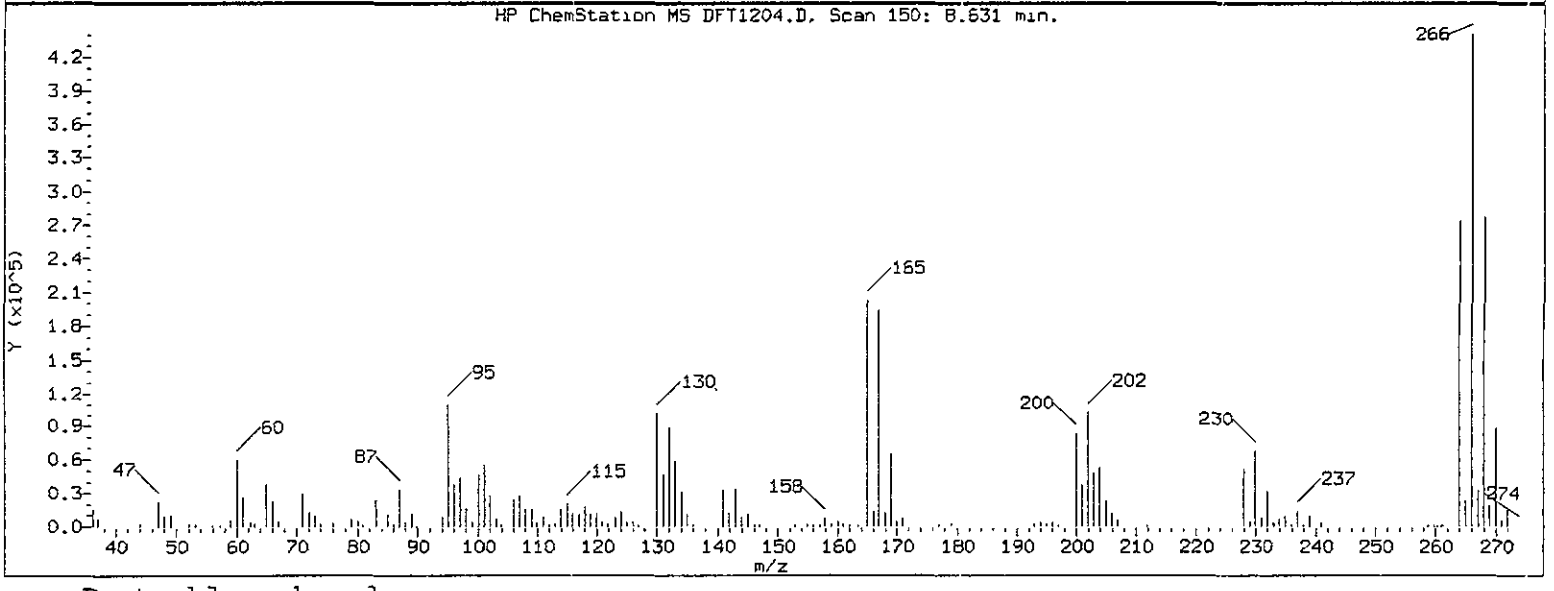
12/6/10

TAILING FACTOR/DEGRADATION SAMPLE AND GRAPHIC REPORT

Report Date: 12/06/2010 09:20

Datafile Analyzed: //SV5/C/chem/sv5.i/120410.B/DFT1204.D/DFT1204.D
Method Used: \\SV5\C\chem\sv5.i\120410.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 04-DEC-2010 10:15 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129





Pentachlorophenol
==========
Exp. RT = 8.662
Found RT = 8.631

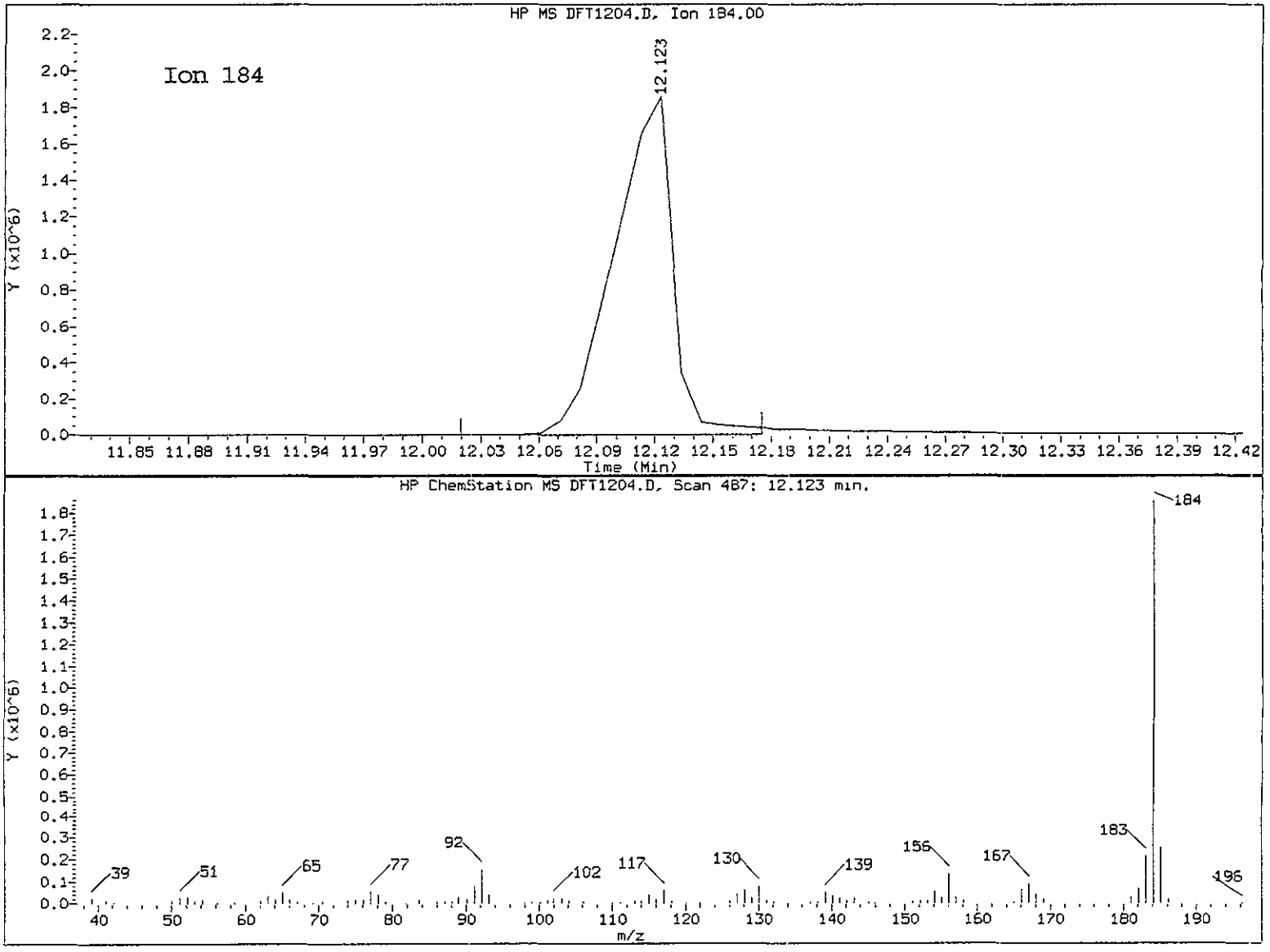
Time1 = 8.600923 Time2 = 8.630883 Time3 = 8.658121
Tailing Factor = (Time3 - Time2)/(Time2 - Time1)

Tailing factor for Pentachlorophenol OK

Tail Factor = 0.909 Maximum Allowed = 5.0

Report Date: 12/06/2010 09:20

Datafile Analyzed: //SV5/C/chem/sv5.i/120410.B/DFT1204.D/DFT1204.D
Method Used: \\SV5\C\chem\sv5.i\120410.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 04-DEC-2010 10:15 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129



Benzidine
=========
Exp. RT = 12.144
Found RT = 12.123

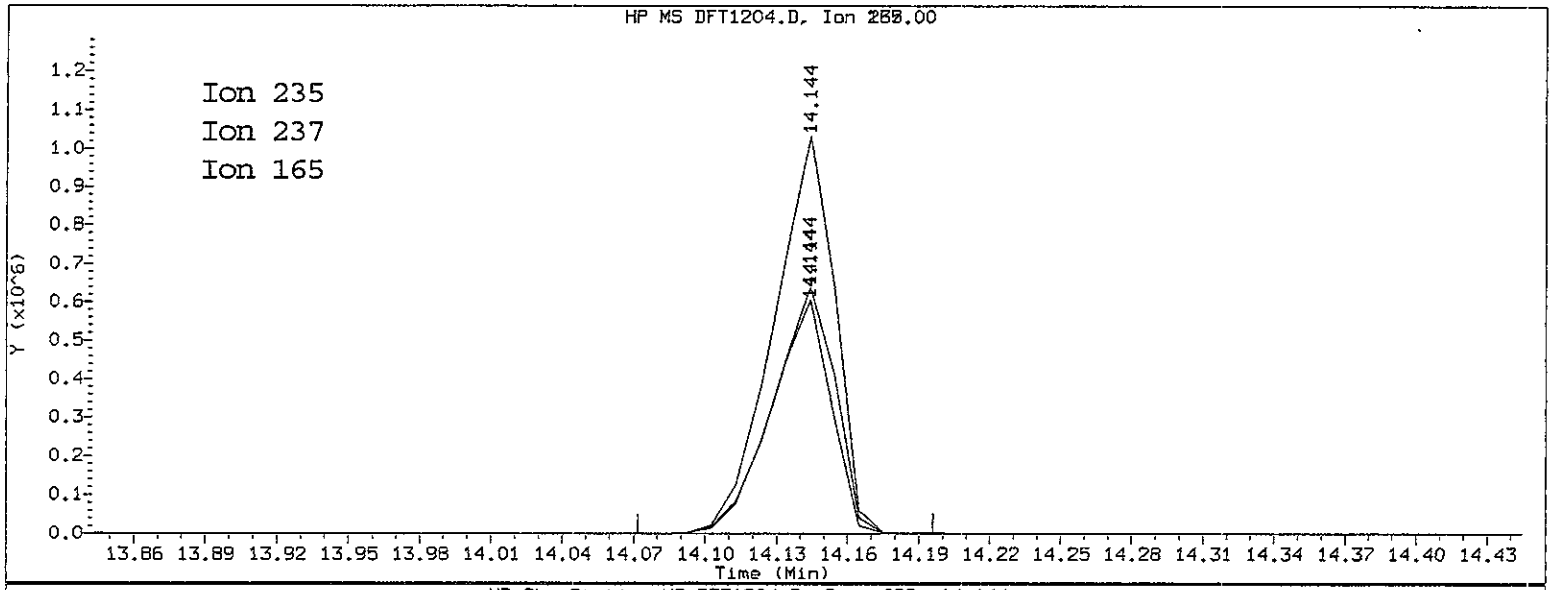
Time1 = 12.07753 Time2 = 12.12323 Time3 = 12.13958
Tailing Factor = (Time3 - Time2)/(Time2 - Time1)

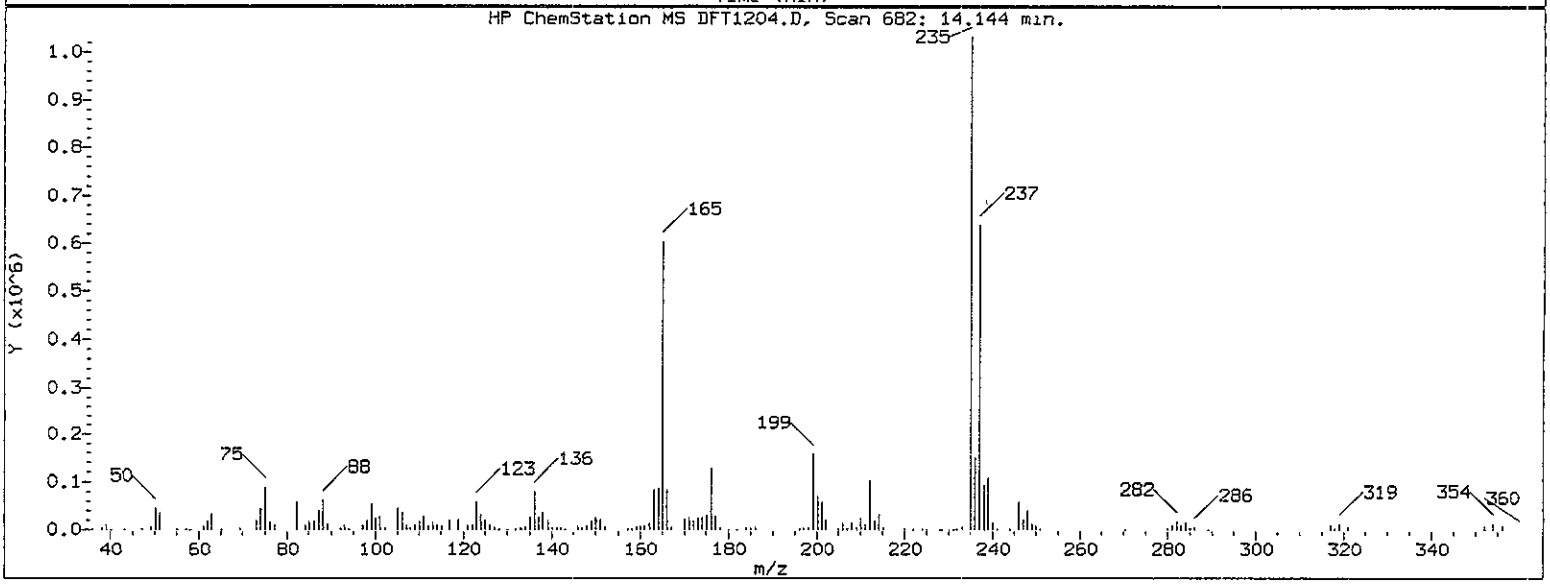
Tailing factor for Benzidine OK

Tail Factor = 0.358 Maximum Allowed = 3.0

Report Date: 12/06/2010 09:20

Datafile Analyzed: //SV5/C/chem/sv5.i/120410.B/DFT1204.D/DFT1204.D
Method Used: \\SV5\C\chem\sv5.i\120410.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 04-DEC-2010 10:15 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129



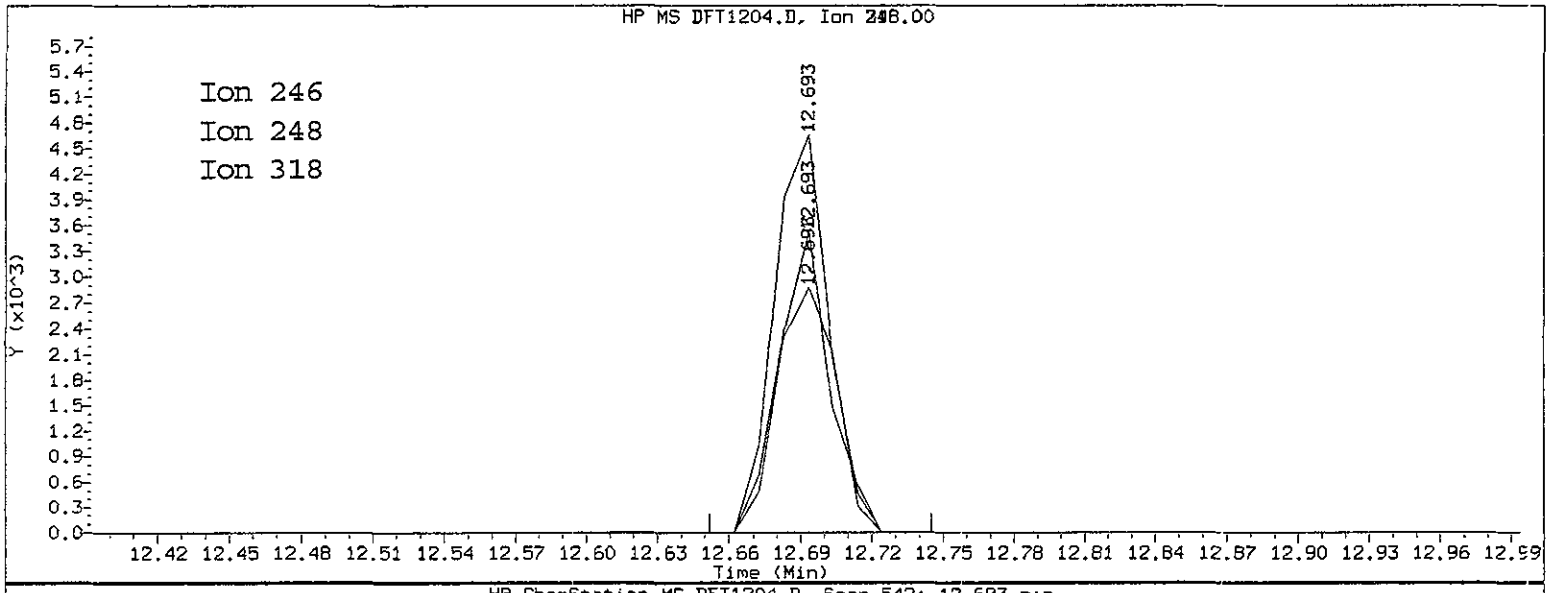


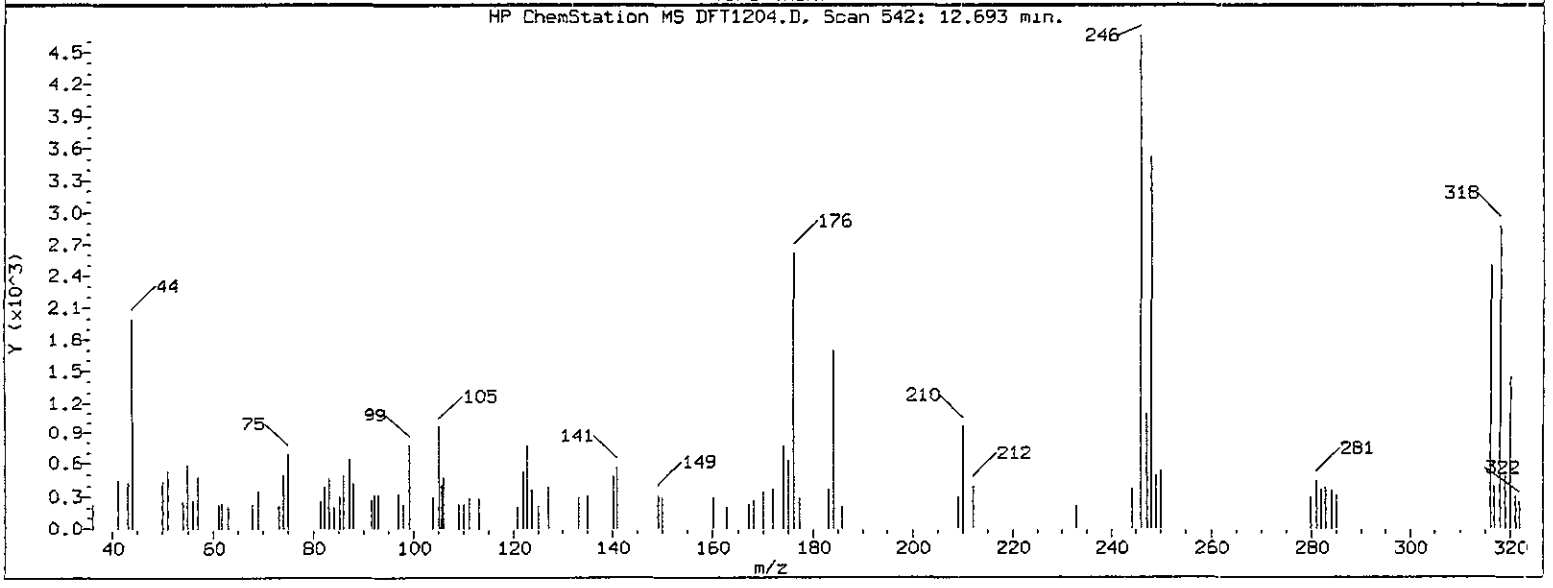
4,4'-DDT
========
Exp. RT = 14.165
Found RT = 14.144

Mass	Area	Ratio
235	1852497	100.00
237	1164121	62.84
165	1063113	57.39

Report Date: 12/06/2010 09:20

Datafile Analyzed: //SV5/C/chem/sv5.i/120410.B/DFT1204.D/DFT1204.D
Method Used: \\SV5\C\chem\sv5.i\120410.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 04-DEC-2010 10:15 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129



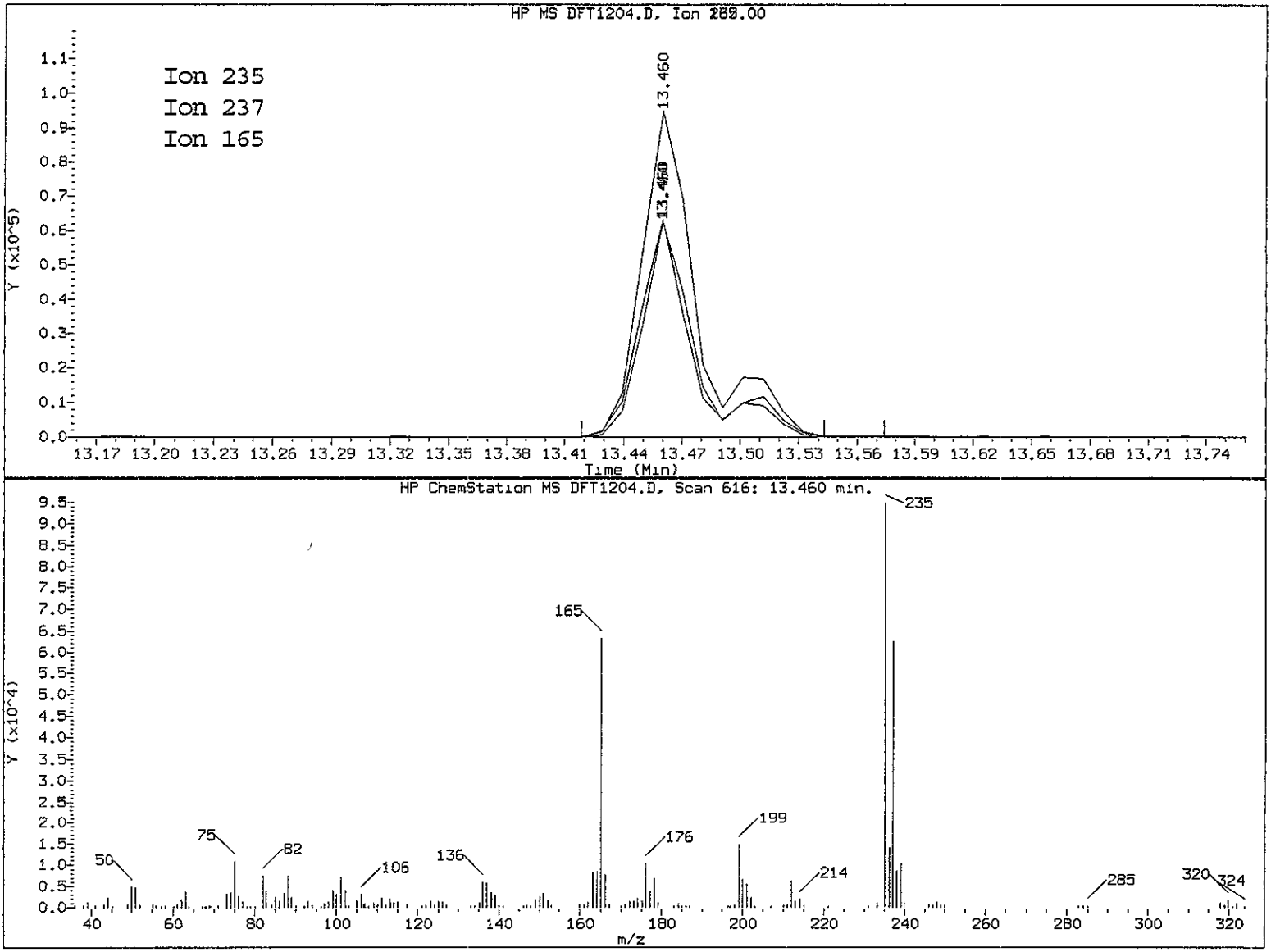


4,4'-DDE
========
Exp. RT = 12.714
Found RT = 12.693

Mass	Area	Ratio
246	7573	100.00
248	5386	71.12
318	5062	66.85

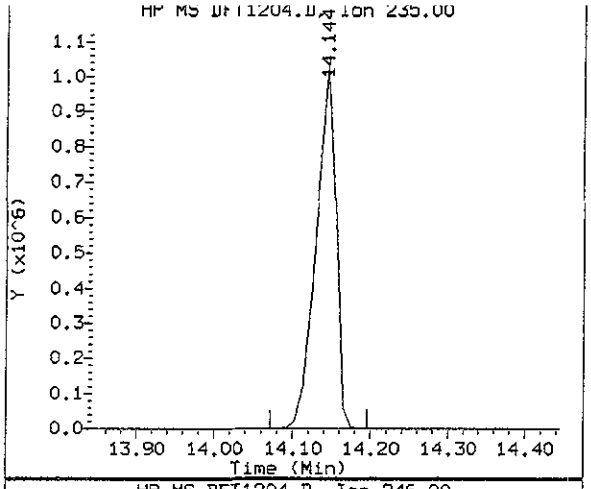
Report Date: 12/06/2010 09:20

Datafile Analyzed: //SV5/C/chem/sv5.i/120410.B/DFT1204.D/DFT1204.D
Method Used: \\SV5\C\chem\sv5.i\120410.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 04-DEC-2010 10:15 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129

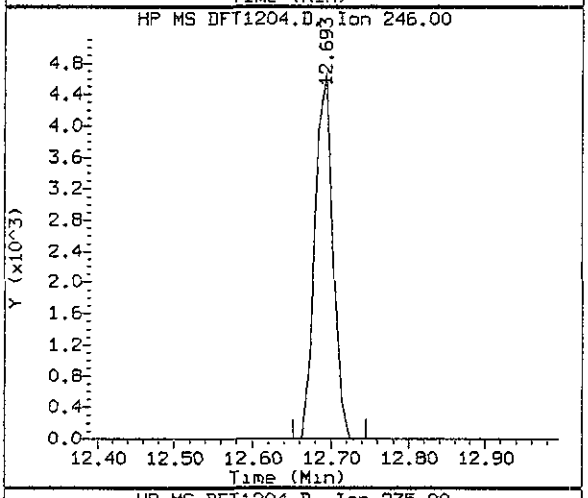


4,4'-DDD
========
 Exp. RT = 13.491
Found RT = 13.460

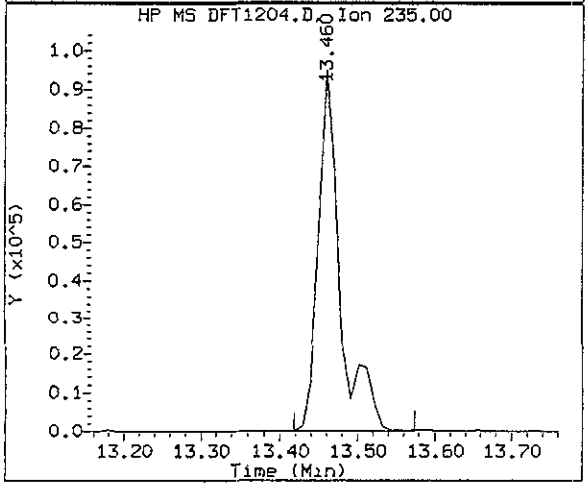
Mass	Area	Ratio
235	190198	100.00
237	120062	63.13
165	118067	62.08



Compound: 4,4'-DDT
Quant Mass: 235
RT: 14.144
Area: 1852497



Compound: 4,4'-DDE
Quant Mass: 246
RT: 12.693
Area: 7573



Compound: 4,4'-DDD
Quant Mass: 235
RT: 13.460
Area: 190198

DDT DEGRADATION BREAKDOWN ANALYSIS SUMMARY

Compound	Response	%Breakdown	Max Allowed	Test
4,4-DDD + DDE	197771	9.6	20.5	PASS

TestAmerica West Sacramento

Data file : \\SV5\C\chem\sv5.i\120410.B\DFT1204.D
Lab Smp Id: DFTPP 50ug/ml
Inj Date : 04-DEC-2010 10:15
Operator : KT Inst ID: sv5.i
Smp Info : DFTPP 50ug/ml;
Misc Info : 50ul DFTPP 10MSSV0129
Comment :
Method : \\SV5\C\chem\sv5.i\120410.B\DFTPP.M
Meth Date : 06-Dec-2010 17:52 onishim Quant Type: ISTD
Cal Date : Cal File:
Als bottle: 96 QC Sample: DFTPP
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14 Sample Matrix: None

					CONCENTRATIONS ON-COL	FINAL		
RT	EXP RT	REL RT	MASS	RESPONSE	(ug/L)	(ug/L)	TARGET RANGE	RATIO
1 dftpp						CAS #: 5074-71-5		
10.102	10.714	(0.000)	198	602112			0.00- 100.00	91.18
10.102	10.714	(0.000)	51	242048			30.00- 60.00	40.20
10.102	10.714	(0.000)	68	4023			0.00- 2.00	1.72
10.102	10.714	(0.000)	69	233472			0.00- 0.00	38.78
10.102	10.714	(0.000)	70	965			0.00- 2.00	0.41
10.102	10.714	(0.000)	127	334016			40.00- 60.00	55.47
10.102	10.714	(0.000)	197	0	0.0	0.0	0.00- 1.00	0.00
10.102	10.714	(0.000)	199	41056			5.00- 9.00	6.82
10.102	10.714	(0.000)	275	145792			10.00- 30.00	24.21
10.102	10.714	(0.000)	365	20320			1.00- 0.00	3.37
10.102	10.714	(0.000)	441	96904			0.01- 99.99	77.40
10.102	10.714	(0.000)	442	660352			40.00- 0.00	109.67
10.102	10.714	(0.000)	443	125192			17.00- 23.00	18.96

12/7/10

Data File: \\SV5\C\chem\sv5.i\120410.B\DFT1204.D

Date : 04-DEC-2010 10:15

Client ID:

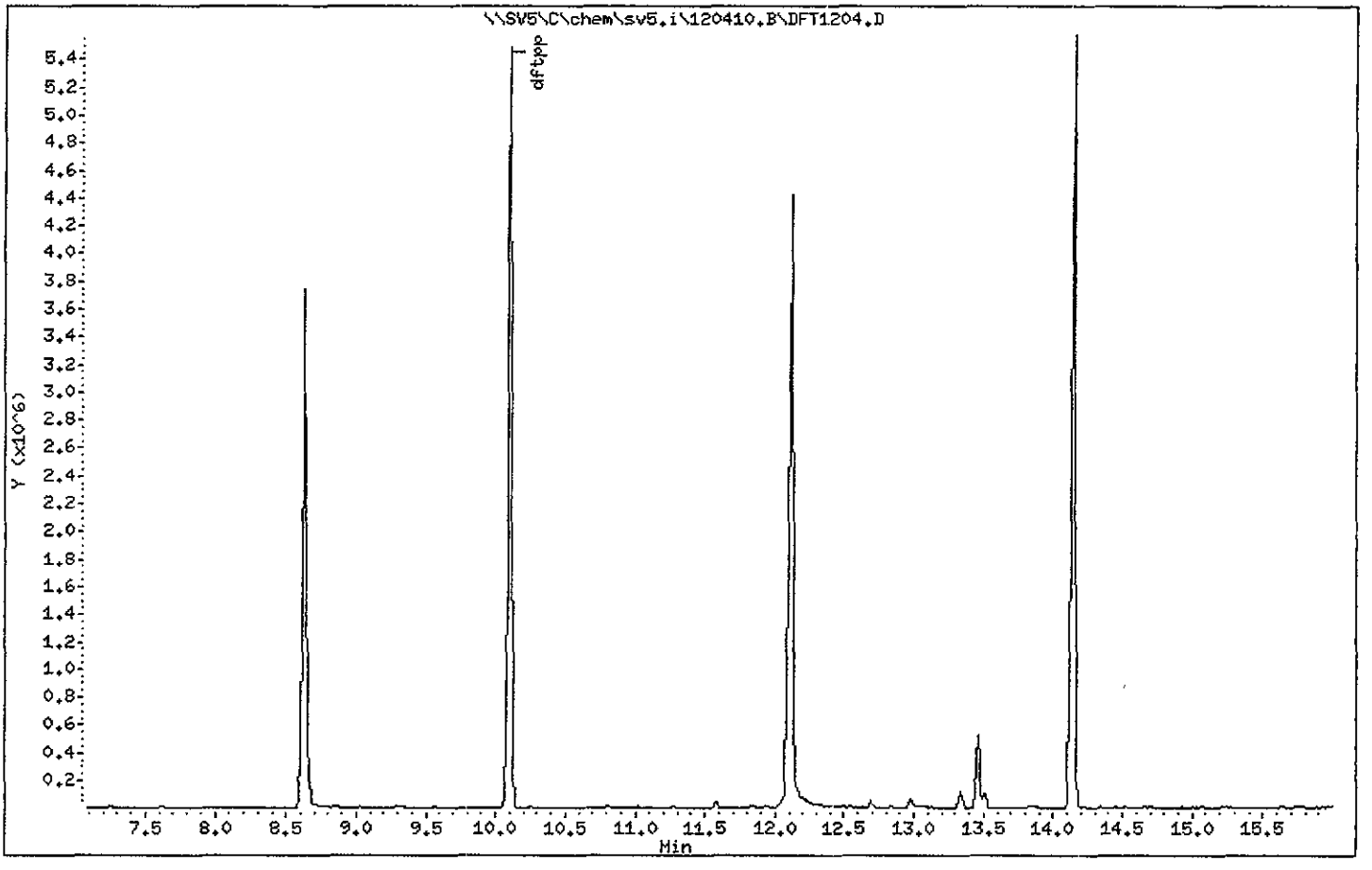
Instrument: sv5.i

Sample Info: DFTPP 50ug/ml;

Operator: KT

Column phase:

Column diameter: 2.00



Data File: \\SV5\C\chem\sv5.i\120410.B\DFT1204.D

Date : 04-DEC-2010 10:15

Client ID:

Instrument: sv5.i

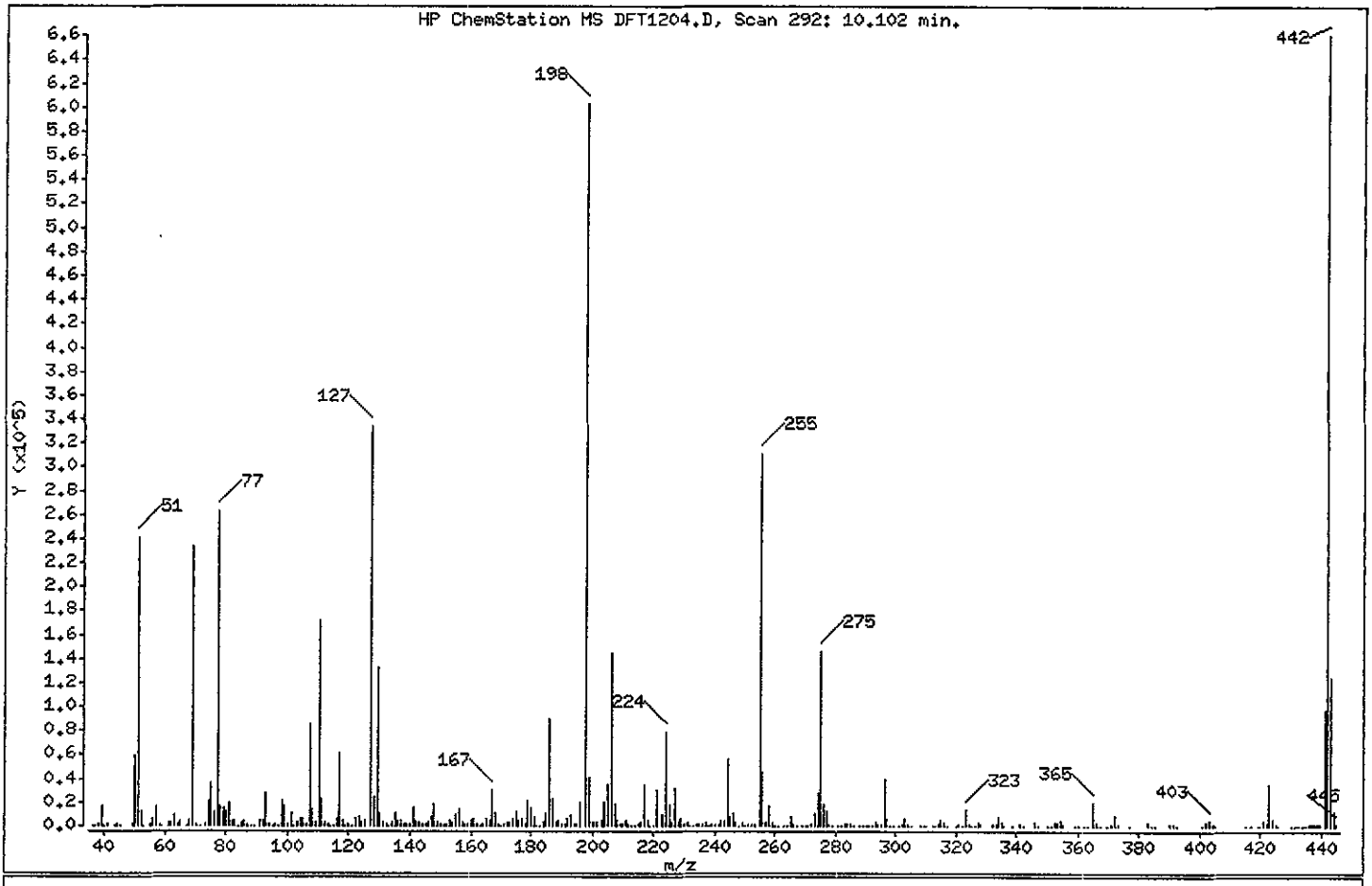
Sample Info: DFTPP 50ug/ml;

Operator: KT

Column phase:

Column diameter: 2.00

1 dftpp



m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
198	Base Peak, 100% relative abundance	100.00
51	30.00 - 60.00% of mass 198	40.20
68	Less than 2.00% of mass 69	0.67 (1.72)
69	Mass 69 relative abundance	38.78
70	Less than 2.00% of mass 69	0.16 (0.41)
127	40.00 - 60.00% of mass 198	55.47
197	Less than 1.00% of mass 198	0.00
199	5.00 - 9.00% of mass 198	6.82
275	10.00 - 30.00% of mass 198	24.21
365	Greater than 1.00% of mass 198	3.37
441	Present, but less than mass 443	16.09
442	Greater than 40.00% of mass 198	109.67
443	17.00 - 23.00% of mass 442	20.79 (18.96)

Data File: \\SV5\C\chem\sv5.i\120410.B\DFT1204.D

Date : 04-DEC-2010 10:15

Client ID:

Instrument: sv5.i

Sample Info: DFTPP 50ug/ml:

Operator: KT

Column phase:

Column diameter: 2.00

Data File: DFT1204.D

Spectrum: HP ChemStation MS DFT1204.D, Scan 292: 10.102 min.

Location of Maximum: 442.00

Number of points: 339

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	282	131.00	2299	217.00	35008	312.20	230
37.00	643	132.10	878	218.00	4584	312.90	363
38.00	2051	133.10	594	219.00	478	314.10	1902
39.10	16247	134.00	2704	219.70	438	315.00	3980
40.00	603	135.00	10782	221.00	31120	316.10	2632
41.00	769	136.00	4073	223.10	8914	317.10	504
43.20	392	137.10	4366	224.10	78848	319.90	314
44.10	1875	138.10	1107	225.10	17672	321.00	1385
45.00	506	139.00	794	226.00	2319	322.10	605
49.10	1928	140.00	1381	227.00	32536	323.10	14145
50.00	59600	141.00	15300	228.00	3654	324.10	2152
51.10	242048	142.00	4763	229.00	6308	325.00	292
52.10	11898	143.00	3119	230.00	1058	326.10	456
53.00	429	144.00	820	231.00	3128	327.00	2717
55.10	1504	145.10	1046	232.10	707	328.00	1253
56.00	6801	146.00	3121	233.10	706	332.00	950
57.00	16760	147.00	8217	234.00	1965	333.00	1578
58.00	810	148.00	17856	235.00	2086	334.00	8020
59.00	456	149.00	3070	236.00	2025	335.00	2458
61.00	2995	150.00	961	237.00	2310	335.80	244
62.00	2955	151.10	2143	238.10	219	338.90	273
63.00	9626	151.60	1146	239.10	1084	341.00	1467
64.00	1222	152.10	1063	240.00	944	342.10	549
65.00	4996	153.00	4760	241.10	2197	346.00	3223
67.10	580	154.00	3725	242.00	4221	347.00	590
68.10	4023	155.10	8778	243.10	4413	349.90	237
69.00	233472	156.10	13502	244.10	56384	351.00	499
70.00	965	157.10	2299	245.10	7252	352.10	3540
71.20	345	158.00	2175	246.00	11214	353.00	3314
73.00	1705	159.00	1919	247.00	2453	354.00	5039
74.10	21304	160.00	4173	248.00	696	355.00	770
75.00	36288	161.00	6612	249.00	2312	359.00	639
76.10	11418	162.00	1584	250.00	753	359.90	247
77.10	263040	163.00	544	251.10	1024	361.70	243
78.10	16904	164.00	791	252.20	800	362.10	221

Date : 04-DEC-2010 10:15

Client ID: Instrument: sv5.i

Sample Info: DFTPP 50ug/ml;

Operator: KT

Column phase: Column diameter: 2.00

Data File: DFT1204.D
Spectrum: HP ChemStation MS DFT1204.D, Scan 292: 10.102 min.
Location of Maximum: 442.00
Number of points: 339

m/z	Y	m/z	Y	m/z	Y	m/z	Y
79.00	14886	165.00	5326	253.10	2145	363.10	258
80.00	11539	166.10	4123	255.00	311232	365.00	20320
81.00	18976	167.00	30456	256.00	45088	366.00	2478
82.10	3990	168.00	10141	257.00	3426	367.00	219
83.00	4602	169.00	1942	258.00	17328	369.90	363
84.10	492	170.10	678	259.00	2798	371.00	1013
85.10	3609	171.10	1130	260.00	379	372.00	8795
86.00	4578	172.00	2700	260.90	614	373.10	2196
86.90	2167	173.00	3318	263.00	275	377.10	250
88.10	740	174.00	5695	263.90	367	383.00	2334
89.10	390	175.10	11501	265.00	7260	384.00	562
91.10	3952	176.00	4423	266.00	1136	385.10	343
92.00	4634	177.00	5885	267.00	285	390.00	1301
93.00	27936	178.00	1773	269.00	228	391.00	1006
94.10	1829	179.00	21016	270.00	307	392.10	493
95.10	323	180.00	14497	270.90	626	400.90	581
96.00	1135	181.00	7240	272.10	1148	402.00	3325
97.10	651	182.00	718	273.00	10518	403.00	4794
98.00	20656	182.90	645	274.00	27240	404.00	1537
99.00	16022	184.10	2499	275.00	145792	405.00	386
100.00	1745	185.00	11060	276.00	18888	415.00	273
101.00	11080	186.00	89032	277.00	11943	417.00	224
102.10	738	187.10	23264	278.00	1987	419.10	233
103.00	3020	188.10	2652	279.00	334	419.60	227
104.00	6559	189.00	4982	281.00	287	421.00	4715
105.00	5539	190.00	768	282.10	374	422.10	3771
106.00	2181	191.00	2137	283.00	1339	423.00	35624
107.00	84768	192.00	6389	284.00	983	424.00	6651
108.00	13676	193.00	8395	285.10	2122	424.90	957
109.10	2829	194.10	2210	286.10	379	425.60	303
110.00	171456	195.10	1405	287.90	227	429.90	366
111.00	23344	196.10	20304	289.00	514	431.30	390
112.00	3077	198.00	602112	290.10	359	431.70	340
113.00	1269	199.00	41056	291.00	300	432.40	366
114.00	388	200.00	3296	292.00	493	433.20	577

Date : 04-DEC-2010 10:15

Client ID: Instrument: sv5.i

Sample Info: DFTPP 50ug/ml:

Operator: KT

Column phase: Column diameter: 2.00

Data File: DFT1204.D
Spectrum: HP ChemStation MS DFT1204.D, Scan 292: 10.102 min.
Location of Maximum: 442.00
Number of points: 339

m/z	Y	m/z	Y	m/z	Y	m/z	Y
114.90	452	201.60	2731	293.00	2809	433.80	408
116.00	5524	203.10	3879	294.00	949	434.50	281
117.00	60416	204.00	19928	295.10	1820	436.00	1026
118.10	4647	205.00	34544	296.00	39752	436.60	916
118.80	450	206.10	144320	297.00	6569	437.10	913
120.00	1298	207.10	18832	298.10	323	437.50	770
121.10	691	208.00	5234	299.10	284	438.20	1061
122.00	6433	209.00	1417	301.00	284	439.10	1171
123.10	8001	210.00	2055	302.10	896	439.70	1680
124.00	3774	211.10	5005	303.00	5368	441.00	96904
125.00	3919	211.70	1342	304.00	1186	442.00	660352
127.00	334016	213.00	234	305.10	218	443.00	125192
128.00	23536	214.20	462	308.00	716	444.00	11472
129.00	131520	215.00	1617	309.10	352	445.00	808
130.00	9935	216.00	3149	310.00	450		

TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\120410.B\S120410.D
Lab Smp Id: MANG51AA G0K30000-3 Client Smp ID: 0334389
Inj Date : 04-DEC-2010 14:41
Operator : KT Inst ID: sv5.i
Smp Info : MANG51AA G0K30000-389B;0;;;1000;;1000;5
Misc Info : 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\120410.B\8270f.m
Meth Date : 06-Dec-2010 09:40 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 8
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: S11JZHCB.SUB
Target Version: 4.14
Processing Host: SACP307UM

Concentration Formula: Amt * DF * Uf * Vt/(Vo * Vi) * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	1000.000	Volume of final extract (uL)
Vo	1000.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected (uL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (NG)	CONCENTRATIONS FINAL (ug/L)	
*	1 1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	111249	40.0000		(Q)
*	2 Naphthalene-d8	136	4.952	4.952	(1.000)	462027	40.0000		
*	3 Acenaphthene-d10	164	7.045	7.045	(1.000)	251894	40.0000		
*	4 Phenanthrene-d10	188	8.900	8.900	(1.000)	419551	40.0000		
*	5 Chrysene-d12	240	13.180	13.190	(1.000)	413584	40.0000		
*	6 Perylene-d12	264	15.543	15.543	(1.000)	426529	40.0000		
$	7 2-Fluorophenol	112	2.340	2.340	(0.661)	247104	63.0157	63.02	
$	8 Phenol-d5	99	3.231	3.231	(0.912)	347094	70.3901	70.39	
$	10 1,2-Dichlorobenzene-d4	152	3.739	3.739	(1.056)	79651	29.0712	29.07	(qR)
$	11 Nitrobenzene-d5	82	4.164	4.174	(0.841)	135320	34.5795	34.58	
$	12 2-Fluorobiphenyl	172	6.257	6.257	(0.888)	295662	36.4372	36.44	
$	13 2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	110415	100.875	100.9	
$	14 Terphenyl-d14	244	11.460	11.460	(0.869)	376860	46.2603	46.26	
	108 Hexachlorobenzene	284	Compound Not Detected.						

12/6/10

QC Flag Legend

Q - Qualifier signal failed the ratio test.
R - Spike/Surrogate failed recovery limits.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: S120410.D
Lab Smp Id: MANG51AA G0K30000-3
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\SV5\C\chem\sv5.i\120410.B\8270F.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M

Calibration Date: 04-DEC-2010
Calibration Time: 10:36
Client Smp ID: 0334389
Level: LOW
Sample Type: AIR

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	111249	-9.28
2 Naphthalene-d8	530514	265257	1061028	462027	-12.91
3 Acenaphthene-d10	282538	141269	565076	251894	-10.85
4 Phenanthrene-d10	462722	231361	925444	419551	-9.33
5 Chrysene-d12	435850	217925	871700	413584	-5.11
6 Perylene-d12	422284	211142	844568	426529	1.01

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.54	3.04	4.04	3.54	-0.00
2 Naphthalene-d8	4.95	4.45	5.45	4.95	0.00
3 Acenaphthene-d10	7.05	6.55	7.55	7.05	0.00
4 Phenanthrene-d10	8.90	8.40	9.40	8.90	0.00
5 Chrysene-d12	13.19	12.69	13.69	13.18	-0.08
6 Perylene-d12	15.54	15.04	16.04	15.54	-0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

TestAmerica West Sacramento

RECOVERY REPORT

Client Name: Client SDG: 090498
Sample Matrix: GAS Fraction: SV
Lab Smp Id: MANG51AA G0K30000-3 Client Smp ID: 0334389
Level: LOW Operator: KT
Data Type: MS DATA SampleType: SAMPLE
SpikeList File: Quant Type: ISTD
Sublist File: S11JZHCB.SUB
Method File: \\sv5\c\chem\sv5.i\120410.B\8270f.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M

SURROGATE COMPOUND	CONC ADDED ug/L	CONC RECOVERED ug/L	% RECOVERED	LIMITS
$ 7 2-Fluorophenol	100.0	63.02	63.02	41-105
$ 8 Phenol-d5	100.0	70.39	70.39	43-122
$ 10 1,2-Dichlorobenzen	50.00	29.07	58.14*	60-120
$ 11 Nitrobenzene-d5	50.00	34.58	69.16	46-118
$ 12 2-Fluorobiphenyl	50.00	36.44	72.87	58-105
$ 13 2,4,6-Tribromophen	100.0	100.9	100.87	61-118
$ 14 Terphenyl-d14	50.00	46.26	92.52	69-110

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\120410.B\S120410.D
Lab Smp Id: MANG51AA G0K30000-3 Client Smp ID: 0334389
Inj Date : 04-DEC-2010 14:41
Operator : KT Inst ID: sv5.i
Smp Info : MANG51AA G0K30000-389B;0;;;1000;;1000;5
Misc Info : 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\120410.B\8270F.m
Meth Date : 06-Dec-2010 09:27 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 8
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: S11JZHCB.SUB
Target Version: 4.14
Processing Host: SV5

Concentration Formula: Amt * DF * Uf * Vt/(Vo * Vi) * CpndVariable

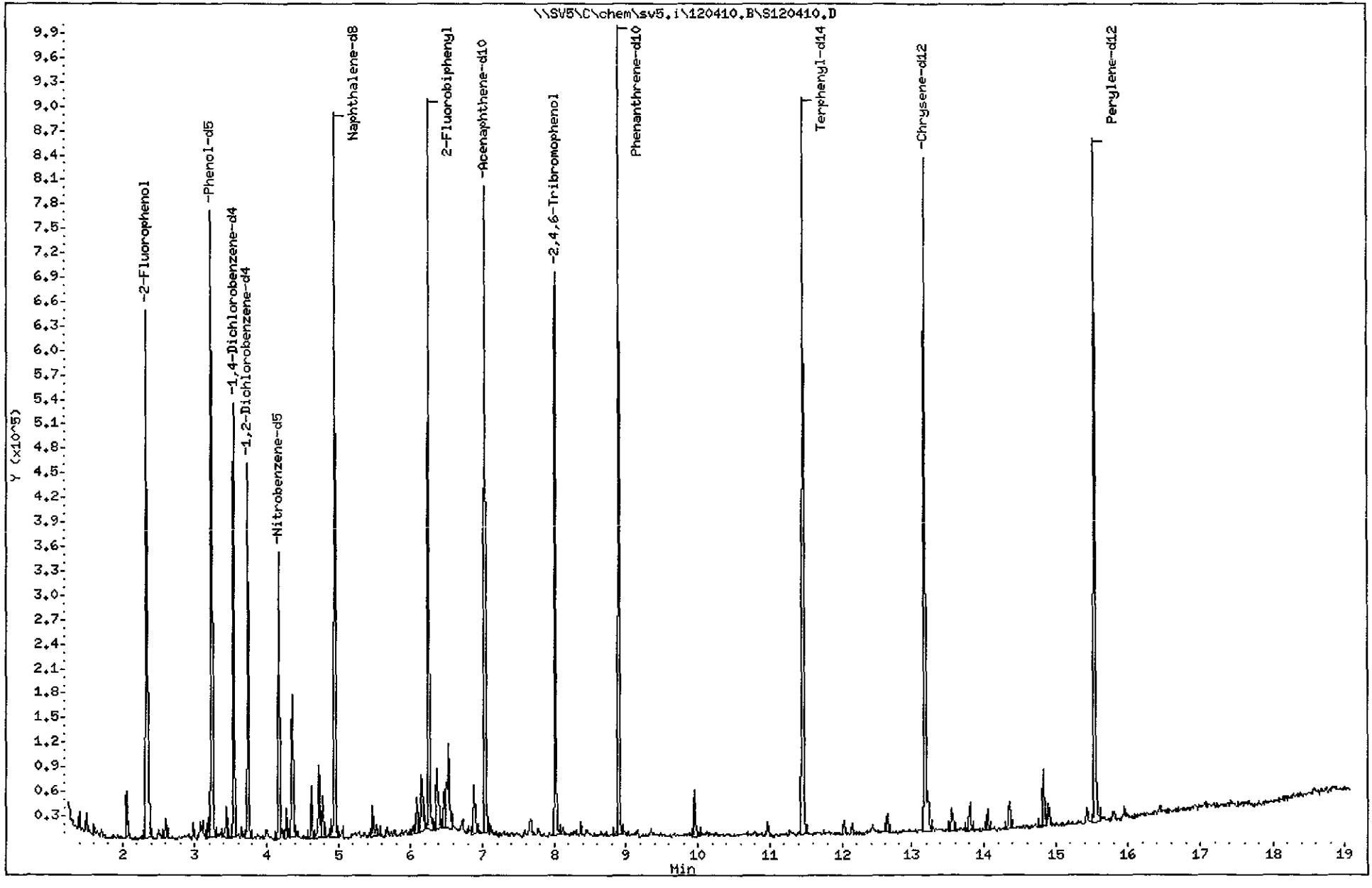
Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	1000.000	Volume of final extract (uL)
Vo	1000.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected (uL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (NG)	CONCENTRATIONS FINAL (ug/L)
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	111249	40.0000	(Q)
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	462027	40.0000	
* 3	Acenaphthene-d10	164	7.045	7.045	(1.000)	251894	40.0000	
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	419551	40.0000	
* 5	Chrysene-d12	240	13.180	13.190	(1.000)	413584	40.0000	
* 6	Perylene-d12	264	15.543	15.543	(1.000)	426529	40.0000	
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	247104	63.0157	63.02
$ 8	Phenol-d5	99	3.231	3.231	(0.912)	347094	70.3901	70.39
$ 10	1,2-Dichlorobenzene-d4	152	3.739	3.739	(1.056)	79651	29.0712	29.07(qR)
$ 11	Nitrobenzene-d5	82	4.164	4.174	(0.841)	135320	34.5795	34.58
$ 12	2-Fluorobiphenyl	172	6.257	6.257	(0.888)	295662	36.4372	36.44
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	110415	100.875	100.9
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	376860	46.2603	46.26
108	Hexachlorobenzene	284	Compound Not Detected.					

QC Flag Legend

Q - Qualifier signal failed the ratio test.
R - Spike/Surrogate failed recovery limits.
q - Qualifier signal exceeded ratio warning limit.

Data File: \\SV5\C\chem\sv5.i\120410.B\S120410.D
Date : 04-DEC-2010 14:41
Client ID: 0334389
Instrument: sv5.i
Sample Info: MANC51AA G0K30000-389B;0;;;1000;;1000;5
Volume Injected (uL): 1.0
Operator: KT
Column phase:
Column diameter: 2.00



TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\120410.B\S120411.D
Lab Smp Id: MANG51AC G0K30000-3
Inj Date : 04-DEC-2010 15:05
Operator : KT Inst ID: sv5.i
Smp Info : MANG51AC G0K30000-389C;3;LCS;;1000;;1000;2
Misc Info : 0;AIR;0;S11JZHCB.SUB;S11JZHCB.SPK;1;;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\120410.B\8270f.m
Meth Date : 06-Dec-2010 09:40 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 9 QC Sample: LCS
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: S11JZHCB.SUB
Target Version: 4.14
Processing Host: SACP307UM

Concentration Formula: Amt * DF * Uf * Vt/(Vo * Vi) * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	1000.000	Volume of final extract (uL)
Vo	1000.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected (uL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (NG)	CONCENTRATIONS FINAL (ug/L)
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	138012	40.0000	
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	611787	40.0000	
* 3	Acenaphthene-d10	164	7.045	7.045	(1.000)	327932	40.0000	
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	544079	40.0000	
* 5	Chrysene-d12	240	13.190	13.190	(1.000)	533027	40.0000	
* 6	Perylene-d12	264	15.543	15.543	(1.000)	541754	40.0000	
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	374216	76.9256	76.92
$ 8	Phenol-d5	99	3.242	3.231	(0.915)	524160	85.6856	85.68
$ 10	1,2-Dichlorobenzene-d4	152	Compound Not Detected.					
$ 11	Nitrobenzene-d5	82	4.174	4.174	(0.843)	221480	42.7423	42.74
$ 12	2-Fluorobiphenyl	172	6.257	6.257	(0.888)	473122	44.7875	44.79
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	155191	108.907	108.9
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	476301	45.3654	45.36
108	Hexachlorobenzene	284	8.496	8.496	(0.955)	290305	97.8735	97.87

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TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: S120411.D
Lab Smp Id: MANG51AC G0K30000-3
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\SV5\C\chem\sv5.i\120410.B\8270F.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;S11JZHCB.SPK;1;;8270F.M

Calibration Date: 04-DEC-2010
Calibration Time: 10:36
Level: LOW
Sample Type: AIR

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	138012	12.55
2 Naphthalene-d8	530514	265257	1061028	611787	15.32
3 Acenaphthene-d10	282538	141269	565076	327932	16.07
4 Phenanthrene-d10	462722	231361	925444	544079	17.58
5 Chrysene-d12	435850	217925	871700	533027	22.30
6 Perylene-d12	422284	211142	844568	541754	28.29

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.54	3.04	4.04	3.54	0.00
2 Naphthalene-d8	4.95	4.45	5.45	4.95	0.00
3 Acenaphthene-d10	7.05	6.55	7.55	7.05	0.00
4 Phenanthrene-d10	8.90	8.40	9.40	8.90	0.00
5 Chrysene-d12	13.19	12.69	13.69	13.19	0.00
6 Perylene-d12	15.54	15.04	16.04	15.54	0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

TestAmerica West Sacramento

RECOVERY REPORT

Client Name: Client SDG: 090498
Sample Matrix: GAS Fraction: SV
Lab Smp Id: MANG51AC G0K30000-3
Level: LOW Operator: KT
Data Type: MS DATA SampleType: LCS
SpikeList File: S11JZHCB.SPK Quant Type: ISTD
Sublist File: S11JZHCB.SUB
Method File: \\sv5\c\chem\sv5.i\120410.B\8270f.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;S11JZHCB.SPK;1;;8270F.M

SPIKE COMPOUND	CONC ADDED ug/L	CONC RECOVERED ug/L	% RECOVERED	LIMITS
108 Hexachlorobenzene	100.0	97.87	97.87	70-100

SURROGATE COMPOUND	CONC ADDED ug/L	CONC RECOVERED ug/L	% RECOVERED	LIMITS
$ 7 2-Fluorophenol	100.0	76.92	76.93	41-105
$ 8 Phenol-d5	100.0	85.68	85.69	43-122
$ 10 1,2-Dichlorobenze	50.00	0.0000	*	60-120
$ 11 Nitrobenzene-d5	50.00	42.74	85.48	46-118
$ 12 2-Fluorobiphenyl	50.00	44.79	89.58	58-105
$ 13 2,4,6-Tribromophen	100.0	108.9	108.91	61-118
$ 14 Terphenyl-d14	50.00	45.36	90.73	69-110

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\120410.B\S120411.D
Lab Smp Id: MANG51AC G0K30000-3
Inj Date : 04-DEC-2010 15:05
Operator : KT Inst ID: sv5.i
Smp Info : MANG51AC G0K30000-389C;3;LCS;;1000;;1000;2
Misc Info : 0;AIR;0;S11JZHCB.SUB;S11JZHCB.SPK;1;;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\120410.B\8270F.m
Meth Date : 06-Dec-2010 09:27 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 9 QC Sample: LCS
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: S11JZHCB.SUB
Target Version: 4.14
Processing Host: SV5

Concentration Formula: Amt * DF * Uf * Vt/(Vo * Vi) * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	1000.000	Volume of final extract (uL)
Vo	1000.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected (uL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (NG)	FINAL (ug/L)
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	138012	40.0000	
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	611787	40.0000	
* 3	Acenaphthene-d10	164	7.045	7.045	(1.000)	327932	40.0000	
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	544079	40.0000	
* 5	Chrysene-d12	240	13.190	13.190	(1.000)	533027	40.0000	
* 6	Perylene-d12	264	15.543	15.543	(1.000)	541754	40.0000	
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	374216	76.9256	76.92
$ 8	Phenol-d5	99	3.242	3.231	(0.915)	524160	85.6856	85.68
$ 10	1,2-Dichlorobenzene-d4	152	3.542	3.739	(1.000)	138011	40.6036	40.60(qR)
$ 11	Nitrobenzene-d5	82	4.174	4.174	(0.843)	221480	42.7423	42.74
$ 12	2-Fluorobiphenyl	172	6.257	6.257	(0.888)	473122	44.7875	44.79
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	155191	108.907	108.9
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	476301	45.3654	45.36
108	Hexachlorobenzene	284	8.496	8.496	(0.955)	290305	97.8735	97.87

QC Flag Legend

R - Spike/Surrogate failed recovery limits.
q - Qualifier signal exceeded ratio warning limit.

Data File: \\SV5\C\chem\sv5.i\120410.B\S120411.D
Date : 04-DEC-2010 15:05
Client ID:
Instrument: sv5.i
Sample Info: MANG51AC G0K30000-389C;3;LCS;;1000;;1000;2
Volume Injected (uL): 1.0
Operator: KT
Column phase:
Column diameter: 2.00

Data File: \\SV5\C\chem\sv5.i\120410.B\S120411.D

Date : 04-DEC-2010 15:05

Client ID:

Instrument: sv5.i

Sample Info: MANG51AC G0K30000-389C;3;LCS;;1000;;1000;2

Volume Injected (uL): 1.0

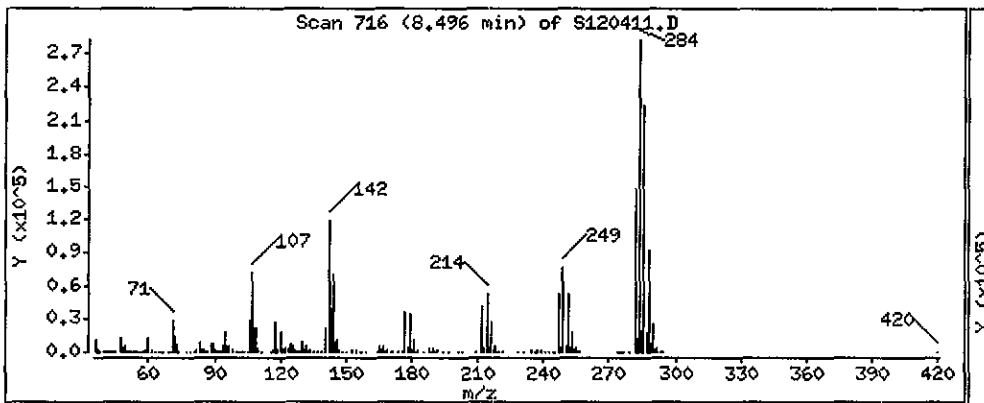
Operator: KT

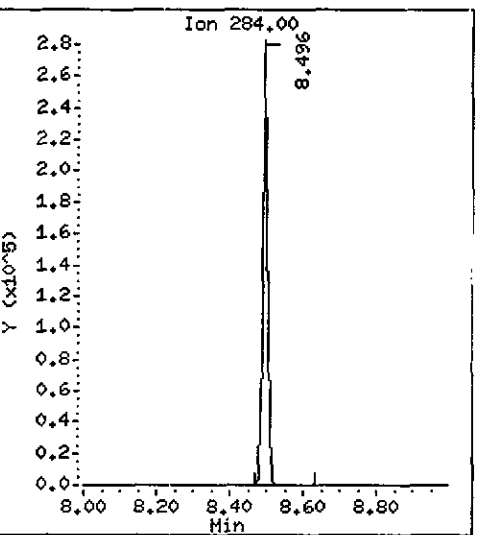
Column phase:

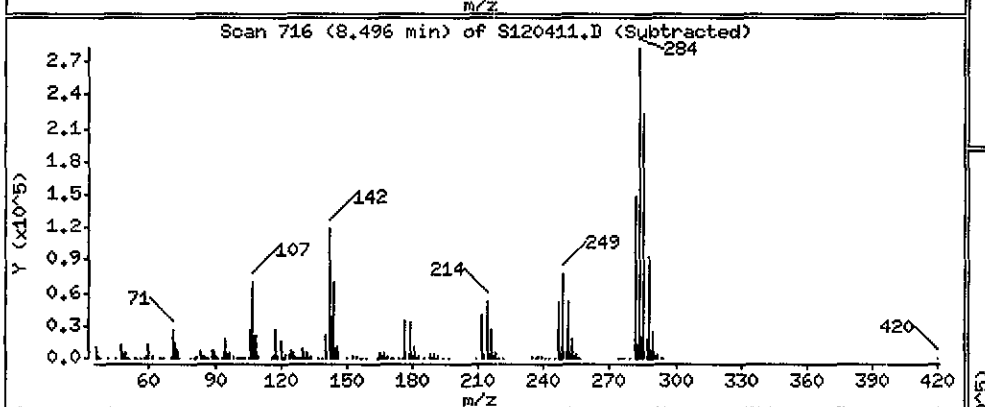
Column diameter: 2.00

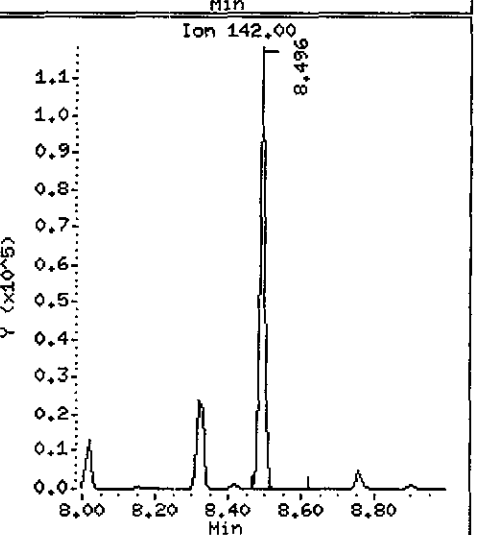
108 Hexachlorobenzene

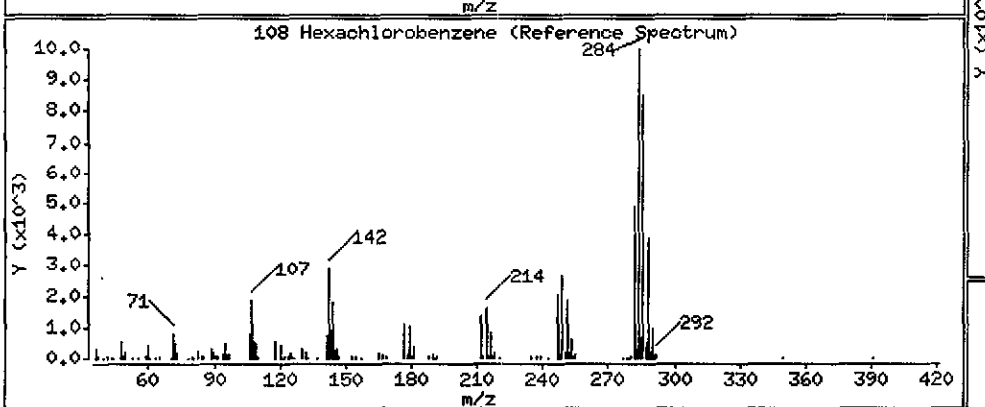
Concentration: 97.87 ug/L

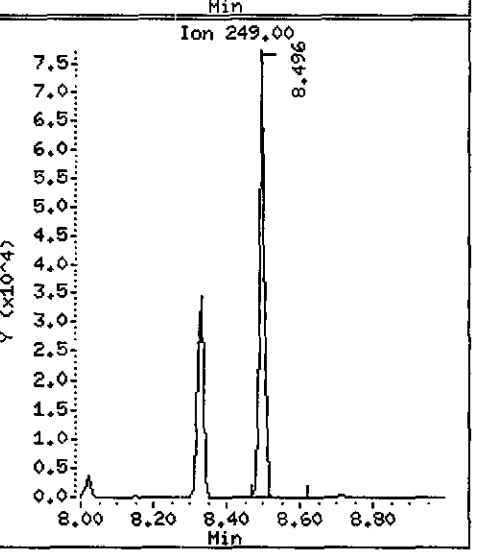


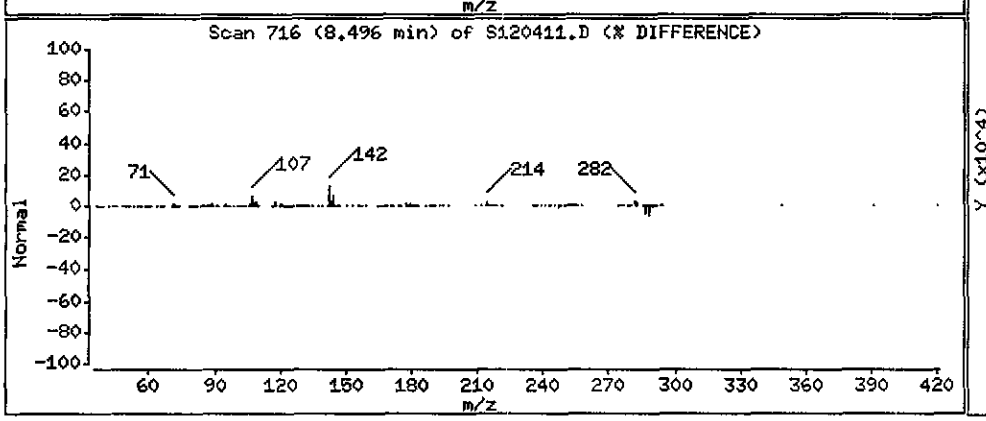












TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\120410.B\S120412.D
Lab Smp Id: MANG51AD G0K30000-3
Inj Date : 04-DEC-2010 15:30
Operator : KT Inst ID: sv5.i
Smp Info : MANG51AD G0K30000-389L;3;LCSD;;1000;;1000;2
Misc Info : 0;AIR;0;S11JZHCB.SUB;S11JZHCB.SPK;1;;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\120410.B\8270f.m
Meth Date : 06-Dec-2010 09:40 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 10 QC Sample: LCSD
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: S11JZHCB.SUB
Target Version: 4.14
Processing Host: SACP307UM

Concentration Formula: Amt * DF * Uf * Vt/(Vo * Vi) * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	1000.000	Volume of final extract (uL)
Vo	1000.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected (uL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (NG)	CONCENTRATIONS FINAL (ug/L)
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	102167	40.0000	
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	426724	40.0000	
* 3	Acenaphthene-d10	164	7.045	7.045	(1.000)	224321	40.0000	
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	371643	40.0000	
* 5	Chrysene-d12	240	13.190	13.190	(1.000)	361730	40.0000	
* 6	Perylene-d12	264	15.543	15.543	(1.000)	380694	40.0000	
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	265284	73.6658	73.66
$ 8	Phenol-d5	99	3.232	3.231	(0.912)	359457	79.3775	79.38
$ 10	1,2-Dichlorobenzene-d4	152	Compound Not Detected.					
$ 11	Nitrobenzene-d5	82	4.175	4.174	(0.843)	151463	41.9067	41.91
$ 12	2-Fluorobiphenyl	172	6.258	6.257	(0.888)	329965	45.6631	45.66
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	109539	112.376	112.4
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	320482	44.9792	44.98
108	Hexachlorobenzene	284	8.496	8.496	(0.955)	196593	97.0319	97.03

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TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: S120412.D
Lab Smp Id: MANG51AD G0K30000-3
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\SV5\C\chem\sv5.i\120410.B\8270F.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;S11JZHCB.SPK;1;;8270F.M

Calibration Date: 04-DEC-2010
Calibration Time: 10:36
Level: LOW
Sample Type: AIR

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	102167	-16.68
2 Naphthalene-d8	530514	265257	1061028	426724	-19.56
3 Acenaphthene-d10	282538	141269	565076	224321	-20.61
4 Phenanthrene-d10	462722	231361	925444	371643	-19.68
5 Chrysene-d12	435850	217925	871700	361730	-17.01
6 Perylene-d12	422284	211142	844568	380694	-9.85

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.54	3.04	4.04	3.54	0.00
2 Naphthalene-d8	4.95	4.45	5.45	4.95	0.00
3 Acenaphthene-d10	7.05	6.55	7.55	7.05	0.00
4 Phenanthrene-d10	8.90	8.40	9.40	8.90	0.00
5 Chrysene-d12	13.19	12.69	13.69	13.19	0.00
6 Perylene-d12	15.54	15.04	16.04	15.54	0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

TestAmerica West Sacramento

RECOVERY REPORT

Client Name: Client SDG: 090498
Sample Matrix: GAS Fraction: SV
Lab Smp Id: MANG51AD G0K30000-3
Level: LOW Operator: KT
Data Type: MS DATA SampleType: LCSD
SpikeList File: S11JZHCB.SPK Quant Type: ISTD
Sublist File: S11JZHCB.SUB
Method File: \\sv5\c\chem\sv5.i\120410.B\8270f.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;S11JZHCB.SPK;1;;8270F.M

SPIKE COMPOUND	CONC ADDED ug/L	CONC RECOVERED ug/L	% RECOVERED	LIMITS
108 Hexachlorobenzene	100.0	97.03	97.03	70-100

SURROGATE COMPOUND	CONC ADDED ug/L	CONC RECOVERED ug/L	% RECOVERED	LIMITS
$ 7 2-Fluorophenol	100.0	73.66	73.67	41-105
$ 8 Phenol-d5	100.0	79.38	79.38	43-122
$ 10 1,2-Dichlorobenze	50.00	0.0000	*	60-120
$ 11 Nitrobenzene-d5	50.00	41.91	83.81	46-118
$ 12 2-Fluorobiphenyl	50.00	45.66	91.33	58-105
$ 13 2,4,6-Tribromophen	100.0	112.4	112.38	61-118
$ 14 Terphenyl-d14	50.00	44.98	89.96	69-110

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\120410.B\S120412.D
Lab Smp Id: MANG51AD G0K30000-3
Inj Date : 04-DEC-2010 15:30
Operator : KT Inst ID: sv5.i
Smp Info : MANG51AD G0K30000-389L;3;LCSD;;1000;;1000;2
Misc Info : 0;AIR;0;S11JZHCB.SUB;S11JZHCB.SPK;1;;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\120410.B\8270F.m
Meth Date : 06-Dec-2010 09:27 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 10 QC Sample: LCSD
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: S11JZHCB.SUB
Target Version: 4.14
Processing Host: SV5

Concentration Formula: Amt * DF * Uf * Vt/(Vo * Vi) * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	1000.000	Volume of final extract (uL)
Vo	1000.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected (uL)
Cpnd Variable		Local Compound Variable

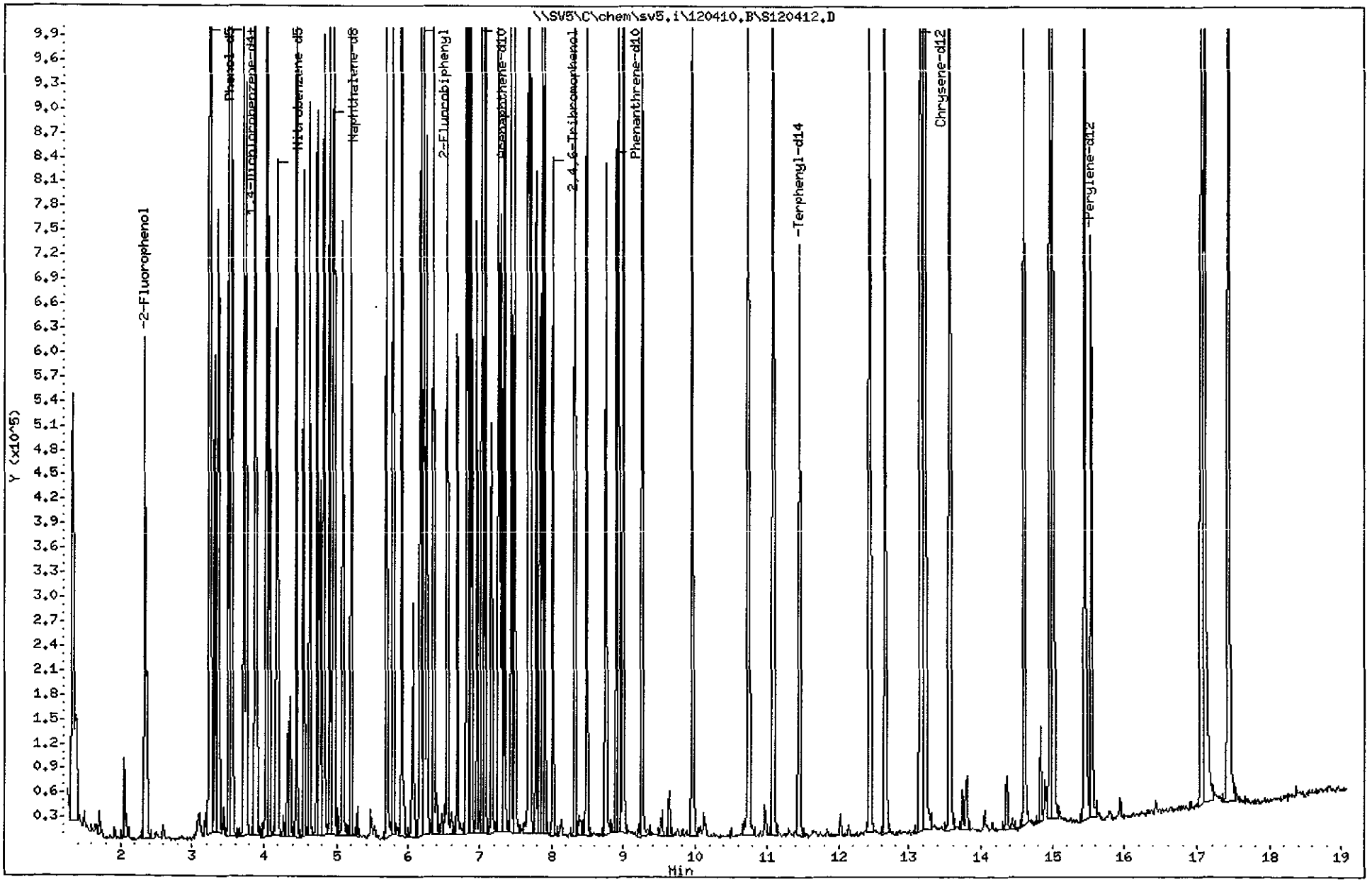
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (NG)	CONCENTRATIONS FINAL (ug/L)
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	102167	40.0000	
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	426724	40.0000	
* 3	Acenaphthene-d10	164	7.045	7.045	(1.000)	224321	40.0000	
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	371643	40.0000	
* 5	Chrysene-d12	240	13.190	13.190	(1.000)	361730	40.0000	
* 6	Perylene-d12	264	15.543	15.543	(1.000)	380694	40.0000	
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	265284	73.6658	73.66
$ 8	Phenol-d5	99	3.232	3.231	(0.912)	359457	79.3775	79.38
$ 10	1,2-Dichlorobenzene-d4	152	3.542	3.739	(1.000)	102167	40.6039	40.60(qR)
$ 11	Nitrobenzene-d5	82	4.175	4.174	(0.843)	151463	41.9067	41.91
$ 12	2-Fluorobiphenyl	172	6.258	6.257	(0.888)	329965	45.6631	45.66
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	109539	112.376	112.4
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	320482	44.9792	44.98
108	Hexachlorobenzene	284	8.496	8.496	(0.955)	196593	97.0319	97.03

QC Flag Legend

R - Spike/Surrogate failed recovery limits.
q - Qualifier signal exceeded ratio warning limit.

Data File: \\SV5\C\chem\sv5.i\120410.B\S120412.D
Date : 04-DEC-2010 15:30
Client ID:
Sample Info: MANG51AD G0K30000-389L;3;LCSD;;1000;;1000;2
Volume Injected (uL): 1.0
Column phase:

Instrument: sv5.i
Operator: KT
Column diameter: 2.00



Data File: \\SV5\C\chem\sv5.i\120410.B\S120412.D

Date : 04-DEC-2010 15:30

Client ID:

Instrument: sv5.i

Sample Info: MANG51AD G0K30000-389L;3;LCSD;;1000;;1000;2

Volume Injected (uL): 1.0

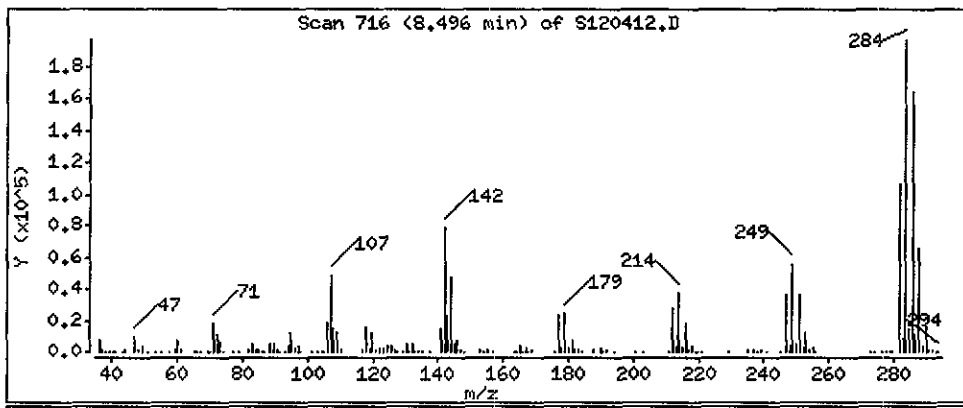
Operator: KT

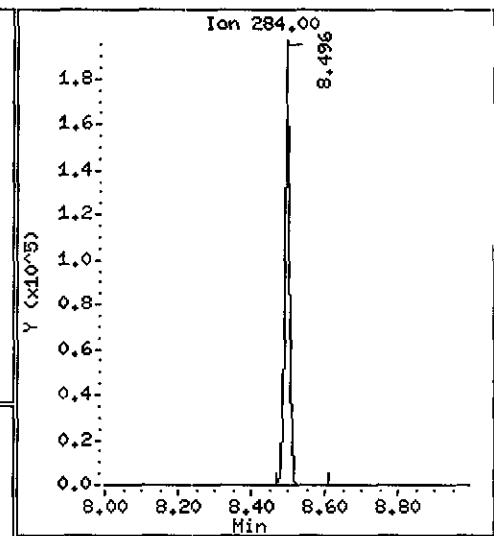
Column phase:

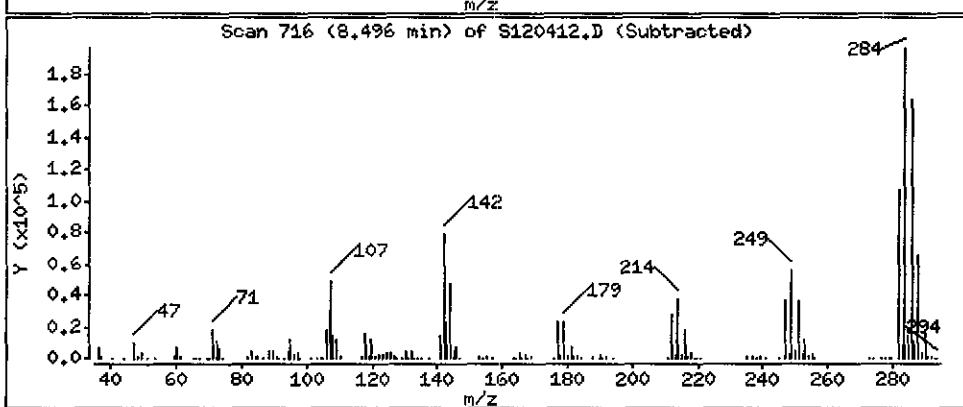
Column diameter: 2.00

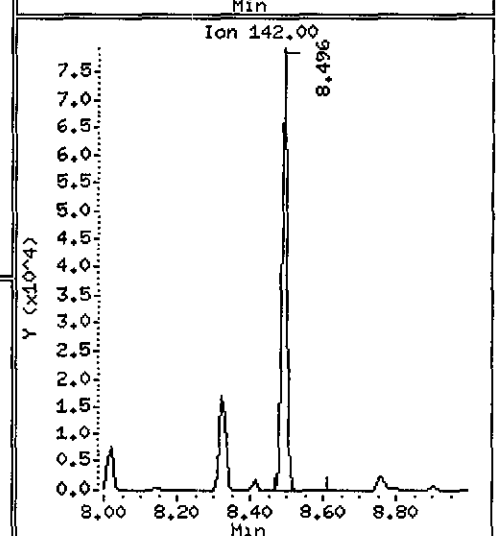
108 Hexachlorobenzene

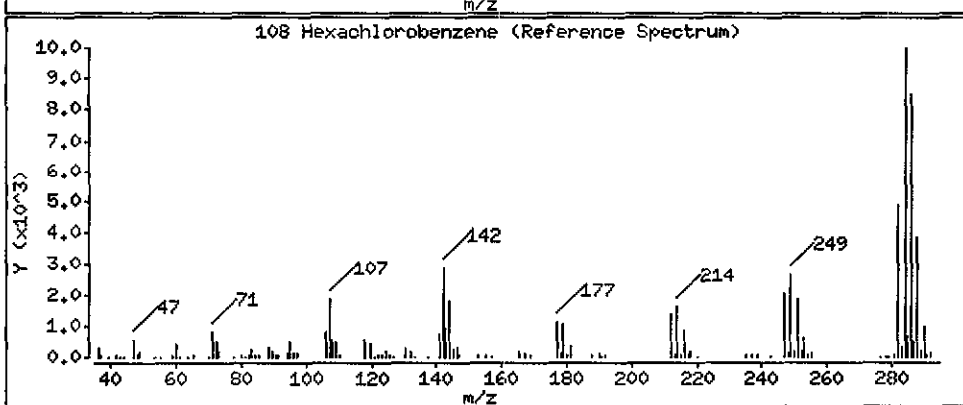
Concentration: 97.03 ug/L

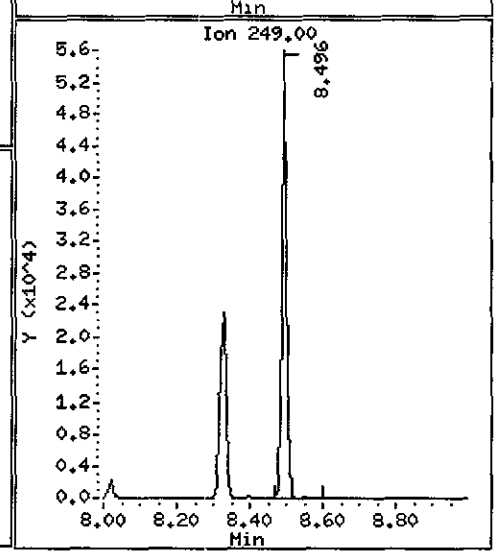


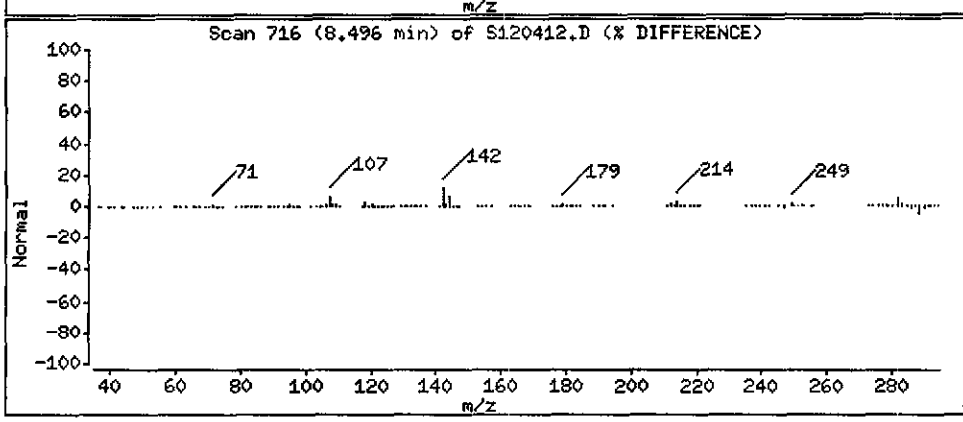












TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\120410.B\S120413.D
Lab Smp Id: MAMLW1AA G0K300434- Client Smp ID: 0334389
Inj Date : 04-DEC-2010 15:54
Operator : KT Inst ID: sv5.i
Smp Info : MAMLW1AA G0K300434-1;0;;;1000;;1000;5
Misc Info : 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\120410.B\8270f.m
Meth Date : 06-Dec-2010 09:40 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 11
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: S11JZHCB.SUB
Target Version: 4.14
Processing Host: SACP307UM

Concentration Formula: Amt * DF * Uf * Vt/(Vo * Vi) * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	1000.000	Volume of final extract (uL)
Vo	1000.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected (uL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (NG)	FINAL (ug/L)
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	109565	40.0000	(Q)
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	493823	40.0000	
* 3	Acenaphthene-d10	164	7.035	7.045	(1.000)	272801	40.0000	
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	465265	40.0000	
* 5	Chrysene-d12	240	13.180	13.190	(1.000)	455506	40.0000	
* 6	Perylene-d12	264	15.543	15.543	(1.000)	467119	40.0000	
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	282871	73.2457	73.24
$ 8	Phenol-d5	99	3.231	3.231	(0.912)	412940	85.0307	85.03
$ 10	1,2-Dichlorobenzene-d4	152	3.739	3.739	(1.056)	95342	35.3330	35.33(q)
$ 11	Nitrobenzene-d5	82	4.164	4.174	(0.841)	170096	40.6674	40.67
$ 12	2-Fluorobiphenyl	172	6.257	6.257	(0.890)	375233	42.6995	42.70
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.140)	126090	106.367	106.4
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	412314	45.9543	45.95
108	Hexachlorobenzene	284	Compound Not Detected.					

QC Flag Legend

Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

12/6/10

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: S120413.D
Lab Smp Id: MAMLW1AA G0K300434-
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\SV5\C\chem\sv5.i\120410.B\8270F.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M

Calibration Date: 04-DEC-2010
Calibration Time: 10:36
Client Smp ID: 0334389
Level: LOW
Sample Type: AIR

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	109565	-10.65
2 Naphthalene-d8	530514	265257	1061028	493823	-6.92
3 Acenaphthene-d10	282538	141269	565076	272801	-3.45
4 Phenanthrene-d10	462722	231361	925444	465265	0.55
5 Chrysene-d12	435850	217925	871700	455506	4.51
6 Perylene-d12	422284	211142	844568	467119	10.62

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.54	3.04	4.04	3.54	0.00
2 Naphthalene-d8	4.95	4.45	5.45	4.95	0.00
3 Acenaphthene-d10	7.05	6.55	7.55	7.04	-0.15
4 Phenanthrene-d10	8.90	8.40	9.40	8.90	0.00
5 Chrysene-d12	13.19	12.69	13.69	13.18	-0.08
6 Perylene-d12	15.54	15.04	16.04	15.54	0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

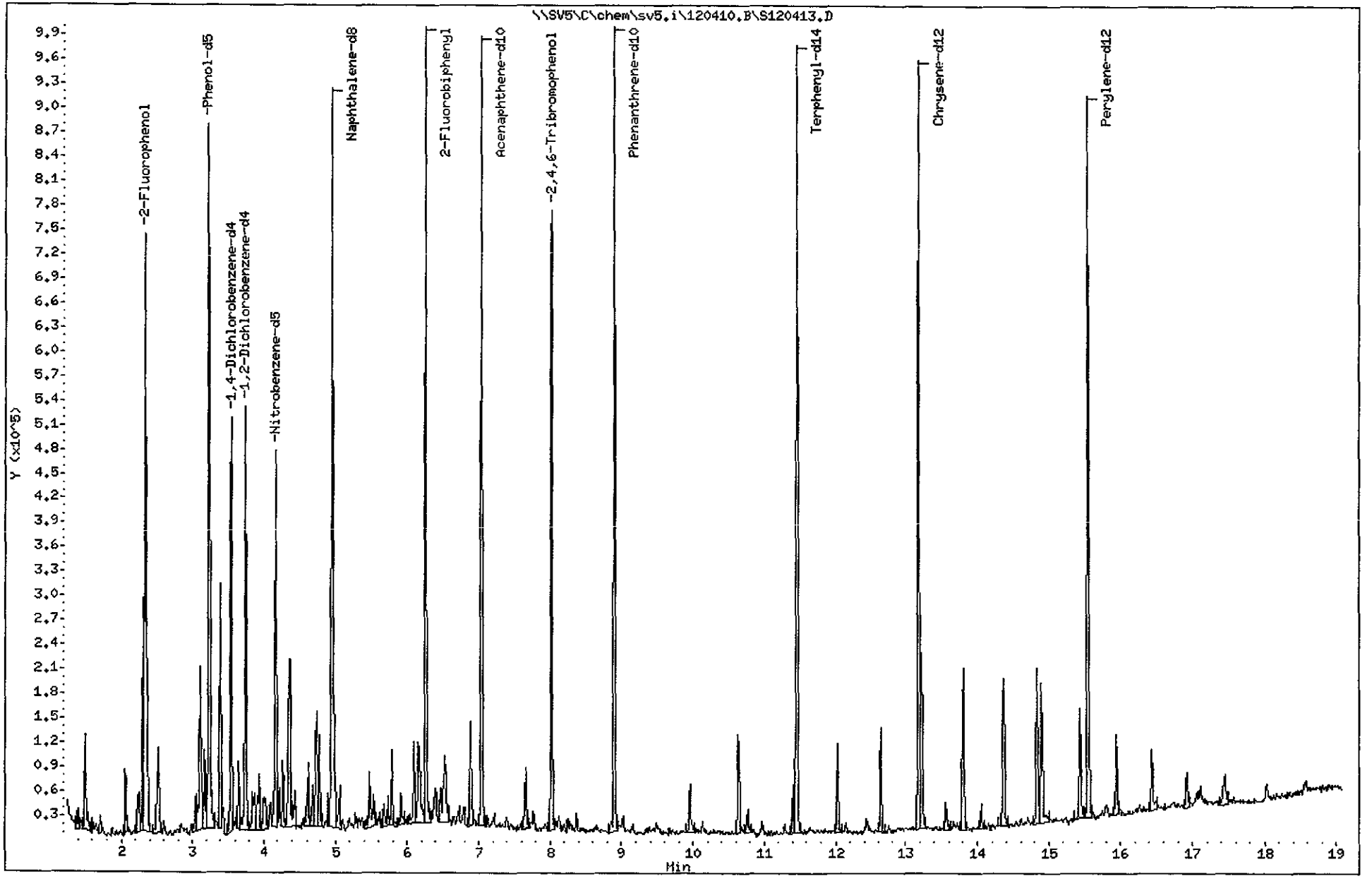
TestAmerica West Sacramento

RECOVERY REPORT

Client Name: Client SDG: 090498
Sample Matrix: GAS Fraction: SV
Lab Smp Id: MAMLW1AA G0K300434- Client Smp ID: 0334389
Level: LOW Operator: KT
Data Type: MS DATA SampleType: SAMPLE
SpikeList File: Quant Type: ISTD
Sublist File: S11JZHCB.SUB
Method File: \\sv5\c\chem\sv5.i\120410.B\8270f.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M

SURROGATE COMPOUND	CONC ADDED ug/L	CONC RECOVERED ug/L	% RECOVERED	LIMITS
$ 7 2-Fluorophenol	100.0	73.24	73.25	41-105
$ 8 Phenol-d5	100.0	85.03	85.03	43-122
$ 10 1,2-Dichlorobenzen	50.00	35.33	70.67	60-120
$ 11 Nitrobenzene-d5	50.00	40.67	81.33	46-118
$ 12 2-Fluorobiphenyl	50.00	42.70	85.40	58-105
$ 13 2,4,6-Tribromophen	100.0	106.4	106.37	61-118
$ 14 Terphenyl-d14	50.00	45.95	91.91	69-110

Data File: \\SV5\C\chem\sv5.i\120410.B\S120413.D
Date : 04-DEC-2010 15:54
Client ID: 0334389
Instrument: sv5.i
Sample Info: MAMLW1AA G0K300434-1;0;;;1000;;1000;5
Volume Injected (uL): 1.0
Operator: KT
Column phase:
Column diameter: 2.00



TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\120410.B\S120414.D
Lab Smp Id: MAML81AA G0K300434- Client Smp ID: 0334389
Inj Date : 04-DEC-2010 16:19
Operator : KT Inst ID: sv5.i
Smp Info : MAML81AA G0K300434-4;0;;;1000;;1000;5
Misc Info : 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\120410.B\8270f.m
Meth Date : 06-Dec-2010 09:40 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 12
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: S11JZHCB.SUB
Target Version: 4.14
Processing Host: SACP307UM

Concentration Formula: Amt * DF * Uf * Vt/(Vo * Vi) * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vt	1000.000	Volume of final extract (uL)
Vo	1000.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected (uL)
Cpnd Variable		Local Compound Variable

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS ON-COLUMN (NG)	FINAL (ug/L)	
* 1	1,4-Dichlorobenzene-d4	152	3.542	3.542	(1.000)	116477	40.0000		(Q)
* 2	Naphthalene-d8	136	4.952	4.952	(1.000)	538761	40.0000		
* 3	Acenaphthene-d10	164	7.045	7.045	(1.000)	305666	40.0000		
* 4	Phenanthrene-d10	188	8.900	8.900	(1.000)	516409	40.0000		
* 5	Chrysene-d12	240	13.180	13.190	(1.000)	511761	40.0000		
* 6	Perylene-d12	264	15.543	15.543	(1.000)	540503	40.0000		
$ 7	2-Fluorophenol	112	2.340	2.340	(0.661)	263497	64.1802	64.18	
$ 8	Phenol-d5	99	3.231	3.231	(0.912)	413699	80.1319	80.13	
$ 10	1,2-Dichlorobenzene-d4	152	3.739	3.739	(1.056)	97568	34.0122	34.01	(q)
$ 11	Nitrobenzene-d5	82	4.164	4.174	(0.841)	167413	36.6874	36.69	
$ 12	2-Fluorobiphenyl	172	6.257	6.257	(0.888)	383099	38.9073	38.91	
$ 13	2,4,6-Tribromophenol	330	8.019	8.019	(1.138)	141369	106.434	106.4	
$ 14	Terphenyl-d14	244	11.460	11.460	(0.869)	468250	46.4518	46.45	
108	Hexachlorobenzene	284	Compound Not Detected.						

12/6/10

QC Flag Legend

Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: S120414.D
Lab Smp Id: MAML81AA G0K300434-
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\SV5\C\chem\sv5.i\120410.B\8270F.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M

Calibration Date: 04-DEC-2010
Calibration Time: 10:36
Client Smp ID: 0334389
Level: LOW
Sample Type: AIR

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	116477	-5.01
2 Naphthalene-d8	530514	265257	1061028	538761	1.55
3 Acenaphthene-d10	282538	141269	565076	305666	8.19
4 Phenanthrene-d10	462722	231361	925444	516409	11.60
5 Chrysene-d12	435850	217925	871700	511761	17.42
6 Perylene-d12	422284	211142	844568	540503	28.00

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.54	3.04	4.04	3.54	0.00
2 Naphthalene-d8	4.95	4.45	5.45	4.95	0.00
3 Acenaphthene-d10	7.05	6.55	7.55	7.05	0.00
4 Phenanthrene-d10	8.90	8.40	9.40	8.90	0.00
5 Chrysene-d12	13.19	12.69	13.69	13.18	-0.08
6 Perylene-d12	15.54	15.04	16.04	15.54	0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

TestAmerica West Sacramento

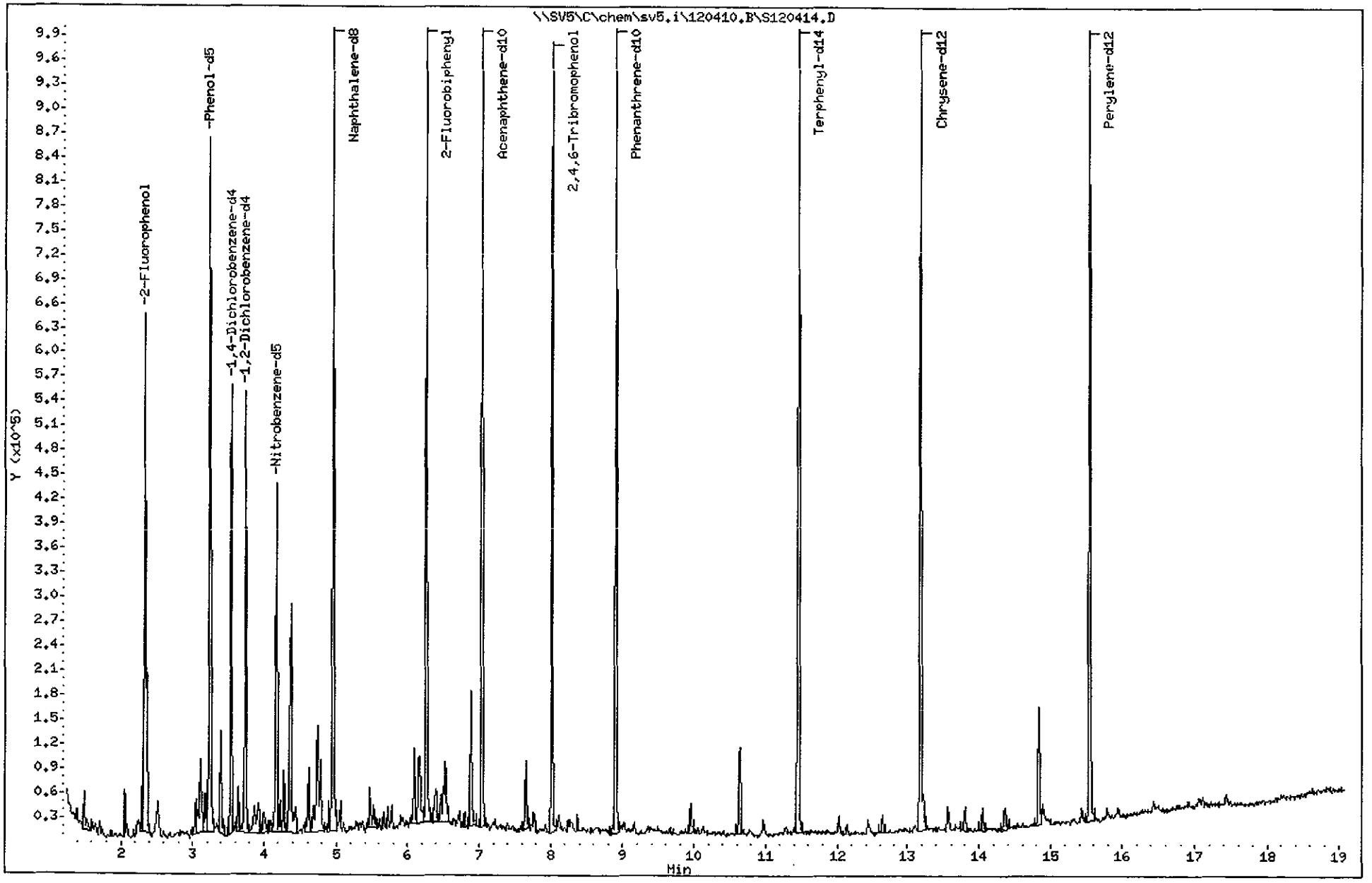
RECOVERY REPORT

Client Name: Client SDG: 090498
Sample Matrix: GAS Fraction: SV
Lab Smp Id: MAML81AA G0K300434- Client Smp ID: 0334389
Level: LOW Operator: KT
Data Type: MS DATA SampleType: SAMPLE
SpikeList File: Quant Type: ISTD
Sublist File: S11JZHCB.SUB
Method File: \\sv5\c\chem\sv5.i\120410.B\8270f.m
Misc Info: 0;AIR;0;S11JZHCB.SUB;;0;0334389;8270F.M

SURROGATE COMPOUND	CONC ADDED ug/L	CONC RECOVERED ug/L	% RECOVERED	LIMITS
$ 7 2-Fluorophenol	100.0	64.18	64.18	41-105
$ 8 Phenol-d5	100.0	80.13	80.13	43-122
$ 10 1,2-Dichlorobenzen	50.00	34.01	68.02	60-120
$ 11 Nitrobenzene-d5	50.00	36.69	73.37	46-118
$ 12 2-Fluorobiphenyl	50.00	38.91	77.81	58-105
$ 13 2,4,6-Tribromophen	100.0	106.4	106.43	61-118
$ 14 Terphenyl-d14	50.00	46.45	92.90	69-110

Data File: \\SV5\C\chem\sv5.i\120410.B\S120414.D
Date : 04-DEC-2010 16:19
Client ID: 0334389
Sample Info: MAML81AA G0K300434-4;0;;;1000;;1000;5
Volume Injected (uL): 1.0
Column phase:

Instrument: sv5.i
Operator: KT
Column diameter: 2.00



Initial Calibration

Includes (as applicable):

runlog

standard raw data

statistical summary

ms tune data

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica West Sacramento
MS SemiVOA ICAL Checklist
Method 8270C

Instrument: SV5

Injection Date: 10/02/10

Initiator/Date: KT-10/03/10

Reviewer/Date: DLZ 10/4/10

DFTPP Mix ID: 10MSSV0129

STD Mix IDs: 10MSSV0307-0313

2nd Source Mix ID: 10MSSV0314, 342

NCM _____

I: SPCCs The SPCC RRFs must be greater than 0.050.

	Initiated	Reviewed		Initiated	Reviewed
N-nitroso-di-n-propylamine	☒	☑	2,4-Dinitrophenol	☒	☑
Hexachlorocyclopentadiene	☒	☑	4-Nitrophenol	☒	☑

II: CCCs The CCC % RSDs must be less than 30%

	Initiated	Reviewed		Initiated	Reviewed
Phenol	☒	☑	Acenaphthene	☒	☑
1,4-Dichlorobenzene	☒	☑	N-nitrosodiphenylamine	☒	☑
2-Nitrophenol	☒	☑	Pentachlorophenol	☒	☑
2,4-Dichlorophenol	☒	☑	Fluoranthene	☒	☑
Hexachlorobutadiene	☒	☑	Di-n-octyl phthalate	☒	☑
4-chloro-3-methylphenol	☒	☑	Benzo(a)pyrene	☒	☑
2,4,6-Trichlorophenol	☒	☑			

III: Other Criteria

	Initiated	Reviewed
The custom.rp shows that the average of the average is less than 15% on the CCV level standard. Avg of AVG:_____	☒	☑
Tailing and degradation criteria are met.	☒	☑
The Tune Documentation is present and meets criteria	☒	☑
All Internal Standards within 50-200% of ICAL mid-point.	☒	☑
Calibration History Included.	☒	☑
Manual re-integrations are checked/initialed and hardcopies included.	☒	☑
Standards analyzed with within 12 hours of Tune time.	☒	☑
Retention time correct for Isomers and all other analytes.	☒	☑
Linear Regressions >0.990 and intercept < ± (½ RL / IS amount)	☒	☑
The second source standard meets the SSCS criteria	☒	☑

File Name: _____

IV: Non-CCC Compounds Over 15% ***(Write compound and %D)***

None

V: Second Source Compounds Over 25% ***(Write compound and %D)***

None

GC/MS INSTRUMENT LOG
SEMI-VOLATILES

Method Key (MTH Column)
QL = EPA 8270C (WS-MS-0005)
JZ = EPA TO-13A (WS-MS-0005).
VX = EPA 8270C-SIM (mod) CWM (WS-MS-0003)
QI = EPA 8270C-SIM (WS-MS-0008)
FX = PAH-SIM Isotope Dilution (WS-MS-0006)
F9 = EPA 8270C-SIM (mod) 1,4-Dioxane (WS-MS-0011)

Inst ID : sv5.i
Batch ID : 100210.B
ICAL Date: See Calib Report
See raw data for standard IDs

Date	Time	USER	Sample ID	File ID	Vol or Wt	Extract Vol	Dilu	MTH	Comments
02-OCT-2010	11:43	KT	Primer	QC001.D	NA	NA	NA		
02-OCT-2010	12:06	KT	DFTPP 50ug/ml	DFT1002.D	NA	NA	NA		
02-OCT-2010	12:27	KT	HSL_005 ug/ml CS-1	HSL1002A.	NA	NA	NA		
02-OCT-2010	12:53	KT	HSL_010 ug/ml CS-2	HSL1002B.	NA	NA	NA		
02-OCT-2010	13:18	KT	HSL_020 ug/ml CS-3	HSL1002C.	NA	NA	NA		
02-OCT-2010	13:44	KT	HSL_050 ug/ml CS-4	HSL1002D.	NA	NA	NA		
02-OCT-2010	14:09	KT	HSL_080 ug/ml CS-5	HSL1002E.	NA	NA	NA		
02-OCT-2010	14:35	KT	HSL_120 ug/ml CS-6	HSL1002F.	NA	NA	NA		
02-OCT-2010	15:00	KT	HSL_160 ug/ml CS-7	HSL1002G.	NA	NA	NA		
02-OCT-2010	16:11	KT	HSL_050 ug/ml ICV	HSL1002H.	NA	NA	NA		
02-OCT-2010	16:36	KT	Benzidines ICV 50ug/mL	HSL1002H1	NA	NA	NA		

SV5 HSL
10/5/10

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:09 onishim

Calibration File Names:
Level 1: \\SV5\C\chem\sv5.i\081710.B\AP90817A.D
Level 2: \\SV5\C\chem\sv5.i\081710.B\AP90817B.D
Level 3: \\SV5\C\chem\sv5.i\081710.B\AP90817C.D
Level 4: \\SV5\C\chem\sv5.i\081710.B\AP90817D.D
Level 5: \\SV5\C\chem\sv5.i\081710.B\AP90817E.D
Level 6: \\SV5\C\chem\sv5.i\081710.B\AP90817F.D
Level 7: \\SV5\C\chem\sv5.i\081710.B\AP90817G.D

Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
15 N-Nitrosodimethylamine	0.92899	0.88268	0.91048	0.91970	0.93146	0.93916					
	0.93833						AVRG		0.92154		2.16207
16 Pyridine	1.67117	1.37423	1.59449	1.56610	1.62399	1.53256					
	1.52523						AVRG		1.54111		5.85560
23 Aniline	2.20796	2.15935	2.19988	2.26058	2.29749	2.33400					
	2.33783						AVRG		2.25673		3.09783
24 Phenol	2.04111	1.96212	2.02834	2.03430	2.06683	2.06089					
	2.06740						AVRG		2.03729		1.80250

Manual correlation for 2,4,5-Trichlorophenol @ Level 3:

$\frac{55529}{328608} \times \frac{60}{20} = 0.33796$ by 10/4/10

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:09 onishim

Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
26 Bis(2-chloroethyl)ether	1.47335	1.38282	1.39491	1.43824	1.42549	1.44300					
	1.44264						AVRG		1.42859		2.17028
27 2-Chlorophenol	1.52099	1.55595	1.56903	1.58168	1.56789	1.58074					
	1.57039						AVRG		1.56381		1.32808
28 1,3-Dichlorobenzene	1.68903	1.69173	1.67754	1.73135	1.68641	1.72299					
	1.72457						AVRG		1.70337		1.29370
29 1,4-Dichlorobenzene	1.77122	1.79861	1.74013	1.76898	1.78200	1.79288					
	1.81444						AVRG		1.78118		1.35229
30 Benzyl Alcohol	1.01643	1.03654	0.99182	1.04980	1.07792	1.08952					
	1.09506						AVRG		1.05101		3.69696
31 1,2-Dichlorobenzene	1.62008	1.63185	1.60455	1.68061	1.63410	1.64415					
	1.64691						AVRG		1.63746		1.45884
32 2-Methylphenol	1.40818	1.38930	1.39110	1.42620	1.45565	1.46154					
	1.47889						AVRG		1.43012		2.60558

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
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Quant Method : ISTD
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:09 onishim

Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
33 2,2'-oxybis(1-Chloropropane)	2.29602	2.22080	2.28329	2.27928	2.27018	2.27830					
	2.28770						AVRG		2.27365		1.08468
34 4-Methylphenol	1.48606	1.48913	1.46270	1.52239	1.52653	1.55886					
	1.58763						AVRG		1.51904		2.88378
36 Hexachloroethane	0.60928	0.60836	0.60573	0.61394	0.60427	0.59381					
	0.60919						AVRG		0.60636		1.04319
37 N-Nitrosodinpropylamine	0.94498	0.97005	1.01302	1.02370	1.04700	1.03627					
	1.04757						AVRG		1.01180		3.92615
42 Nitrobenzene	0.32855	0.32602	0.32543	0.33083	0.33379	0.33450					
	0.33901						AVRG		0.33116		1.48904
44 Isophorone	0.63431	0.62291	0.61160	0.63344	0.63648	0.66468					
	0.65411						AVRG		0.63679		2.81109
45 2-Nitrophenol	0.18608	0.18833	0.18840	0.20021	0.20022	0.20702					
	0.20808						AVRG		0.19648		4.42274

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

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Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:09 onishim

Compound	5.0000 Level 1 / 160.0000 Level 7	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
46 2,4-Dimethyphenol	0.34459	0.34167	0.34307	0.34912	0.34788	0.35962					
	0.35785						AVRG		0.34911		2.02786
47 Bis(2-chloroethoxy)methane	0.41146	0.37494	0.38565	0.38249	0.38500	0.39859					
	0.38545						AVRG		0.38908		3.10601
49 2,4-Dichlorophenol	0.25434	0.26318	0.27019	0.27037	0.27274	0.28180					
	0.27809						AVRG		0.27010		3.39345
50 Benzoic Acid	0.16747	0.16266	0.17423	0.19357	0.21024	0.22272					
	0.22180						AVRG		0.19324		13.25202
51 1,2,4-Trichlorobenzene	0.29430	0.28827	0.28475	0.29747	0.29189	0.29959					
	0.29091						AVRG		0.29246		1.75989
52 Naphthalene	1.09939	1.12462	1.07435	1.09325	1.09870	1.13821					
	1.10247						AVRG		1.10443		1.89960
54 4-Chloroaniline	0.40751	0.42534	0.43264	0.43910	0.43781	0.44905					
	0.43867						AVRG		0.43288		3.06843

TestAmerica West Sacramento

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Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
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Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
57 Hexachlorobutadiene	0.14295	0.13812	0.14428	0.14415	0.14385	0.14379					
	0.14473						AVRG		0.14313		1.58904
60 4-Chloro-3-Methylphenol	0.29329	0.28866	0.29079	0.30972	0.30295	0.31766					
	0.30839						AVRG		0.30164		3.64422
63 2-Methylnaphthalene	0.68483	0.68064	0.68080	0.70067	0.70860	0.71172					
	0.69217						AVRG		0.69378		1.79740
66 Hexachlorocyclopentadiene	0.26878	0.27757	0.28896	0.29704	0.30236	0.32262					
	0.33186						AVRG		0.29846		7.64489
69 2,4,6-Trichlorophenol	0.31186	0.29820	0.30223	0.31996	0.32305	0.34225					
	0.33638						AVRG		0.31913		5.18654
70 2,4,5-Trichlorphenol	0.30823	0.32892	0.33796	0.36298	0.35236	0.35480					
	0.36135						AVRG		0.34380		5.80662
71 2-Chloronaphthalene	1.13629	1.09411	1.10012	1.14181	1.11220	1.14447					
	1.15096						AVRG		1.13571		2.05054

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Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
73 2-Nitroaniline	0.31576	0.31759	0.33397	0.35205	0.34821	0.35794					
	0.36278						AVRG		0.34119		5.87334
76 Dimethylphthalate	1.23388	1.28191	1.29803	1.34568	1.31165	1.32891					
	1.30237						AVRG		1.29606		3.09317
77 Acenaphthylene	1.86531	1.91304	1.91818	2.01646	1.98204	1.99786					
	2.02968						AVRG		1.96037		3.15026
79 2,6-Dinitrotoluene	0.28347	0.27378	0.29890	0.31220	0.31294	0.32140					
	0.31106						AVRG		0.30197		5.78579
80 3-Nitroaniline	0.35362	0.34622	0.35978	0.40036	0.38674	0.39559					
	0.39603						AVRG		0.37691		6.06861
81 Acenaphthene	1.25874	1.22468	1.26733	1.27046	1.21141	1.24781					
	1.25463						AVRG		1.24787		1.76776
82 2,4-Dinitrophenol	4083	7537	23799	58864	110384	199007					
	266655						QUAD	0.10620	5.22413	-0.71963	0.99812

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Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
83 Dibenzofuran	1.57786	1.62124	1.65200	1.69530	1.65117	1.68460					
	1.71077						AVRG		1.65612		2.77923
84 4-Nitrophenol	0.12712	0.14148	0.15316	0.16076	0.17130	0.16653					
	0.17404						AVRG		0.15634		10.90920
86 2,4-Dinitrotoluene	0.34360	0.36989	0.38479	0.42164	0.41035	0.42305					
	0.43110						AVRG		0.39633		8.61892
91 Fluorene	1.34667	1.33840	1.34292	1.39902	1.38899	1.37835					
	1.40640						AVRG		1.37139		2.08557
92 Diethylphthalate	1.22240	1.29889	1.31549	1.37912	1.31873	1.37345					
	1.38087						AVRG		1.32699		4.31889
93 4-Chlorophenyl-phenylether	0.54964	0.56917	0.56887	0.59265	0.56708	0.57698					
	0.57695						AVRG		0.57019		2.42913
94 4-Nitroaniline	0.33346	0.33747	0.37329	0.38337	0.39216	0.39102					
	0.40452						AVRG		0.37361		7.42395

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Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
97 4,6-Dinitro-2-methylphenol	5780	11282	32982	76137	134784	236477					
	324244						LINR	0.10840	0.15581		0.99840
98 N-Nitrosodiphenylamine	0.87756	0.59736	0.60533	0.60433	0.62172	0.61801					
	0.61968						AVRG		0.60628		2.57715
100 Azobenzene	0.77527	0.76965	0.77321	0.79522	0.80064	0.81892					
	0.77331						AVRG		0.78660		2.37146
101 4-Bromophenyl-phenylether	0.18964	0.18507	0.19281	0.19931	0.19607	0.20581					
	0.19815						AVRG		0.19527		3.48752
108 Hexachlorobenzene	0.22958	0.22064	0.20740	0.21605	0.21731	0.21704					
	0.21854						AVRG		0.21807		3.00928
110 Pentachlorophenol	5949	10551	30451	67882	126397	215360					
	293184						LINR	0.09816	0.14122		0.99845
114 Phenanthrene	1.30347	1.26007	1.25408	1.24163	1.24375	1.25610					
	1.26611						AVRG		1.26074		1.64308

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Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
115 Anthracene	1.25034	1.21759	1.24206	1.25982	1.27529	1.30214					
	1.26950						AVRG		1.25955		2.12888
118 Carbazole	1.13211	1.12547	1.13694	1.14260	1.17067	1.18192					
	1.16456						AVRG		1.15061		1.87826
120 Di-n-Butylphthalate	1.28492	1.32287	1.36193	1.38164	1.41474	1.43847					
	1.48636						AVRG		1.38442		4.97257
126 Fluoranthene	1.03840	1.07611	1.17216	1.10620	1.15861	1.18294					
	1.17440						AVRG		1.12969		5.01774
127 Benzidine	0.78175	0.76431	0.75260	0.82658	0.82201	0.86375					
	0.86381						AVRG		0.81067		5.60614
128 Pyrene	1.25791	1.33783	1.17078	1.28684	1.25586	1.28463					
	1.25794						AVRG		1.25025		3.12172
134 3,3'-dimethylbenzidine	0.68472	0.64388	0.67361	0.70756	0.73630	0.79414					
	0.79926						AVRG		0.71564		8.88815

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Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
136 Butylbenzylphthalate	0.64964	0.60187	0.59142	0.62586	0.61590	0.65233					
	0.64920						AVRG		0.62663		3.95034
138 Benzo(a)Anthracene	1.10169	0.99731	1.03245	1.04489	1.06449	1.10831					
	1.10920						AVRG		1.06648		4.05847
139 Chrysene	1.05284	1.10175	1.06320	1.09705	1.06985	1.12241					
	1.12246						AVRG		1.08994		2.59426
140 3,3'-Dichlorobenzidine	0.39148	0.37695	0.39090	0.39906	0.40353	0.42717					
	0.42415						AVRG		0.40189		4.83885
141 bis(2-ethylhexyl)Phthalate	0.91826	0.80897	0.84032	0.86193	0.84371	0.89539					
	0.88354						AVRG		0.86316		4.34816
142 Di-n-octylphthalate	1.34838	1.23185	1.35627	1.34433	1.39356	1.47616					
	1.50770						AVRG		1.37975		6.66055
144 Benzo(b)fluoranthene	0.81012	0.81077	0.82747	0.99930	0.98373	0.91132					
	1.02672						AVRG		0.90549		10.05836

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Compound	5.0000 Level 1 / 160.0000 Level 7	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	Coefficients m1	Coefficients m2	%RSD or R^2
145 Benzo(k)fluoranthene	1.22939	1.16528	1.20022	1.09895	1.14223	1.19597					
	1.10447						AVRG		1.16236		4.27893
147 Benzo(e)pyrene	0.90394	0.92734	0.90757	0.95977	0.96997	0.96929					
	0.97186						AVRG		0.94425		3.22007
148 Benzo(a)pyrene	0.98300	0.97686	0.99402	1.02789	1.07610	1.06275					
	1.06523						AVRG		1.02655		4.11137
151 Indeno(1,2,3-cd)pyrene	0.73783	0.73267	0.73671	0.84698	0.84057	0.93730					
	0.97995						AVRG		0.83029		12.15083
152 Dibenzo(a,h)anthracene	0.88099	0.84384	0.87256	0.92240	0.95990	1.00944					
	1.00392						AVRG		0.92756		7.07091
153 Benzo(g,h,i)perylene	0.96026	0.98487	0.97380	0.99974	1.01731	1.05397					
	1.04026						AVRG		1.00427		3.45188
M 162 benzo b,k Fluoranthene Totals	2.03951	1.97605	2.02770	2.09825	2.09596	2.10729					
	2.13019						AVRG		2.06785		2.64859

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Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
$ 7 2-Fluorophenol	1.44503	1.30436	1.38373	1.44170	1.43635	1.42292					
	1.43635						AVRG		1.40992		3.61494
$ 8 Phenol-d5	1.72227	1.67335	1.74151	1.79006	1.80863	1.83864					
	1.83627						AVRG		1.77296		3.52001
$ 9 2-Chlorophenol-d4	1.47770	1.55530	1.53916	1.59414	1.57486	1.57967					
	1.57804						AVRG		1.55698		2.52388
$ 10 1,2-Dichlorobenzene-d4	0.95776	0.98111	0.99827	0.98914	0.99518	0.98547					
	0.98696						AVRG		0.98513		1.36559
$ 11 Nitrobenzene-d5	0.33028	0.34256	0.33065	0.34108	0.33606	0.35127					
	0.33970						AVRG		0.33879		2.16217
$ 12 2-Fluorobiphenyl	1.28499	1.26007	1.27668	1.34206	1.25854	1.29723					
	1.30010						AVRG		1.28852		2.22622
$ 13 2,4,6-Tribromophenol	0.15034	0.16527	0.17466	0.17926	0.17825	0.18501					
	0.18390						AVRG		0.17381		7.05197

Report Date : 03-Oct-2010 11:10

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INITIAL CALIBRATION DATA

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Last Edit : 03-Oct-2010 11:09 onishim

Compound	5.0000 Level 1	10.0000 Level 2	20.0000 Level 3	50.0000 Level 4	80.0000 Level 5	120.0000 Level 6	Curve	Coefficients b	m1	m2	%RSD or R^2
	160.0000 Level 7										
$ 14 Terphenyl-d14	0.78508	0.78616	0.73917	0.80441	0.78047	0.81889					
	0.80107						AVRG		0.78789		3.21384

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Curve	Formula	Units
Averaged	Amt = Rsp/m1	Response
Linear	Amt = b + Rsp/m1	Response
Quad	Amt = b + m1*Rsp + m2*Rsp^2	Response

Signal Calibration Report

Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit: 04-Oct-2010 09:00 onishim
Compound : 82 2,4-Dinitrophenol
Mass: 184.00
Istd Compound: * 3 Acenaphthene-d10

Calibration Formulas

Calibration Mode: by Response

Curve Type: Averaged
Origin: None
Amt = Rsp/m1
m1 = 0.15933171100000
RSD: 26.349

Initial Calibration Table

Lvl	RT	Amount	Response	RT	Istd Amount	Istd Response	Response Factor
1	7.572	5.00000	4083	7.468	40.000	321839	0.10149173965865
2	7.572	10.00000	7537	7.468	40.000	272639	0.110578457227732
3	7.572	20.00000	23799	7.468	40.000	328608	0.144847356512036
4	7.582	50.00000	58864	7.468	40.000	282538	0.16667209366528
5	7.572	80.00000	110384	7.468	40.000	300315	0.183780363951118
6	7.582	120.00000	199007	7.468	40.000	322596	0.205630778641160
7	7.582	160.00000	265655	7.478	40.000	328259	0.202321185405443

Lvl	Sublist	Calibration File
1	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002A
2	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002B
3	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002C
4	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002D
5	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002E
6	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002F
7	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002G

Continuing Calibration Table

Ind	RT	Amount	Response	RT	Istd Amount	Istd Response	Response Factor

1	7.582	50.000	50142	7.468	40.000	236662	0.16949742670982
2	7.572	50.000	58864	7.468	40.000	282538	0.16667209366528
3	7.582	50.000	56608	7.468	40.000	239304	0.18924213552636
4	7.589	50.000	98553	7.485	40.000	440855	0.17883975456783
5	7.599	50.000	81881	7.485	40.000	371846	0.17616109894957
6	7.599	50.000	55069	7.495	40.000	283828	0.15521794889863
7	7.599	50.000	52896	7.496	40.000	256342	0.16507946415336
8	7.599	50.000	50586	7.495	40.000	224545	0.18022578993075
9	7.610	50.000	31559	7.506	40.000	165705	0.15236233064784
10	7.610	50.000	50181	7.506	40.000	226619	0.17714666456625
11	7.610	50.000	44092	7.506	40.000	201923	0.17468837130986
12	7.620	50.000	81056	7.516	40.000	329174	0.19699247206645
13	7.620	50.000	93793	7.516	40.000	378407	0.19829020076267
14	7.630	50.000	68549	7.516	40.000	271629	0.20189007801082
15	7.630	50.000	54835	7.516	40.000	219680	0.19969045884926
16	7.630	50.000	67628	7.527	40.000	267569	0.20219980640508
17	7.630	50.000	94376	7.527	40.000	349016	0.21632475301992
18	7.635	50.000	51607	7.532	40.000	209252	0.19730086211840
19	7.635	50.000	62563	7.531	40.000	260404	0.19220288474631
20	7.646	50.000	80386	7.542	40.000	334425	0.19229662854153
21	7.645	50.000	25473	7.542	40.000	302573	0.06735035842590
22	7.645	50.000	17649	7.542	40.000	223404	0.06320030080034
23	7.646	50.000	68382	7.542	40.000	292758	0.18686286967393
24	7.656	50.000	97952	7.552	40.000	390143	0.20085353319168
25	7.656	50.000	63647	7.552	40.000	289221	0.17605084001507
26	7.666	50.000	79703	7.563	40.000	331752	0.19219899201813
27	7.677	50.000	59624	7.573	40.000	245725	0.19411618679418
28	7.687	50.000	60561	7.583	40.000	237909	0.20364425053277
29	7.687	50.000	42226	7.583	40.000	172923	0.19535168832370
30	7.687	50.000	51997	7.583	40.000	208221	0.19977619932668
31	7.697	50.000	51275	7.594	40.000	202822	0.20224630454151
32	7.697	50.000	65531	7.594	40.000	250339	0.20941523294413
33	7.760	50.000	76785	7.656	40.000	344524	0.17829817371214

34	7.759	50.000	68725	7.656	40.000	303207	0.18132826748723
35	7.770	50.000	66249	7.666	40.000	308864	0.17159397016162
36	7.780	50.000	63983	7.677	40.000	288883	0.17718730420274
37	7.780	50.000	61267	7.677	40.000	292290	0.16768825481542
38	7.791	50.000	56069	7.687	40.000	238922	0.18773993186061
39	7.791	50.000	50573	7.687	40.000	243613	0.166076523000986
40	7.791	50.000	55930	7.687	40.000	256301	0.17457598682799
41	7.791	50.000	55930	7.687	40.000	256301	0.17457598682799
42	7.791	50.000	43995	7.687	40.000	215682	0.16318468856928
43	7.801	50.000	55663	7.697	40.000	269061	0.16550299002828
44	7.801	50.000	52406	7.697	40.000	242418	0.17294425331452
45	7.801	50.000	49689	7.697	40.000	246748	0.161100393392417
46	7.801	50.000	83728	7.697	40.000	361851	0.18511044601231
47	7.801	50.000	69470	7.697	40.000	316865	0.17539330629763
48	7.811	50.000	98764	7.708	40.000	448001	0.17636389204488
49	7.811	50.000	65199	7.708	40.000	319060	0.16347771579013
50	7.811	50.000	63819	7.708	40.000	326041	0.15659134894078
51	7.811	50.000	69420	7.708	40.000	325539	0.17059707131864
52	7.822	50.000	66513	7.718	40.000	295770	0.17990465564459
53	7.822	50.000	58901	7.718	40.000	274779	0.17148617616339
54	7.822	50.000	58321	7.718	40.000	264752	0.17622831933281
55	7.816	50.000	90734	7.713	40.000	414154	0.17526620532459
56	7.858	50.000	49564	7.754	40.000	260934	0.151958732859651
57	7.858	50.000	63475	7.754	40.000	318667	0.15935129774969
58	7.889	50.000	58884	7.785	40.000	318462	0.14792094504211
59	7.889	50.000	52456	7.796	40.000	304639	0.13775255302177
60	7.889	50.000	44855	7.796	40.000	283970	0.126365461140261
61	7.889	50.000	40711	7.785	40.000	264293	0.123229900148701
Avg	7.719	50.000	61661	7.615	40.000	4333	0.173642339865731

Ind	Sublist	Calibration File
1	1_8270STD	\\sv5\c\chem\sv5.i\100210.B\HSL1002H

2	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002D
3	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\QC001
4	1_8270STD	\\SV5\C\chem\sv5.i\100110.B\HSL1001
5	1_8270STD	\\SV5\C\chem\sv5.i\093010.B\HSL0930
6	1_8270STD	\\sv5\c\chem\sv5.i\092910A.B\HSL0929A
7	1_8270STD	\\SV5\C\chem\sv5.i\092910.B\HSL0929
8	1_8270STD	\\SV5\C\chem\sv5.i\092910.B\QC001
9	1_8270STD	\\SV5\C\chem\sv5.i\092810A.B\HSL0928
10	1_8270STD	\\SV5\C\chem\sv5.i\092810.B\HSL0928
11	1_8270STD	\\SV5\C\chem\sv5.i\092710.B\HSL0927
12	1_8270STD	\\SV5\C\chem\sv5.i\092510.B\QC001
13	1_8270STD	\\sv5\c\chem\sv5.i\092510.B\HSL0925
14	1_8270STD	\\SV5\C\chem\sv5.i\092410.B\QC001
15	1_8270STD	\\SV5\C\chem\sv5.i\092410.B\HSL0924
16	1_8270STD	\\SV5\C\chem\sv5.i\092310A.B\HSL0923A
17	1_8270STD	\\SV5\C\chem\sv5.i\092310A.B\QC001
18	1_8270STD	\\SV5\C\chem\sv5.i\092310.B\QC001
19	1_8270STD	\\SV5\C\chem\sv5.i\092310.B\HSL0923
20	1_8270STD	\\SV5\C\chem\sv5.i\092210.B\HSL0922a
21	1_8270STD	\\SV5\C\chem\sv5.i\092210.B\HSL0922
22	1_8270STD	\\SV5\C\chem\sv5.i\092210.B\QC001
23	1_8270STD	\\SV5\C\chem\sv5.i\092110.B\HSL0921
24	1_8270STD	\\SV5\C\chem\sv5.i\092010.B\QC001
25	1_8270STD	\\SV5\C\chem\sv5.i\092010.B\HSL0920
26	1_8270STD	\\SV5\C\chem\sv5.i\091910a.B\HSL0919a
27	1_8270STD	\\SV5\C\chem\sv5.i\091910.B\HSL0919
28	1_8270STD	\\SV5\C\chem\sv5.i\091910.B\QC001
29	1_8270STD	\\SV5\C\chem\sv5.i\091710.B\HSL0917
30	1_8270STD	\\SV5\C\chem\sv5.i\091710.B\QC001
31	1_8270STD	\\SV5\C\chem\sv5.i\091510b.B\HSL0915b
32	1_8270STD	\\SV5\C\chem\sv5.i\091510b.B\QC003
33	1_8270STD	\\sv5\c\chem\sv5.i\091010.B\HSL0910
34	1_8270STD	\\SV5\C\chem\sv5.i\091010.B\QC001

35	1_8270STD	\\sv5\c\chem\sv5.i\090918a.B\ESL0909a
36	1_8270STD	\\SV5\C\chem\sv5.i\090910.B\HSL0909
37	1_8270STD	\\SV5\C\chem\sv5.i\090910.B\QC001
38	1_8270STD	\\SV5\C\chem\sv5.i\090810.B\HSL0908
39	1_8270STD	\\SV5\C\chem\sv5.i\090810.B\Primer
40	1_8270STD	\\sv5\c\chem\sv5.i\090710.B\HSL0907
41	1_8270STD	\\SV5\C\chem\sv5.i\090710.B\HSL0907
42	1_8270STD	\\sv5\c\chem\sv5.i\090110.B\HSL0901
43	1_8270STD	\\SV5\C\chem\sv5.i\083110.B\HSL0831
44	1_8270STD	\\sv5\c\chem\sv5.i\083010.B\QC001
45	1_8270STD	\\sv5\c\chem\sv5.i\083010.B\HSL0830
46	1_8270STD	\\SV5\C\chem\sv5.i\082710.B\QC001
47	1_8270STD	\\sv5\c\chem\sv5.i\082710.B\ESL0827
48	1_8270STD	\\SV5\C\chem\sv5.i\082610.B\HSL0826
49	1_8270STD	\\SV5\C\chem\sv5.i\082610.B\QC001
50	1_8270STD	\\SV5\C\chem\sv5.i\082510.B\QC001
51	1_8270STD	\\SV5\C\chem\sv5.i\082510.B\HSL0825
52	1_8270STD	\\sv5\c\chem\sv5.i\082310B.B\HSL0823
53	1_8270STD	\\sv5\c\chem\sv5.i\082310B.B\HSL0823H
54	1_8270STD	\\sv5\c\chem\sv5.i\082310B.B\HSL0823D
55	1_8270STD	\\SV5\C\chem\sv5.i\082310A.B\HSL0823A
56	1_8270STD	\\SV5\C\chem\sv5.i\082010.B\HSL0820
57	1_8270STD	\\sv5\c\chem\sv5.i\082010.B\QC001
58	1_8270STD	\\sv5\c\chem\sv5.i\081810A.B\HSL0818A
59	1_8270STD	\\sv5\c\chem\sv5.i\081810.B\HSL0818
60	1_8270STD	\\SV5\C\chem\sv5.i\081710.B\HSL0817D
61	1_8270STD	\\SV5\C\chem\sv5.i\081710.B\HSL0817H

Signal Calibration Report

Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit: 04-Oct-2010 09:00 onishim
Compound : 110 Pentachlorophenol
Mass: 266.00
Istd Compound: * 4 Phenanthrene-d10

Calibration Formulas

Calibration Mode: by Response

Curve Type: Averaged
Origin: None
Amt = Rsp/m1
m1 = 0.119308974000000
RSD: 15.221

Initial Calibration Table

Lvl	RT	Amount	Response	RT	Istd Amount	Istd Response	Response Factor
1	9.240	5.00000	5849	9.406	40.000	496356	0.094271047393340
2	9.240	10.00000	10551	9.406	40.000	428440	0.098506208570630
3	9.240	20.00000	30451	9.406	40.000	525834	0.115819821464570
4	9.240	50.00000	67882	9.406	40.000	462722	0.117361180147040
5	9.240	80.00000	126397	9.406	40.000	477777	0.132276145565820
6	9.240	120.00000	215360	9.406	40.000	515607	0.139227486567610
7	9.250	160.00000	293184	9.406	40.000	532284	0.137700926573030

Lvl	Sublist	Calibration File
1	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002A
2	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002B
3	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002C
4	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002D
5	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002E
6	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002F
7	1_8270STD	\\SV5\C\chem\sv5.i\100210.B\HSL1002G

Continuing Calibration Table

Ind	RT	Amount	Response	RT	Istd Amount	Istd Response	Response Factor

1	9.240	50.000	62906	9.406	40.000	380734	0.132178371251321
2	9.240	50.000	67882	9.406	40.000	462722	0.117361180147041
3	9.257	50.000	111129	9.423	40.000	692643	0.128352567422181
4	9.257	50.000	88353	9.423	40.000	569627	0.124085410277251
5	9.267	50.000	65176	9.433	40.000	444572	0.117283139738891
6	9.268	50.000	60910	9.433	40.000	402268	0.121133174898231
7	9.278	50.000	51724	9.433	40.000	342388	0.120854702851741
8	9.278	50.000	37406	9.444	40.000	257561	0.116185292027911
9	9.278	50.000	56153	9.444	40.000	367144	0.122356350641711
10	9.278	50.000	49979	9.444	40.000	316244	0.126431489609291
11	9.299	50.000	89278	9.465	40.000	533339	0.133915577146991
12	9.288	50.000	102299	9.454	40.000	604130	0.135466207604321
13	9.299	50.000	74887	9.464	40.000	434948	0.137739683824271
14	9.299	50.000	61171	9.465	40.000	350214	0.139733991216801
15	9.309	50.000	72641	9.475	40.000	436116	0.133250786488001
16	9.309	50.000	99213	9.475	40.000	545533	0.145491473476401
17	9.314	50.000	56050	9.480	40.000	341600	0.131264637002341
18	9.314	50.000	67187	9.480	40.000	410196	0.131033944748361
19	9.324	50.000	90596	9.490	40.000	530756	0.136553896705831
20	9.324	50.000	32043	9.490	40.000	484990	0.052855522794281
21	9.324	50.000	22238	9.490	40.000	346959	0.051275222720841
22	9.324	50.000	81528	9.490	40.000	462218	0.141107442808371
23	9.335	50.000	103580	9.511	40.000	589949	0.140459599050091
24	9.335	50.000	72155	9.501	40.000	446339	0.129327708311401
25	9.355	50.000	91662	9.521	40.000	517550	0.141686020674331
26	9.366	50.000	67431	9.532	40.000	396847	0.135933495780491
27	9.366	50.000	71407	9.542	40.000	407176	0.140297070554261
28	9.366	50.000	49946	9.532	40.000	298933	0.133664734238111
29	9.366	50.000	58621	9.542	40.000	335623	0.139730590573351
30	9.386	50.000	53858	9.552	40.000	329730	0.130671761744461
31	9.387	50.000	69993	9.552	40.000	399673	0.140100532185061
32	9.459	50.000	87217	9.625	40.000	539077	0.129431602535441
33	9.459	50.000	77540	9.625	40.000	458679	0.135240549491041

34	9.470	50.000	75232	9.646	40.000	482971	0.131241006188781
35	9.480	50.000	75075	9.656	40.000	465501	0.129022279221741
36	9.480	50.000	69872	9.656	40.000	435300	0.128411670112571
37	9.490	50.000	60626	9.656	40.000	378611	0.128101930477461
38	9.490	50.000	60476	9.666	40.000	383533	0.126145077477041
39	9.490	50.000	68275	9.656	40.000	401081	0.136181968230861
40	9.490	50.000	68275	9.656	40.000	401081	0.136181968230861
41	9.490	50.000	51783	9.666	40.000	337799	0.122636242262411
42	9.501	50.000	70205	9.677	40.000	425699	0.131933596273421
43	9.511	50.000	60939	9.677	40.000	381025	0.127947510005911
44	9.501	50.000	61157	9.677	40.000	380328	0.128640541848091
45	9.500	50.000	98266	9.676	40.000	586969	0.133930071264411
46	9.500	50.000	82460	9.677	40.000	500580	0.131783131567381
47	9.511	50.000	117721	9.687	40.000	687233	0.137037656806351
48	9.511	50.000	77582	9.687	40.000	485585	0.127816139295901
49	9.511	50.000	77449	9.687	40.000	498103	0.124390336938341
50	9.511	50.000	85917	9.687	40.000	500311	0.137381748552401
51	9.521	50.000	80098	9.697	40.000	460974	0.139006538329711
52	9.521	50.000	71155	9.697	40.000	428920	0.132714725356711
53	9.521	50.000	72603	9.697	40.000	415811	0.139684616328091
54	9.526	50.000	108254	9.702	40.000	650674	0.133097680251551
55	9.568	50.000	64139	9.744	40.000	411802	0.124601628938181
56	9.578	50.000	85309	9.754	40.000	511730	0.133365642037791
57	9.599	50.000	78595	9.785	40.000	486034	0.129365435339911
58	9.609	50.000	72755	9.785	40.000	467607	0.124472045970231
59	9.609	50.000	67958	9.785	40.000	451801	0.120332624319111
60	9.609	50.000	63635	9.785	40.000	418038	0.121778402920311
Avg	9.411	50.000	72233	9.581	40.000	6967	0.128494282418101

Ind	Sublist	Calibration File
1	1_8270STD	\\svs\c\chem\svs.i\100210.B\HSL1002H
2	1_8270STD	\\SVS\C\chem\svs.i\100210.B\HSL1002D

3	1_8270STD	\\SV5\C\chem\sv5.i\100110.B\HSL1001
4	1_8270STD	\\SV5\C\chem\sv5.i\053010.B\HSL0930
5	1_8270STD	\\sv5\c\chem\sv5.i\092910A.B\HSL0929A
6	1_8270STD	\\SV5\C\chem\sv5.i\092910.B\HSL0929
7	1_8270STD	\\SV5\C\chem\sv5.i\092910.B\QC001
8	1_8270STD	\\SV5\C\chem\sv5.i\092810A.B\HSL0928
9	1_8270STD	\\SV5\C\chem\sv5.i\092810.B\HSL0928
10	1_8270STD	\\SV5\C\chem\sv5.i\092710.B\HSL0927
11	1_8270STD	\\SV5\C\chem\sv5.i\092510.B\QC001
12	1_8270STD	\\sv5\c\chem\sv5.i\092510.B\HSL0925
13	1_8270STD	\\SV5\C\chem\sv5.i\092410.B\QC001
14	1_8270STD	\\SV5\C\chem\sv5.i\092410.B\HSL0924
15	1_8270STD	\\SV5\C\chem\sv5.i\092310A.B\HSL0923A
16	1_8270STD	\\SV5\C\chem\sv5.i\092310A.B\QC001
17	1_8270STD	\\SV5\C\chem\sv5.i\092310.B\QC001
18	1_8270STD	\\SV5\C\chem\sv5.i\092310.B\HSL0923
19	1_8270STD	\\SV5\C\chem\sv5.i\092210.B\HSL0922a
20	1_8270STD	\\SV5\C\chem\sv5.i\092210.B\HSL0922
21	1_8270STD	\\SV5\C\chem\sv5.i\092210.B\QC001
22	1_8270STD	\\SV5\C\chem\sv5.i\092110.B\HSL0921
23	1_8270STD	\\SV5\C\chem\sv5.i\092010.B\QC001
24	1_8270STD	\\SV5\C\chem\sv5.i\092010.B\HSL0920
25	1_8270STD	\\SV5\C\chem\sv5.i\091910a.B\HSL0919a
26	1_8270STD	\\SV5\C\chem\sv5.i\091910.B\HSL0919
27	1_8270STD	\\SV5\C\chem\sv5.i\091910.B\QC001
28	1_8270STD	\\SV5\C\chem\sv5.i\091710.B\HSL0917
29	1_8270STD	\\SV5\C\chem\sv5.i\091710.B\QC001
30	1_8270STD	\\SV5\C\chem\sv5.i\091510b.B\HSL0915b
31	1_8270STD	\\SV5\C\chem\sv5.i\091510b.B\QC003
32	1_8270STD	\\sv5\c\chem\sv5.i\091010.B\HSL0910
33	1_8270STD	\\SV5\C\chem\sv5.i\091010.B\QC001
34	1_8270STD	\\sv5\c\chem\sv5.i\090910a.B\HSL0909a
35	1_8270STD	\\SV5\C\chem\sv5.i\090910.B\HSL0909

36	1_8270STD	\\SV5\C\chem\sv5.i\090910.B\QC001
37	1_8270STD	\\SV5\C\chem\sv5.i\090810.B\HSL0908
38	1_8270STD	\\SV5\C\chem\sv5.i\090810.B\Primer
39	1_8270STD	\\sv5\c\chem\sv5.i\090710.B\HSL0907
40	1_8270STD	\\SV5\C\chem\sv5.i\090710.B\HSL0907
41	1_8270STD	\\sv5\c\chem\sv5.i\090110.B\HSL0901
42	1_8270STD	\\SV5\C\chem\sv5.i\083110.B\HSL0831
43	1_8270STD	\\sv5\c\chem\sv5.i\083010.B\QC001
44	1_8270STD	\\sv5\c\chem\sv5.i\083010.B\HSL0830
45	1_8270STD	\\SV5\C\chem\sv5.i\082710.B\QC001
46	1_8270STD	\\sv5\c\chem\sv5.i\082710.B\HSL0827
47	1_8270STD	\\SV5\C\chem\sv5.i\082610.B\HSL0826
48	1_8270STD	\\SV5\C\chem\sv5.i\082610.B\QC001
49	1_8270STD	\\SV5\C\chem\sv5.i\082510.B\QC001
50	1_8270STD	\\SV5\C\chem\sv5.i\082510.B\HSL0825
51	1_8270STD	\\sv5\c\chem\sv5.i\082310B.B\HSL0823
52	1_8270STD	\\sv5\c\chem\sv5.i\082310B.B\HSL0823E
53	1_8270STD	\\sv5\c\chem\sv5.i\082310B.B\HSL0823D
54	1_8270STD	\\SV5\C\chem\sv5.i\082310A.B\HSL0823A
55	1_8270STD	\\SV5\C\chem\sv5.i\082010.B\HSL0820
56	1_8270STD	\\sv5\c\chem\sv5.i\082010.B\QC001
57	1_8270STD	\\sv5\c\chem\sv5.i\081810A.B\HSL0818A
58	1_8270STD	\\sv5\c\chem\sv5.i\081810.B\HSL0818
59	1_8270STD	\\SV5\C\chem\sv5.i\081710.B\HSL0817D
60	1_8270STD	\\SV5\C\chem\sv5.i\081710.B\HSL0817H

TAILING FACTOR/DEGRADATION SUMMARY RESULTS

TAILING ANALYSIS SUMMARY

Compound	Tail Factor	Max Allowed	Test
Pentachlorophenol	0.6825896	5.000	PASS
Benzidine	0.6244503	3.000	PASS

DDT DEGRADATION BREAKDOWN ANALYSIS SUMMARY

Compound	Response	%Breakdown	Max Allowed	Test
4,4-DDD + DDE	189907	8.9	20.5	PASS

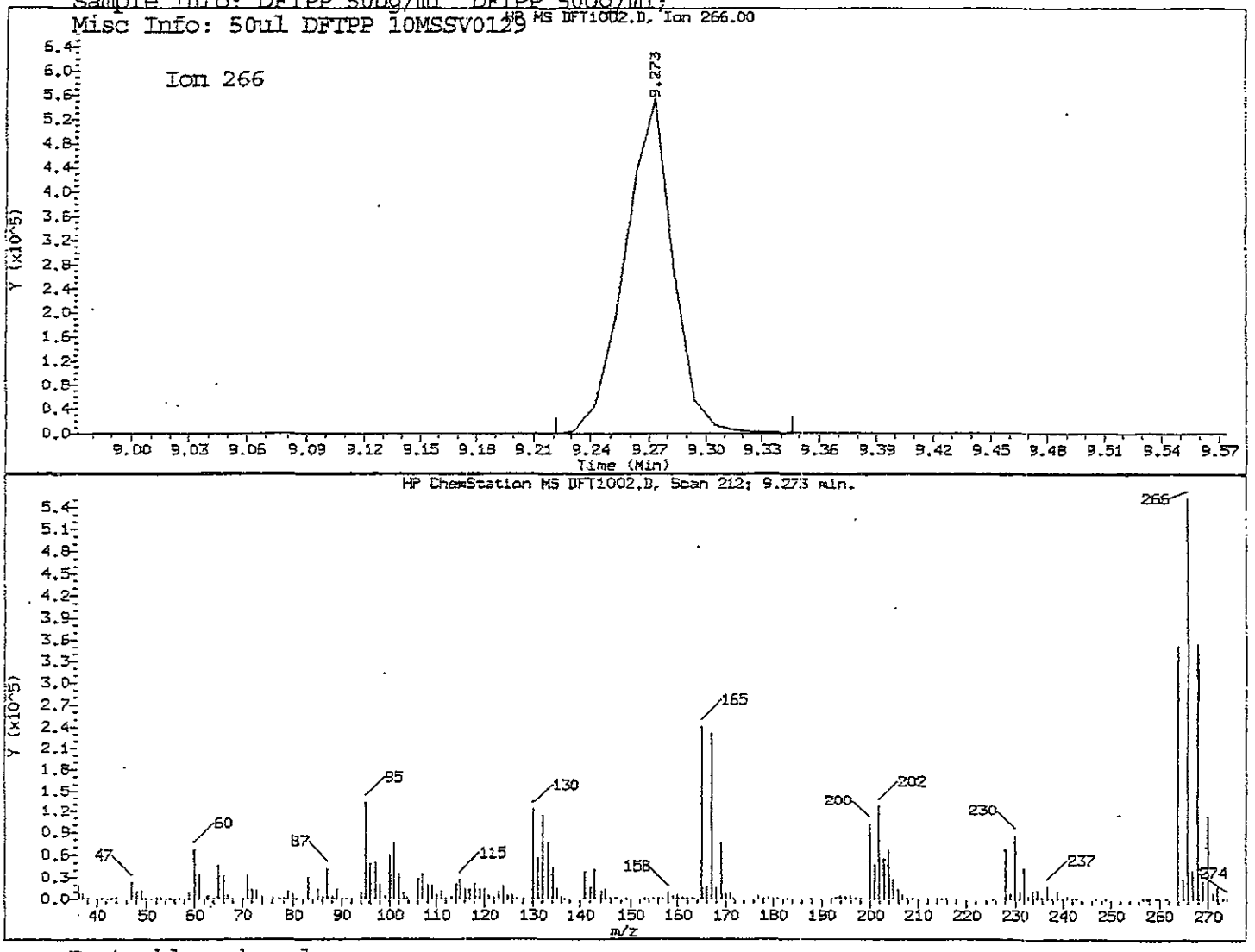
Sample //SV5/C/chem/sv5.i/100210.B/DFT1002.D/DFT1002.D

****
*** PASSED ***
****

TAILING FACTOR/DEGRADATION SAMPLE AND GRAPHIC REPORT

Report Date: 10/03/2010 11:04

Datafile Analyzed: //SV5/C/chem/sv5.i/100210.B/DFT1002.D/DFT1002.D
Method Used: \\SV5\C\chem\sv5.i\100210.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 02-OCT-2010 12:06 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129



Pentachlorophenol
==========
Exp. RT = 9.387
Found RT = 9.273

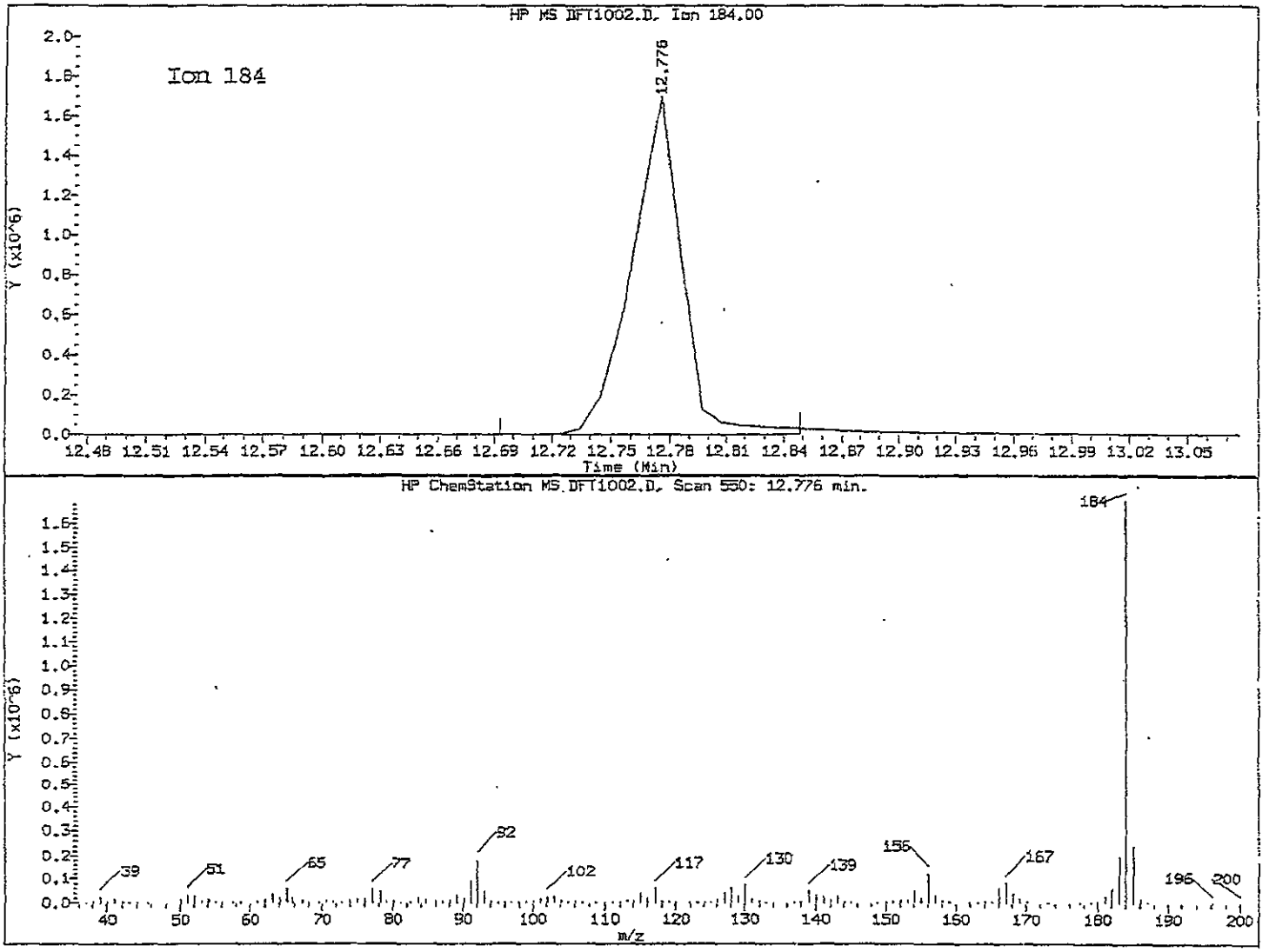
Time1 = 9.243001 Time2 = 9.273333 Time3 = 9.294038
Tailing Factor = (Time3 - Time2)/(Time2 - Time1)

Tailing factor for Pentachlorophenol OK

Tail Factor = 0.683 Maximum Allowed = 5.0

Report Date: 10/03/2010 11:04

Datafile Analyzed: //SV5/C/chem/sv5.i/100210.B/DFT1002.D/DFT1002.D
Method Used: \\SV5\C\chem\sv5.i\100210.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 02-OCT-2010 12:06 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129



Benzidine
=========
Exp. RT = 12.911
Found RT = 12.776

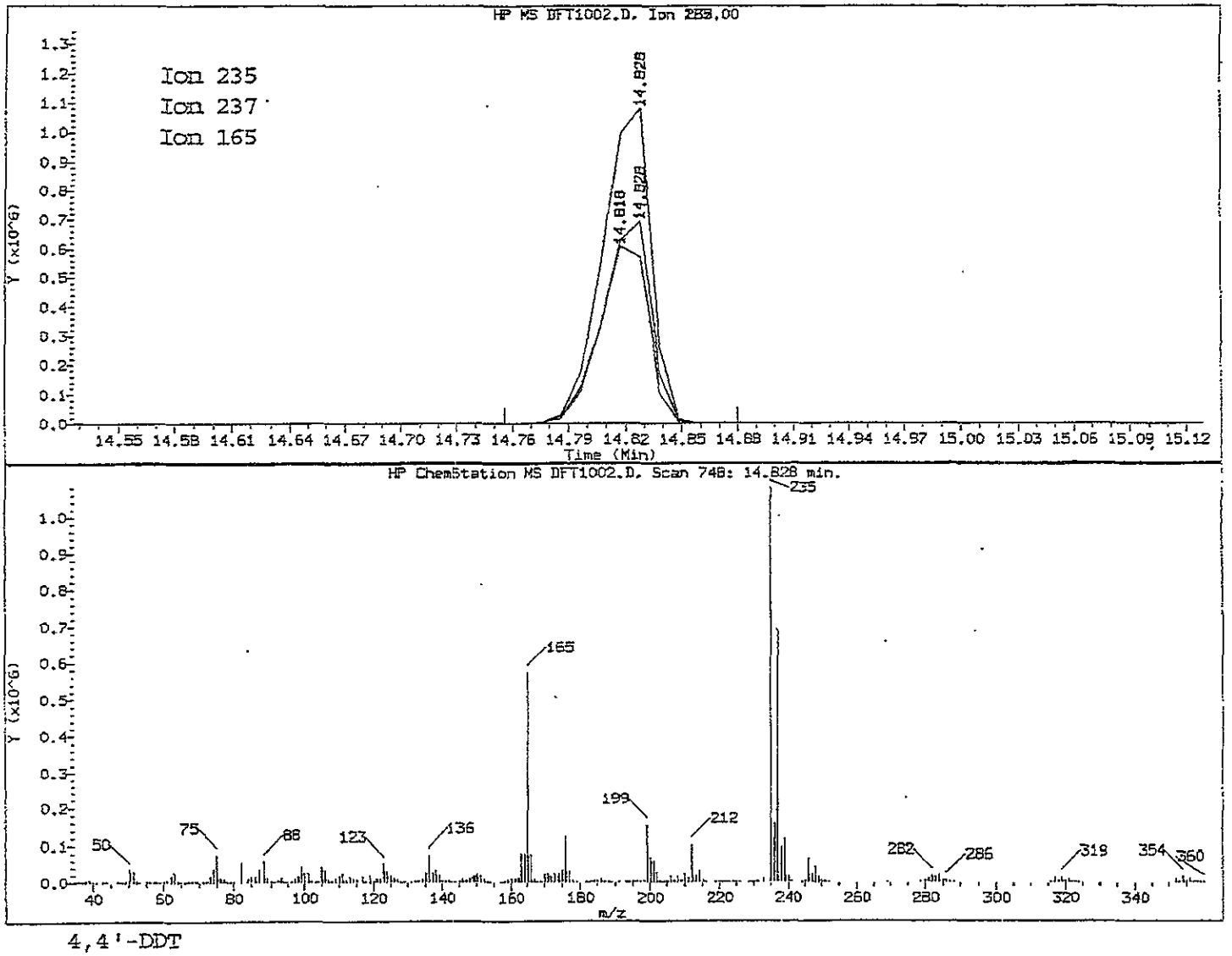
Time1 = 12.74377 Time2 = 12.77603 Time3 = 12.79618
Tailing Factor = (Time3 - Time2)/(Time2 - Time1)

Tailing factor for Benzidine OK

Tail Factor = 0.624 Maximum Allowed = 3.0

Report Date: 10/03/2010 11:04

Datafile Analyzed: //SV5/C/chem/sv5.i/100210.B/DFT1002.D/DFT1002.D
Method Used: \\SV5\C\chem\sv5.i\100210.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 02-OCT-2010 12:06 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129

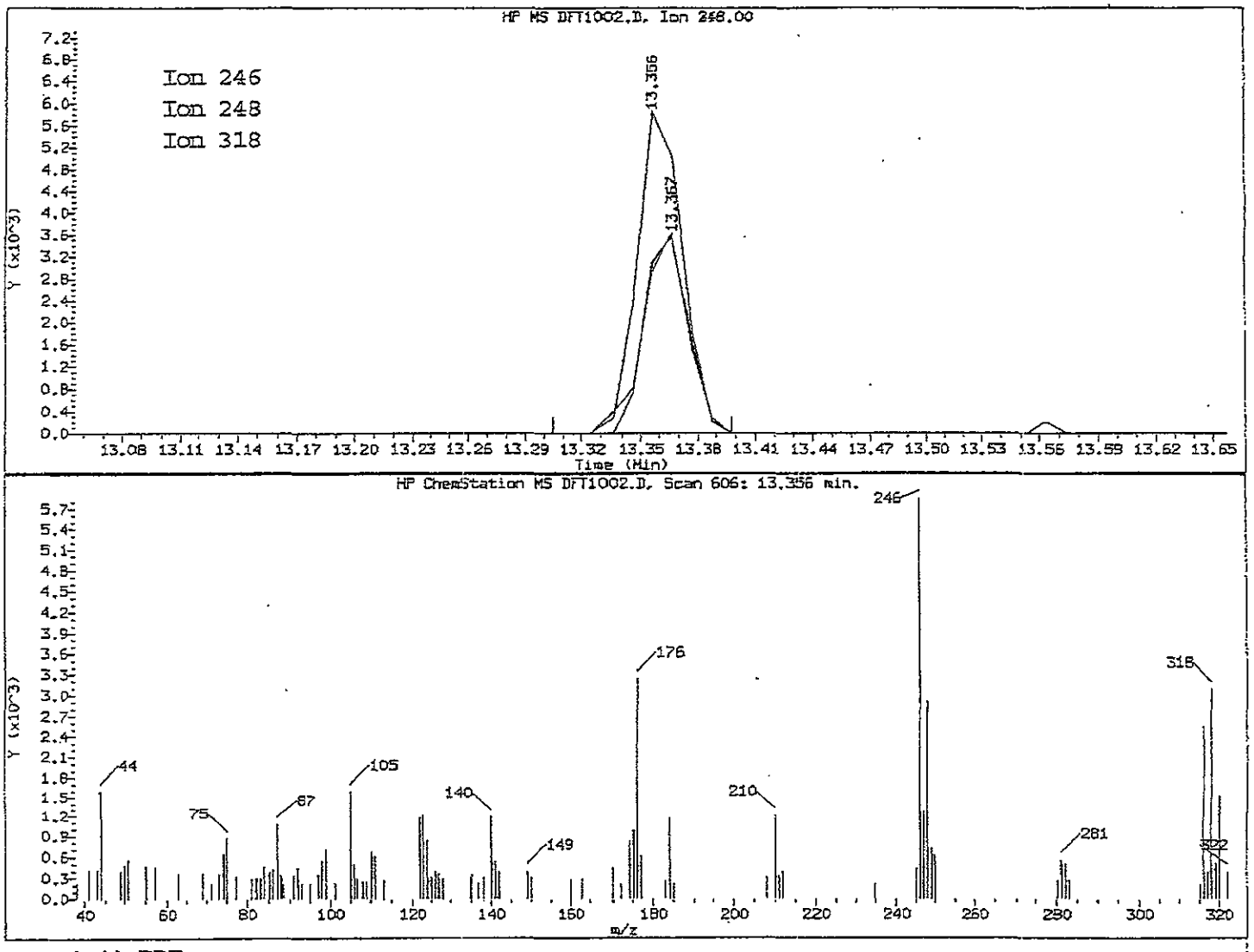


4,4'-DDT
========
Exp. RT = 14.942
Found RT = 14.828

Mass	Area	Ratio
235	1937042	100.00
237	1226081	63.30
165	1111108	57.36

Report Date: 10/03/2010 11:04

Datafile Analyzed: //SV5/C/chem/sv5.i/100210.B/DFT1002.D/DFT1002.D
Method Used: \\SV5\C\chem\sv5.i\100210.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 02-OCT-2010 12:06 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129



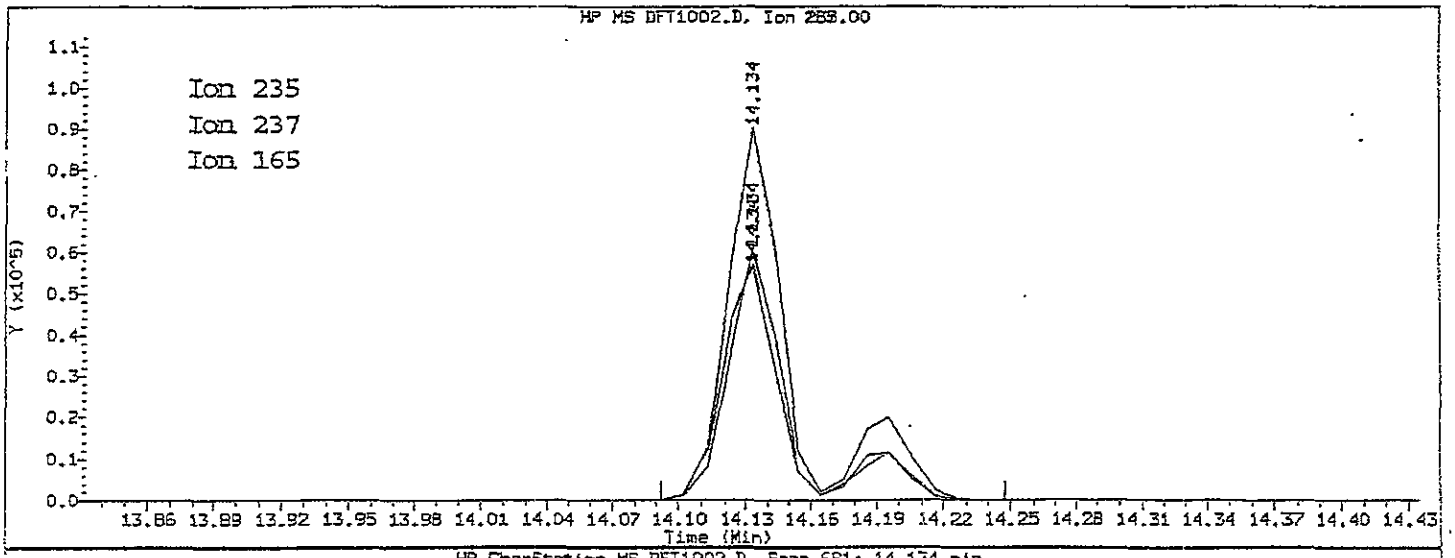
4,4'-DDE

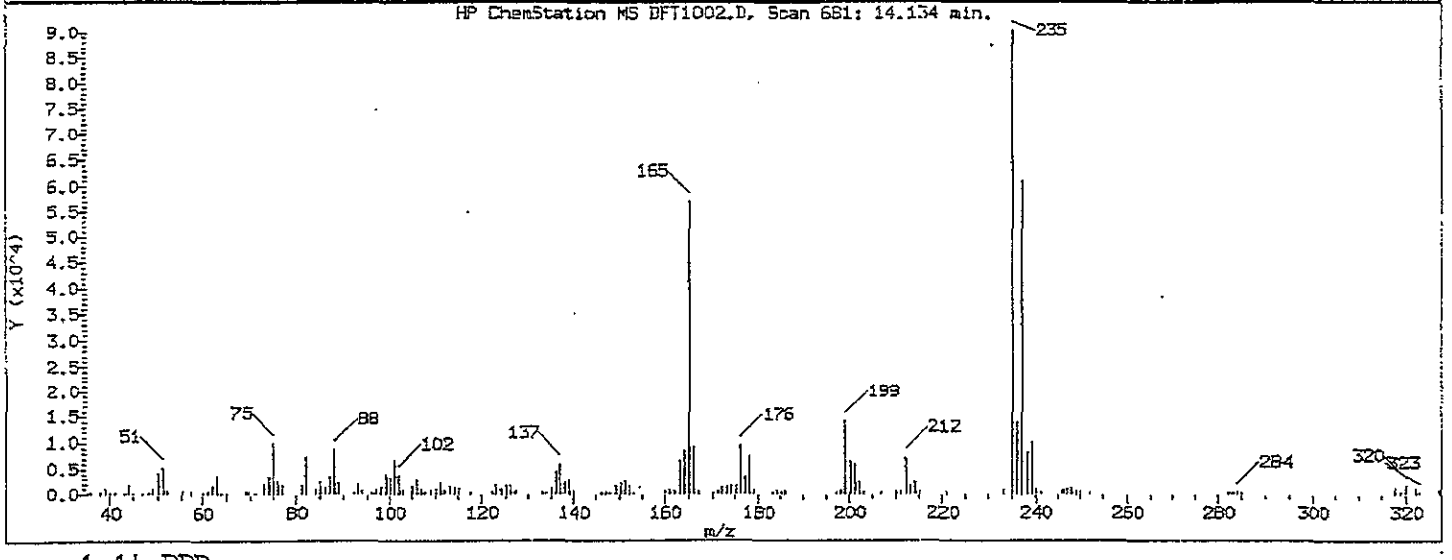
Exp. RT = 13.470
Found RT = 13.356

Mass	Area	Ratio
246	9630	100.00
248	5964	61.93
318	0	0.00

Report Date: 10/03/2010 11:04

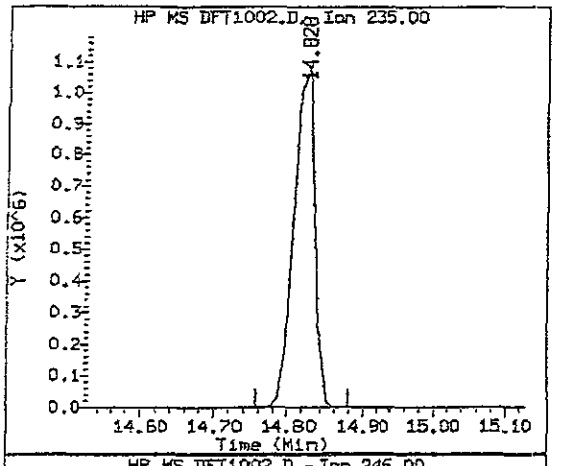
Datafile Analyzed: //SV5/C/chem/sv5.i/100210.B/DFT1002.D/DFT1002.D
Method Used: \\SV5\C\chem\sv5.i\100210.B\DFTPP.M\resol.m Inst: sv5
Injection Date: 02-OCT-2010 12:06 Operator: KT
Sample Info: DFTPP 50ug/ml DFTPP 50ug/ml;
Misc Info: 50ul DFTPP 10MSSV0129



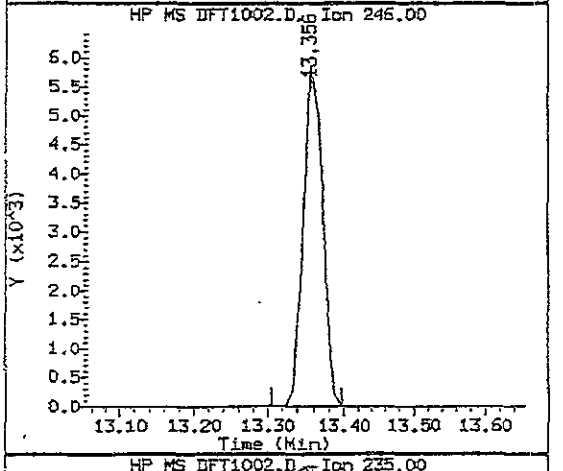


4,4'-DDD
========
Exp. RT = 14.248
Found RT = 14.134

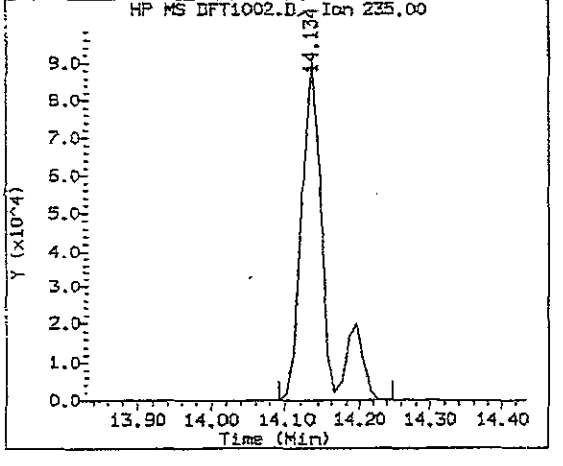
Mass	Area	Ratio
235	180277	100.00
237	115795	64.23
165	113090	62.73



Compound: 4,4'-DDT
Quant Mass: 235
RT: 14.828
Area: 1937042



Compound: 4,4'-DDE
Quant Mass: 246
RT: 13.356
Area: 9630



Compound: 4,4'-DDD
Quant Mass: 235
RT: 14.134
Area: 180277

DDT DEGRADATION BREAKDOWN ANALYSIS SUMMARY

Compound	Response	%Breakdown	Max Allowed	Test
4,4-DDD + DDE	189907	8.9	20.5	PASS

TestAmerica West Sacramento

Data file : \\SV5\C\chem\sv5.i\100210.B\DFT1002.D
Lab Smp Id: DFTPP 50ug/ml
Inj Date : 02-OCT-2010 12:06
Operator : KT Inst ID: sv5.i
Smp Info : DFTPP 50ug/ml;
Misc Info : 50ul DFTPP 10MSSV0129
Comment :
Method : \\SV5\C\chem\sv5.i\100210.B\DFTPP.m
Meth Date : 17-Aug-2010 14:10 scotts Quant Type: ISTD
Cal Date : Cal File:
Als bottle: 96 QC Sample: DFTPP
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: all.sub
Target Version: 4.14 Sample Matrix: None
Processing Host: SV5

RT	EXP RT	REL RT	MASS	RESPONSE	CONCENTRATIONS ON-COL (ug/L)	CONCENTRATIONS FINAL (ug/L)	TARGET RANGE	RATIO
1 dftpp						CAS #: 5074-71-5		
0.000	11.201	(0.000)	198	746688			0.00- 100.00	100.00
0.000	11.201	(0.000)	51	320640			30.00- 80.00	42.94
0.000	11.201	(0.000)	68	4826			0.00- 2.00	1.62
0.000	11.201	(0.000)	69	298048			0.00- 0.00	39.92
0.000	11.201	(0.000)	70	1913			0.00- 2.00	0.64
0.000	11.201	(0.000)	127	406528			25.00- 75.00	54.44
0.000	11.201	(0.000)	197	0	0.0	0.0	0.00- 1.00	0.00
0.000	11.201	(0.000)	199	49104			5.00- 9.00	6.58
0.000	11.201	(0.000)	275	170816			10.00- 30.00	22.88
0.000	11.201	(0.000)	365	20496			0.75- 0.00	2.74
0.000	11.201	(0.000)	441	100984			0.01- 99.99	74.22
0.000	11.201	(0.000)	442	702528			40.00- 110.00	94.09
0.000	11.201	(0.000)	443	136064			15.00- 24.00	19.37

Data File: \\SV5\C\chem\sv5.i\100210.B\DFT1002.D

Date : 02-OCT-2010 12:06

Client ID:

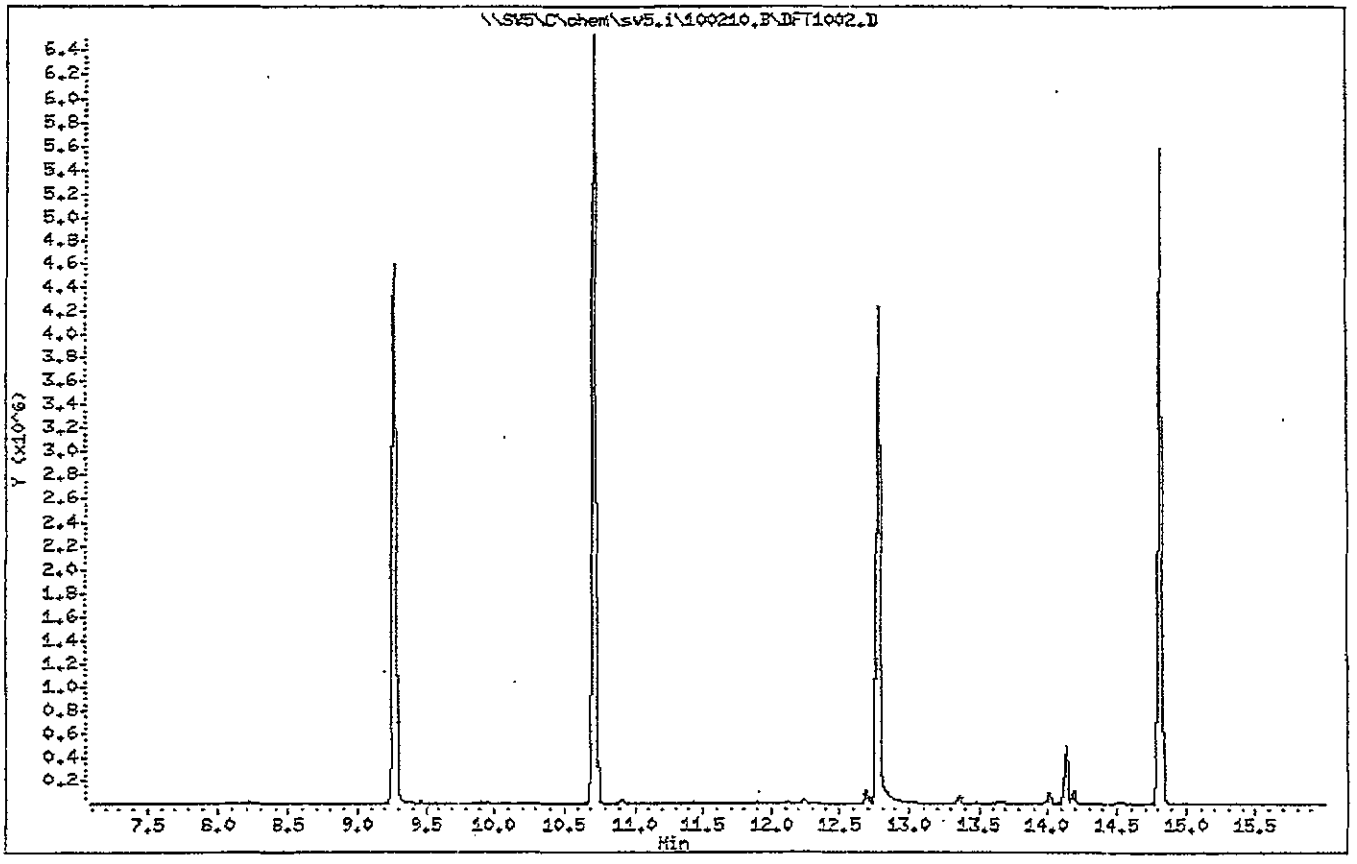
Instrument: sv5.i

Sample Info: DFTPP 50ug/ml

Operator: KT

Column phase:

Column diameter: 2.00



Data File: \\SV5\C\chem\sv5.i\100210.B\DFT1002.D

Date : 02-OCT-2010 12:06

Client ID:

Instrument: sv5.i

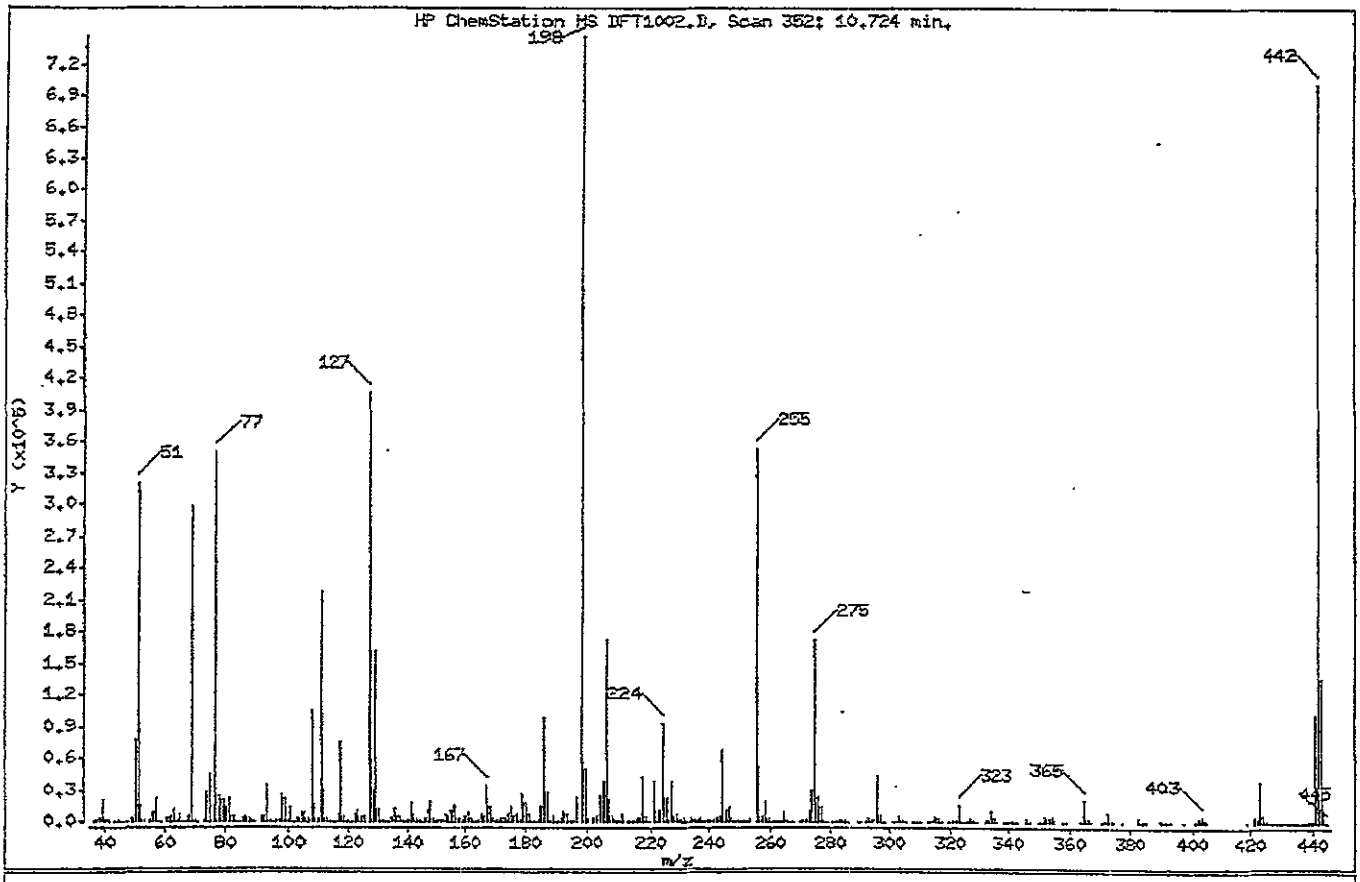
Sample Info: DFTPP 50ug/ml

Operator: KT

Column phase:

Column diameter: 2.00

1 dftpp



m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
198	Base Peak, 100% relative abundance	100.00
51	30.00 - 80.00% of mass 198	42.94
68	Less than 2.00% of mass 69	0.65 (1.62)
69	Mass 69 relative abundance	39.92
70	Less than 2.00% of mass 69	0.26 (0.64)
127	25.00 - 75.00% of mass 198	54.44
197	Less than 1.00% of mass 198	0.00
199	5.00 - 9.00% of mass 198	6.58
275	10.00 - 30.00% of mass 198	22.88
365	Greater than 0.75% of mass 198	2.74
441	Present, but less than mass 443	13.52
442	40.00 - 110.00% of mass 198	94.09
443	15.00 - 24.00% of mass 442	18.22 (19.37)

Data File: \\SV5\C\chem\sv5.i\100210.B\DFT1002.D

Date : 02-OCT-2010 12:06

Client ID:

Instrument: sv5.i

Sample Info: DFTPP 50ug/ml

Operator: KT

Column phase:

Column diameter: 2.00

Data File: DFT1002.D

Spectrum: HP ChemStation MS DFT1002.D, Scan 352: 10.724 min.

Location of Maximum: 198.00

Number of points: 340

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.10	203	130.00	12809	219.20	447	321.00	1763
37.10	1216	131.00	2287	221.00	37608	322.10	913
38.10	3314	132.00	1225	223.10	9674	323.10	16294
39.10	21392	133.00	620	224.10	93432	324.10	2245
40.00	1076	134.00	3794	225.10	21544	324.80	382
41.10	949	135.10	11378	226.10	1736	326.00	507
43.10	352	136.00	4886	227.00	37976	327.00	2789
44.00	922	137.00	5203	228.00	4945	328.00	1262
45.00	428	138.00	1265	229.00	7548	329.10	343
47.00	204	139.00	791	230.00	1024	331.90	894
49.10	2676	140.00	2233	231.10	2757	333.00	1455
50.10	77024	141.00	17480	232.00	528	334.10	9590
51.10	320640	142.00	7259	233.00	641	335.00	2774
52.10	16189	143.00	3921	234.00	2909	336.00	291
53.10	963	144.00	1378	235.00	2419	339.00	369
55.00	1815	145.10	829	236.10	1608	340.00	399
56.00	8872	146.00	3251	237.00	3192	341.00	2042
57.00	22504	147.00	9463	238.00	581	342.10	852
58.00	755	148.00	18744	239.00	1185	343.20	220
59.10	372	149.00	4031	240.00	1065	346.00	2819
61.00	3888	150.10	1094	241.00	1870	346.90	608
62.00	4800	151.20	2277	242.00	3682	350.30	205
63.10	11199	152.10	1506	243.10	4924	351.00	283
64.10	1448	153.00	6113	244.10	66488	352.00	5049
65.10	6509	154.00	5445	245.10	9865	353.10	3110
66.00	498	155.00	10151	246.00	14573	354.00	5432
67.10	461	156.10	14866	247.00	3022	355.00	1087
68.00	4826	157.10	3676	248.10	618	358.00	241
69.00	298048	158.10	3734	249.00	2441	359.00	574
70.10	1913	159.00	2313	250.00	627	363.50	249
71.10	410	160.00	5246	250.90	1000	365.00	20496
73.10	2021	161.10	8666	252.00	756	366.00	3166
74.00	28000	162.00	2863	253.10	2603	367.00	225
75.00	45304	163.10	562	255.00	353024	370.10	477
76.10	15795	164.00	1067	256.00	51440	370.90	1541

Data File: \\SV5\C\chem\sv5.i\100210.B\DFT1002.D

Date : 02-OCT-2010 12:06

Client ID:

Instrument: sv5.i

Sample Info: DFTPP 50ug/ml;

Operator: KT

Column phase:

Column diameter: 2.00

Data File: DFT1002.D

Spectrum: HP ChemStation MS DFT1002.D, Scan 352: 10.724 min.

Location of Maximum: 198.00

Number of points: 340

m/z	Y	m/z	Y	m/z	Y	m/z	Y
77.10	349952	165.00	6962	257.00	4474	372.10	8489
78.10	23464	166.00	5717	258.00	19504	373.10	1814
79.00	20048	167.00	33648	259.10	3095	373.80	348
80.00	14146	168.00	13682	260.00	645	377.10	263
81.00	22008	169.00	2802	261.10	797	383.00	2624
82.00	5822	170.00	1014	262.20	249	383.90	598
83.00	5093	171.00	1339	263.00	269	385.00	289
84.00	614	172.00	3224	264.10	532	390.00	1367
85.00	3848	173.00	4109	265.00	7904	391.00	754
86.00	5985	174.00	7189	266.00	1181	392.10	664
87.00	2652	175.10	13638	267.20	204	393.20	281
88.00	1078	176.10	4293	267.60	232	397.00	230
89.00	472	177.00	6577	270.00	489	400.90	335
91.00	5074	178.10	1572	271.00	901	402.00	3464
92.00	5292	179.00	25912	272.10	1129	403.00	5568
93.00	34848	180.00	16984	273.00	10963	404.10	1777
94.00	2386	181.00	7182	274.00	30032	405.00	292
95.00	749	182.00	1363	275.00	170816	418.90	259
96.00	1660	183.00	559	276.10	22944	421.00	5400
97.10	1007	184.10	2227	277.00	13493	422.00	4183
98.00	25944	185.10	13301	278.10	2251	423.00	37592
99.00	21688	186.00	97584	279.00	648	424.00	6802
100.00	1844	187.10	27792	281.10	266	425.00	930
101.00	13609	188.10	2556	282.00	217	426.50	251
102.10	646	189.00	5094	283.00	1957	427.30	338
103.00	3748	189.90	756	284.00	1097	428.40	200
104.00	8390	191.10	2995	285.10	2569	429.20	300
105.00	8359	192.00	7909	286.10	444	430.20	272
106.10	3007	193.00	7605	289.00	691	431.10	404
107.00	104896	194.10	1998	290.10	589	431.50	324
108.00	17616	195.10	1321	292.10	763	432.20	298
109.00	3545	196.00	22448	293.00	3141	432.50	326
110.00	218112	198.00	746688	294.10	1275	433.30	317
111.00	30736	199.00	49104	296.00	42616	433.70	342
112.00	4281	200.00	4038	297.00	6196	434.30	362

Data File: \\SV5\C\chem\sv5.i\100210.B\DFT1002.D

Date : 02-OCT-2010 12:06

Client ID:

Instrument: sv5.i

Sample Info: DFTPP 50ug/ml

Operator: KT

Column phase:

Column diameter: 2.00

Data File: DFT1002.D

Spectrum: HP ChemStation MS DFT1002.D, Scan 352: 10.724 min.

Location of Maximum: 198.00

Number of points: 340

m/z	Y	m/z	Y	m/z	Y	m/z	Y
113.00	1310	201.60	4029	298.00	465	434.90	650
114.40	467	203.00	4788	301.00	504	435.90	530
115.00	646	204.00	23416	302.00	695	436.50	586
116.10	6327	205.00	38288	303.10	5810	436.90	846
117.00	75520	206.10	172352	304.00	2035	437.50	628
118.00	5507	207.10	21328	305.10	290	438.20	1136
119.00	839	208.00	5487	308.00	764	439.30	1287
120.10	1180	209.00	2186	309.10	446	441.00	100934
121.00	807	210.00	2002	310.00	839	442.00	702528
122.00	6408	211.10	7473	312.20	271	443.00	136064
123.00	10302	213.00	410	312.90	292	444.00	12344
124.00	4600	214.10	372	314.00	2431	445.10	689
125.00	4447	215.10	1837	315.00	5363		
127.00	406528	216.00	3226	316.00	2900		
128.00	28392	217.00	41648	317.10	363		
129.00	161024	218.00	5388	319.80	287		

TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002A.D
Lab Smp Id: HSL_005 ug/ml CS-1 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 12:27
Operator : KT Inst ID: sv5.i
Smp Info : HSL_005 ug/ml CS-1;1;;1;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0307;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:09 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 1 Calibration Sample, Level: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SACP307UM

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
*	1 1,4-Dichlorobenzene-d4	152	3.955	3.955	(1.000)	141539	40.0000		(Q)
*	2 Naphthalene-d8	136	5.374	5.374	(1.000)	605687	40.0000		
*	3 Acenaphthene-d10	164	7.468	7.468	(1.000)	321839	40.0000		
*	4 Phenanthrene-d10	188	9.406	9.405	(1.000)	496356	40.0000		
*	5 Chrysene-d12	240	13.779	13.779	(1.000)	453007	40.0000		
*	6 Perylene-d12	264	16.162	16.162	(1.000)	445119	40.0000		
$	7 2-Fluorophenol	112	2.742	2.732	(0.693)	25566	5.00000	5.124	
$	8 Phenol-d5	99	3.613	3.613	(0.914)	30471	5.00000	4.857	
$	9 2-Chlorophenol-d4	132	3.758	3.758	(0.950)	26144	5.00000	4.745	
$	10 1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	16945	5.00000	4.861	
$	11 Nitrobenzene-d5	82	4.576	4.576	(0.852)	25006	5.00000	4.874	(M)
$	12 2-Fluorobiphenyl	172	6.680	6.680	(0.895)	51695	5.00000	4.986	
$	13 2,4,6-Tribromophenol	330	8.473	8.473	(1.135)	6048	5.00000	4.325	
$	14 Terphenyl-d14	244	12.017	12.017	(0.872)	44456	5.00000	4.982	
	15 N-Nitrosodimethylamine	74	1.716	1.706	(0.434)	16436	5.00000	5.040	(q)
	16 Pyridine	79	1.737	1.726	(0.439)	29567	5.00000	5.422	(q)
	23 Aniline	93	3.654	3.654	(0.924)	39064	5.00000	4.892	(Q)
	24 Phenol	94	3.623	3.623	(0.916)	36112	5.00000	5.009	(Q)
	26 Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	26067	5.00000	5.157	
	27 2-Chlorophenol	128	3.768	3.768	(0.953)	26910	5.00000	4.863	
	28 1,3-Dichlorobenzene	146	3.923	3.923	(0.992)	29883	5.00000	4.958	
	29 1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	31337	5.00000	4.972	
	30 Benzyl Alcohol	108	4.120	4.120	(1.042)	17983	5.00000	4.835	
	31 1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	28663	5.00000	4.947	
	32 2-Methylphenol	108	4.255	4.255	(1.076)	24914	5.00000	4.923	
	33 2,2'-oxybis(1-Chloropropane)	45	4.297	4.297	(1.086)	40622	5.00000	5.049	
	34 4-Methylphenol	108	4.421	4.421	(1.118)	26292	5.00000	4.891	
	36 Hexachloroethane	117	4.504	4.504	(1.139)	10779	5.00000	5.024	
	37 N-Nitrosodipropylamine	70	4.442	4.442	(1.123)	16719	5.00000	4.670	
	42 Nitrobenzene	77	4.597	4.597	(0.855)	24875	5.00000	4.960	
	44 Isophorone	82	4.856	4.856	(0.904)	48024	5.00000	4.980	
	45 2-Nitrophenol	139	4.960	4.960	(0.923)	14088	5.00000	4.735	
	46 2,4-Dimethyphenol	107	5.012	5.012	(0.933)	26089	5.00000	4.935	

10-7-10

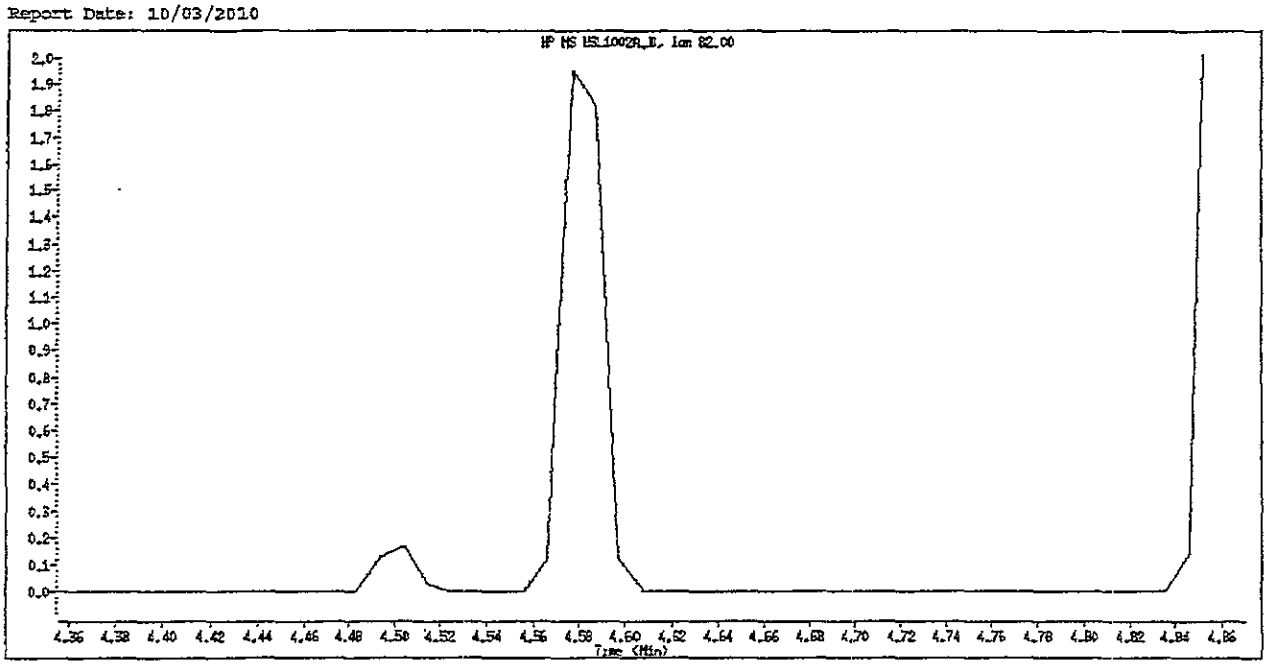
Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)	
47 Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.954)	31152	5.00000	5.288	
49 2,4-Dichlorophenol	162	5.229	5.229	(0.973)	19256	5.00000	4.708	
50 Benzoic Acid	122	5.084	5.115	(0.946)	12679	5.00000	4.333	
51 1,2,4-Trichlorobenzene	180	5.322	5.322	(0.990)	22282	5.00000	5.032	
52 Naphthalene	128	5.395	5.395	(1.004)	83236	5.00000	4.977	
54 4-Chloroaniline	127	5.488	5.488	(1.021)	30853	5.00000	4.707	
57 Hexachlorobutadiene	225	5.613	5.613	(1.044)	10823	5.00000	4.994	
60 4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	22205	5.00000	4.862	
63 2-Methylnaphthalene	142	6.203	6.203	(1.154)	51849	5.00000	4.936	
66 Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	10813	5.00000	4.503	
69 2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	12546	5.00000	4.886	
70 2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	12400	5.00000	4.483	
71 2-Chloronaphthalene	162	6.784	6.784	(0.908)	45713	5.00000	5.047	
73 2-Nitroaniline	65	6.949	6.949	(0.931)	12703	5.00000	4.627	
76 Dimethylphthalate	163	7.219	7.229	(0.967)	49639	5.00000	4.760	
77 Acenaphthylene	152	7.281	7.281	(0.975)	75041	5.00000	4.758	
79 2,6-Dinitrotoluene	165	7.291	7.302	(0.976)	11404	5.00000	4.694 (QM)	
80 3-Nitroaniline	138	7.447	7.447	(0.997)	14226	5.00000	4.691 (Q)	
81 Acenaphthene	153	7.509	7.509	(1.006)	50639	5.00000	5.044	
82 2,4-Dinitrophenol	184	7.571	7.572	(1.014)	4083	5.00000	6.945 (q)	
83 Dibenzofuran	168	7.696	7.706	(1.031)	63477	5.00000	4.764	
84 4-Nitrophenol	109	7.675	7.675	(1.028)	5114	5.00000	4.065 (Q)	
86 2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	13823	5.00000	4.335 (q)	
91 Fluorene	166	8.131	8.131	(1.089)	54136	5.00000	4.906	
92 Diethylphthalate	149	8.100	8.100	(1.085)	49177	5.00000	4.606	
93 4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	22112	5.00000	4.820	
94 4-Nitroaniline	138	8.214	8.214	(1.100)	13415	5.00000	4.463	
97 4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	5780	5.00000	7.325 (q)	
98 N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	41998	5.86000	5.582	
100 Azobenzene	77	8.348	8.348	(0.888)	48101	5.00000	4.928	
101 4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	11766	5.00000	4.856	
108 Hexachlorobenzene	284	8.981	8.981	(0.955)	14244	5.00000	5.264	
110 Pentachlorophenol	266	9.240	9.240	(0.982)	5849	5.00000	7.264	
114 Phenanthrene	178	9.437	9.437	(1.003)	80873	5.00000	5.169	
115 Anthracene	178	9.499	9.499	(1.010)	77577	5.00000	4.963	
118 Carbazole	167	9.768	9.768	(1.039)	70241	5.00000	4.920	
120 Di-n-Butylphthalate	149	10.463	10.463	(1.112)	78722	5.00000	4.641	
126 Fluoranthene	202	11.302	11.302	(1.202)	64427	5.00000	4.596	
127 Benzidine	184	11.571	11.571	(0.840)	44267	5.00000	4.822	
128 Pyrene	202	11.665	11.665	(0.847)	71230	5.00000	5.030	
134 3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	37074	5.00000	4.574	
136 Butylbenzylphthalate	149	12.991	12.991	(0.943)	36798	5.00000	5.185	
138 Benzo(a)Anthracene	228	13.758	13.758	(0.998)	62384	5.00000	5.170	
139 Chrysene	228	13.820	13.831	(1.003)	59618	5.00000	4.830	
140 3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	22168	5.00000	4.870	
141 bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.024)	51997	5.00000	5.319	
142 Di-n-octylphthalate	149	15.157	15.167	(1.100)	76353	5.00000	4.886	
144 Benzo(b)fluoranthene	252	15.572	15.582	(0.963)	45075	5.00000	4.473 (Q)	
145 Benzo(k)fluoranthene	252	15.613	15.623	(0.966)	68403	5.00000	5.288 (q)	
147 Benzo(e)pyrene	252	15.996	16.007	(0.990)	50295	5.00000	4.786	
148 Benzo(a)pyrene	252	16.069	16.079	(0.994)	54694	5.00000	4.788	
151 Indeno(1,2,3-cd)pyrene	276	17.789	17.800	(1.101)	41053	5.00000	4.443	
152 Dibenzo(a,h)anthracene	278	17.841	17.841	(1.104)	49018	5.00000	4.749	
153 Benzo(g,h,i)perylene	276	18.224	18.235	(1.128)	53428	5.00000	4.781	

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				113478	5.00000	4.931(A)

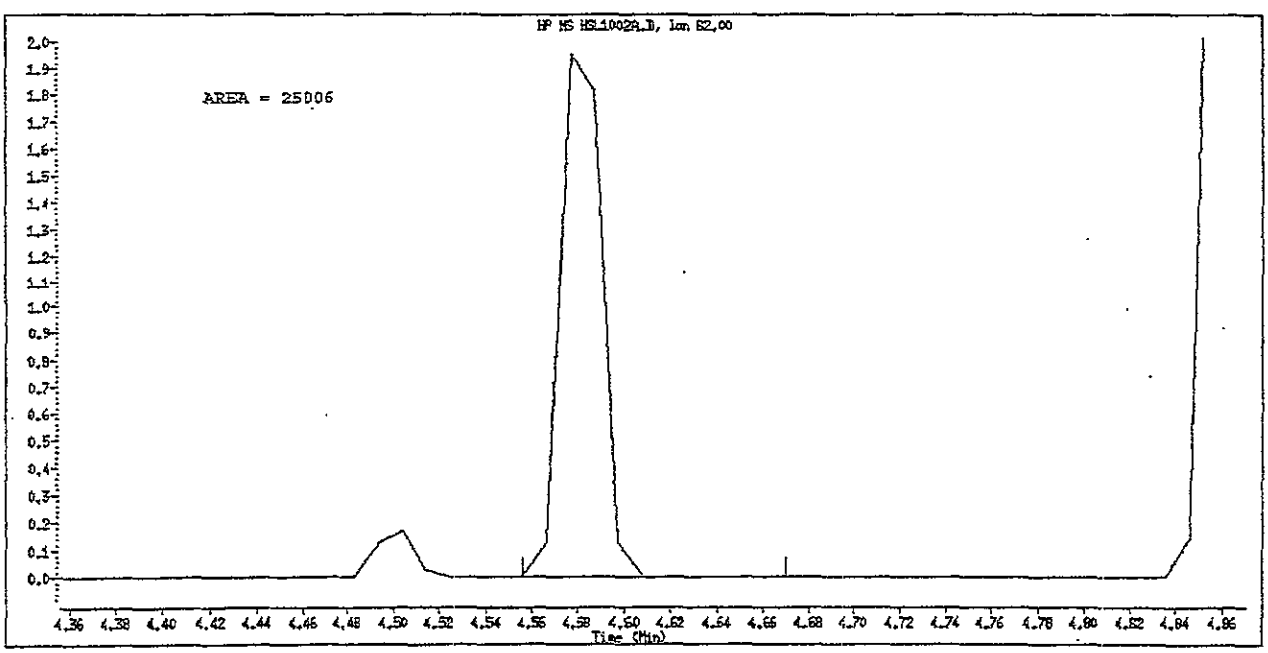
QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
M - Compound response manually integrated.
q - Qualifier signal exceeded ratio warning limit.

Data File Name: HSL1002A.D
Inj. Date and Time: 02-OCT-2010 12:27
Instrument ID: sv5.i
Client ID: 8270F.M
Compound Name: Nitrobenzene-d5
CAS #: 4165-60-0
Report Date: 10/03/2010



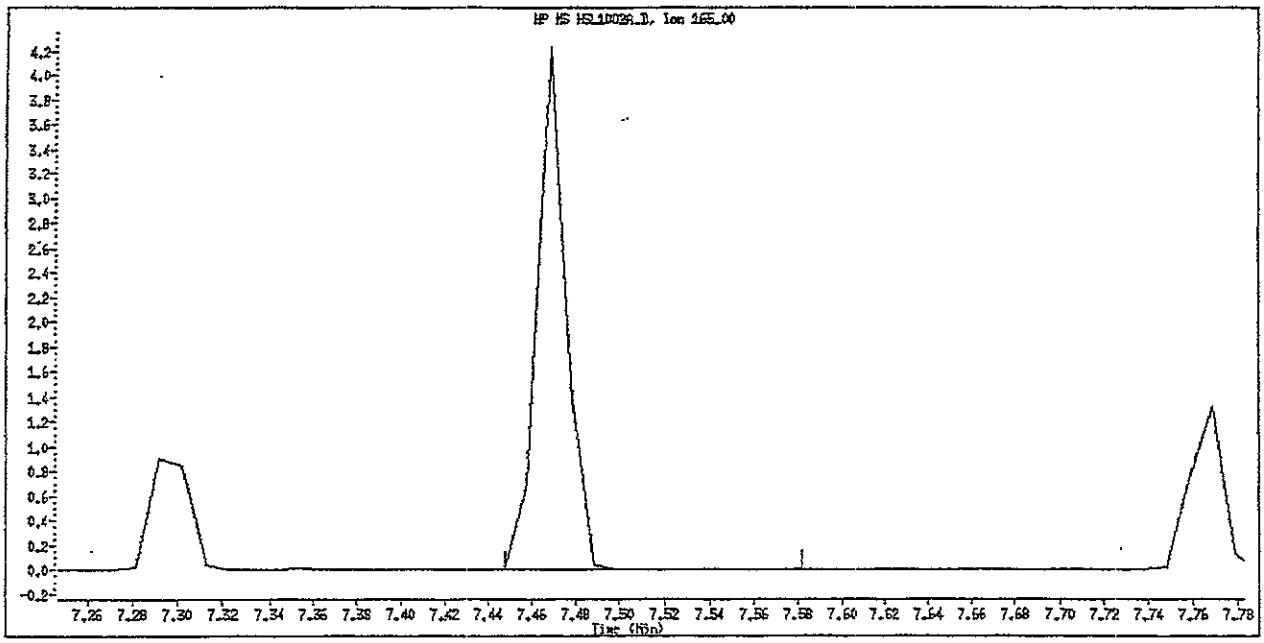
Original Integration



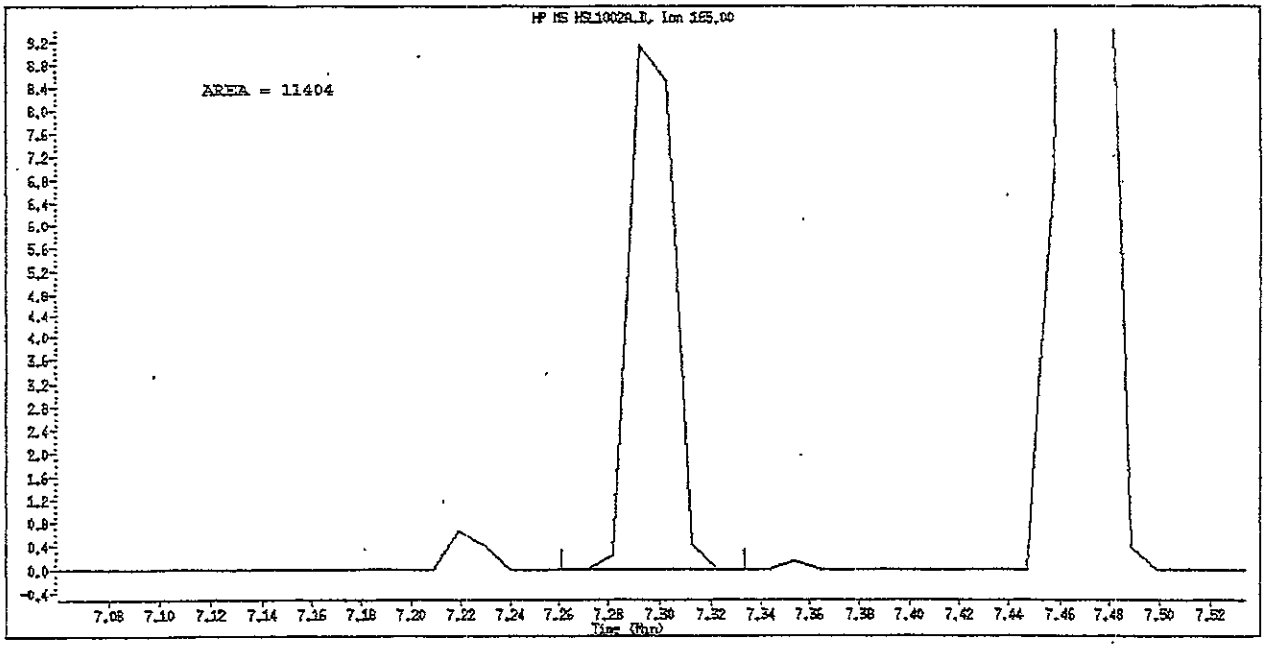
Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Peak Not Found

Data File Name: HSL1002A.D
Inj. Date and Time: 02-OCT-2010 12:27
Instrument ID: sv5.i
Client ID: 8270P.M
Compound Name: 2,6-Dinitrotoluene
CAS #: 606-20-2
Report Date: 10/03/2010



Original Integration



Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Wrong Peak

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\100210.B\HSL1002A.D
Lab Smp Id: HSL_005 ug/ml CS-1 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 12:27
Operator : KT Inst ID: sv5.i
Smp Info : HSL_005 ug/ml CS-1;1;;1;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0307;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Meth Date : 02-Oct-2010 16:57 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 1 Calibration Sample, Level: 1
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
*	1 1,4-Dichlorobenzene-d4	152	3.955	3.955	(1.000)	141539	40.0000		(Q)
*	2 Naphthalene-d8	136	5.374	5.374	(1.000)	605687	40.0000		
*	3 Acenaphthene-d10	164	7.468	7.468	(1.000)	321839	40.0000		
*	4 Phenanthrene-d10	188	9.406	9.405	(1.000)	496356	40.0000		
*	5 Chrysene-d12	240	13.779	13.779	(1.000)	453007	40.0000		
*	6 Perylene-d12	264	16.162	16.162	(1.000)	445119	40.0000		
$	7 2-Fluorophenol	112	2.742	2.732	(0.693)	25566	5.00000	4.894	
$	8 Phenol-d5	99	3.613	3.613	(0.914)	30471	5.00000	4.587	
$	9 2-Chlorophenol-d4	132	3.758	3.758	(0.950)	26144	5.00000	4.616	
$	10 1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	16965	5.00000	4.793	
$	11 Nitrobenzene-d5	82	Compound Not Detected.						
$	12 2-Fluorobiphenyl	172	6.680	6.680	(0.895)	51695	5.00000	5.015	
$	13 2,4,6-Tribromophenol	330	8.473	8.473	(1.135)	6048	5.00000	4.760	
$	14 Terphenyl-d14	244	12.017	12.017	(0.872)	44456	5.00000	5.032	
	15 N-Nitrosodimethylamine	74	1.716	1.706	(0.436)	16436	5.00000	4.767	(q)
	16 Pyridine	79	1.737	1.726	(0.439)	29567	5.00000	5.146	
	23 Aniline	93	3.654	3.654	(0.924)	39064	5.00000	4.689	(Q)
	24 Phenol	94	3.623	3.623	(0.916)	36112	5.00000	5.111	(Q)
	26 Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	26057	5.00000	4.856	
	27 2-Chlorophenol	128	3.768	3.768	(0.953)	26910	5.00000	4.813	
	28 1,3-Dichlorobenzene	146	3.923	3.923	(0.992)	29883	5.00000	4.837	
	29 1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	31337	5.00000	5.017	
	30 Benzyl Alcohol	108	4.120	4.120	(1.042)	17983	5.00000	4.681	
	31 1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	28663	5.00000	4.842	
	32 2-Methylphenol	108	4.255	4.255	(1.076)	24914	5.00000	4.770	
	33 2,2'-oxybis(1-Chloropropane)	45	4.297	4.297	(1.086)	40622	5.00000	4.077	
	34 4-Methylphenol	108	4.421	4.421	(1.118)	26292	5.00000	4.723	
	36 Hexachloroethane	117	4.504	4.504	(1.139)	10779	5.00000	4.891	
	37 N-Nitrosodinpropylamine	70	4.442	4.442	(1.123)	16719	5.00000	4.290	
	42 Nitrobenzene	77	4.597	4.597	(0.855)	24875	5.00000	4.659	
	44 Isophorone	82	4.856	4.856	(0.904)	48024	5.00000	4.744	
	45 2-Nitrophenol	139	4.960	4.960	(0.923)	14088	5.00000	4.833	
	46 2,4-Dimethyphenol	107	5.012	5.012	(0.933)	26089	5.00000	4.820	

	Compounds	QUANT STG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.954)	31152	5.00000	5.169
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	19256	5.00000	4.834
50	Benzoic Acid	122	5.084	5.115	(0.946)	12679	5.00000	4.202
51	1,2,4-Trichlorobenzene	180	5.322	5.322	(0.990)	22282	5.00000	5.160
52	Naphthalene	128	5.395	5.395	(1.004)	83236	5.00000	4.937
54	4-Chloroaniline	127	5.488	5.488	(1.021)	30853	5.00000	4.652
57	Hexachlorobutadiene	225	5.613	5.613	(1.044)	10823	5.00000	5.267
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	22205	5.00000	4.844
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	51849	5.00000	5.040
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	10813	5.00000	4.405
69	2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	12546	5.00000	5.149
70	2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	12400	5.00000	4.633
71	2-Chloronaphthalene	162	6.784	6.784	(0.908)	45713	5.00000	5.066
73	2-Nitroaniline	65	6.949	6.949	(0.931)	12703	5.00000	4.204
76	Dimethylphthalate	163	7.219	7.229	(0.967)	49639	5.00000	4.763
77	Acenaphthylene	152	7.281	7.281	(0.975)	75041	5.00000	4.757
79	2,6-Dinitrotoluene	165	7.468	7.302	(1.000)	39415	5.00000	16.89 (Q)
80	3-Nitroaniline	138	7.447	7.447	(0.997)	14226	5.00000	4.597 (Q)
81	Acenaphthene	153	7.509	7.509	(1.006)	50639	5.00000	5.038
82	2,4-Dinitrophenol	184	7.571	7.571	(1.014)	4083	5.00000	5.740 (q)
83	Dibenzofuran	168	7.696	7.706	(1.031)	63477	5.00000	4.780
84	4-Nitrophenol	109	7.675	7.675	(1.028)	5114	5.00000	3.785 (Q)
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	13823	5.00000	4.422 (q)
91	Fluorene	166	8.131	8.131	(1.089)	54136	5.00000	4.976
92	Diethylphthalate	149	8.100	8.100	(1.085)	49177	5.00000	4.514
93	4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	22112	5.00000	4.930
94	4-Nitroaniline	138	8.214	8.214	(1.100)	13415	5.00000	4.635
97	4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	5780	5.00000	8.076 (q)
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	41998	5.86000	5.430
100	Azobenzene	77	8.348	8.348	(0.888)	48101	5.00000	4.470
101	4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	11766	5.00000	4.905
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	14244	5.00000	5.498
110	Pentachlorophenol	266	9.240	9.240	(0.982)	5849	5.00000	3.762
114	Phenanthrene	178	9.437	9.437	(1.003)	80873	5.00000	5.224
115	Anthracene	178	9.499	9.499	(1.010)	77577	5.00000	4.979
118	Carbazole	167	9.768	9.768	(1.039)	70241	5.00000	4.847
120	Di-n-Butylphthalate	149	10.463	10.463	(1.112)	79722	5.00000	4.549
126	Fluoranthene	202	11.302	11.302	(1.202)	64427	5.00000	4.624
127	Benzidine	184	11.571	11.571	(0.840)	44267	5.00000	4.759
128	Pyrene	202	11.665	11.665	(0.847)	71230	5.00000	5.029
134	3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	37074	5.00000	4.644
136	Butylbenzylphthalate	149	12.991	12.991	(0.943)	36798	5.00000	5.084
138	Benzo(a)Anthracene	228	13.758	13.758	(0.998)	62384	5.00000	5.220
139	Chrysene	228	13.820	13.831	(1.003)	59618	5.00000	4.801
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	22168	5.00000	5.069
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.024)	51997	5.00000	5.218
142	Di-n-octylphthalate	149	15.157	15.167	(1.100)	76353	5.00000	4.752
144	Benzo(b)fluoranthene	252	15.572	15.582	(0.963)	45075	5.00000	4.270 (Q)
145	Benzo(k)fluoranthene	252	15.613	15.623	(0.966)	68403	5.00000	5.546 (q)
147	Benzo(e)pyrene	252	15.996	16.007	(0.990)	50295	5.00000	4.807
148	Benzo(a)pyrene	252	16.069	16.079	(0.994)	54694	5.00000	4.761
151	Indeno(1,2,3-cd)pyrene	276	17.789	17.800	(1.101)	41053	5.00000	4.039
152	Dibenzo(a,h)anthracene	278	17.841	17.841	(1.104)	49018	5.00000	4.706
153	Benzo(g,h,i)perylene	276	18.224	18.235	(1.128)	53428	5.00000	4.784

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				113478	5.00000	4.958 (A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002A.D
Lab Smp Id: HSL_005 ug/ml CS-1
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\sv5\c\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0307;0;8270F.M

Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	141539	15.42
2 Naphthalene-d8	530514	265257	1061028	605687	14.17
3 Acenaphthene-d10	282538	141269	565076	321839	13.91
4 Phenanthrene-d10	462722	231361	925444	496356	7.27
5 Chrysene-d12	435850	217925	871700	453007	3.94
6 Perylene-d12	422284	211142	844568	445119	5.41

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.96	3.46	4.46	3.96	0.00
2 Naphthalene-d8	5.37	4.87	5.87	5.37	0.00
3 Acenaphthene-d10	7.47	6.97	7.97	7.47	0.00
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	0.00
5 Chrysene-d12	13.78	13.28	14.28	13.78	0.00
6 Perylene-d12	16.16	15.66	16.66	16.16	0.00

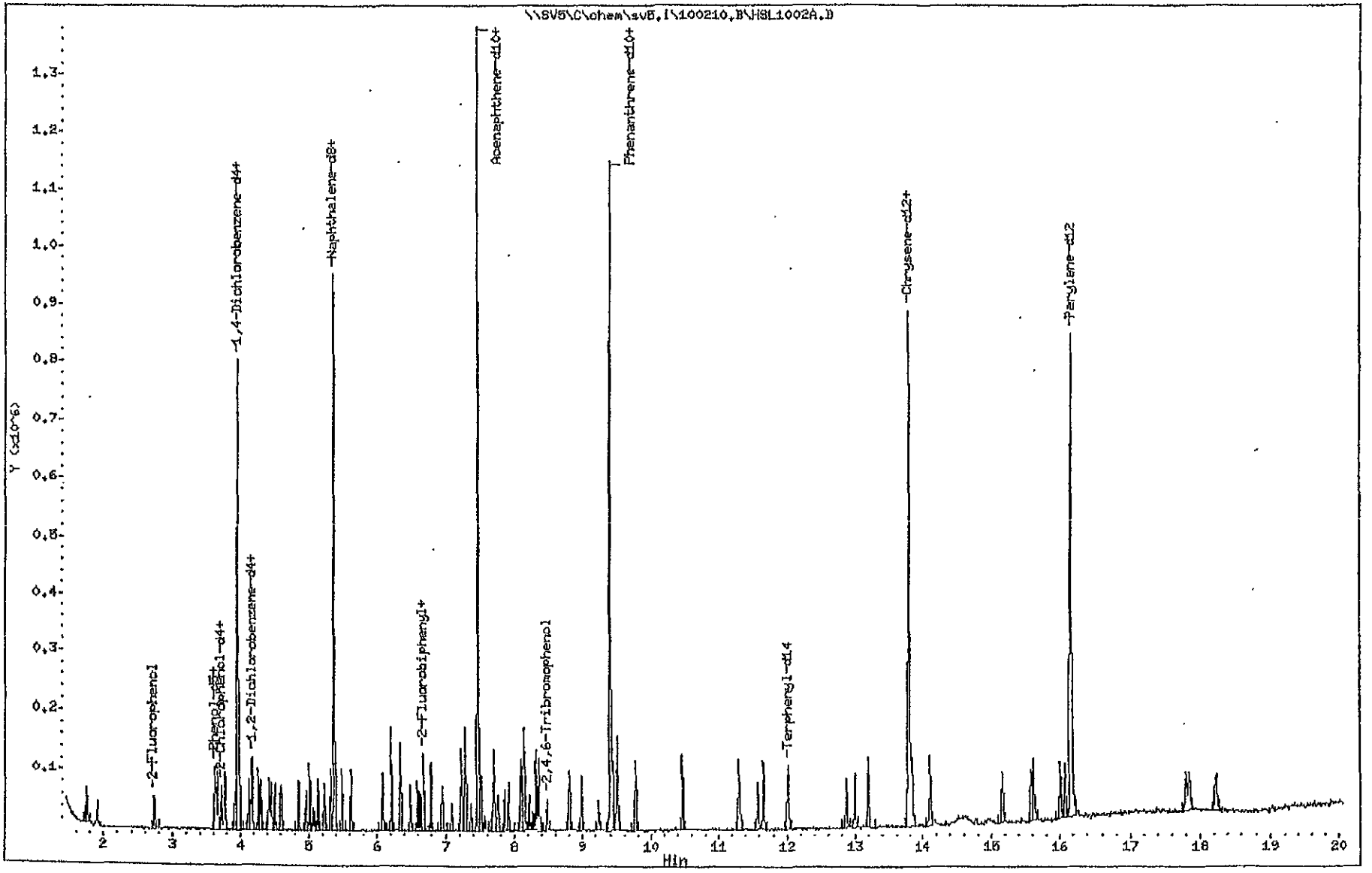
AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV5\C\chem\sv5.i\100210.B\HSL1002A.D
Date : 02-OCT-2010 12:27
Client ID: 8270F.M
Sample Info: HSL_005 ug/ml CS-1;1;;1;;;;;4

Instrument: sv5.i

Operator: KT
Column diameter: 2.00

Column phase:



TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002B.D
Lab Smp Id: HSL_010 ug/ml CS-2 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 12:53
Operator : KT Inst ID: sv5.i
Smp Info : HSL_010 ug/ml CS-2;1;;2;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0308;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:09 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 2 Calibration Sample, Level: 2
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SACP307UM

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
*	1 1,4-Dichlorobenzene-d4	152	3.955	3.955	(1.000)	116839	40.0000	(Q)
*	2 Naphthalene-d8	136	5.364	5.374	(1.000)	493196	40.0000	
*	3 Acenaphthene-d10	164	7.468	7.468	(1.000)	272639	40.0000	
*	4 Phenanthrene-d10	188	9.406	9.405	(1.000)	428440	40.0000	
*	5 Chrysene-d12	240	13.779	13.779	(1.000)	412260	40.0000	
*	6 Perylene-d12	264	16.162	16.162	(1.000)	419005	40.0000	
$	7 2-Fluorophenol	112	2.732	2.732	(0.691)	38100	10.0000	9.251
$	8 Phenol-d5	99	3.613	3.613	(0.914)	48878	10.0000	9.438
$	9 2-Chlorophenol-d4	132	3.747	3.758	(0.948)	45430	10.0000	9.989
$	10 1,2-Dichlorobenzene-d4	152	4.151	4.162	(1.050)	28658	10.0000	9.959
$	11 Nitrobenzene-d5	82	4.576	4.576	(0.853)	42237	10.0000	10.11 (QM)
$	12 2-Fluorobiphenyl	172	6.680	6.680	(0.895)	85886	10.0000	9.779
$	13 2,4,6-Tribromophenol	330	8.473	8.473	(1.135)	11265	10.0000	9.508
$	14 Terphenyl-d14	244	12.017	12.017	(0.872)	81626	10.0000	9.978
	15 N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	25783	10.0000	9.578 (Q)
	16 Pyridine	79	1.737	1.726	(0.439)	40141	10.0000	8.917 (Q)
	23 Aniline	93	3.654	3.654	(0.924)	63074	10.0000	9.568 (Q)
	24 Phenol	94	3.623	3.623	(0.916)	57313	10.0000	9.631 (Q)
	26 Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	40363	10.0000	9.677
	27 2-Chlorophenol	128	3.768	3.768	(0.953)	45449	10.0000	9.950
	28 1,3-Dichlorobenzene	146	3.913	3.923	(0.990)	49415	10.0000	9.932
	29 1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	52537	10.0000	10.10
	30 Benzyl Alcohol	108	4.120	4.120	(1.042)	30277	10.0000	9.862
	31 1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	47666	10.0000	9.966
	32 2-Methylphenol	108	4.255	4.255	(1.076)	40581	10.0000	9.714
	33 2,2'-oxybis(1-Chloropropane)	45	4.297	4.297	(1.086)	64869	10.0000	9.768
	34 4-Methylphenol	108	4.421	4.421	(1.118)	43497	10.0000	9.803
	36 Hexachloroethane	117	4.504	4.504	(1.139)	17770	10.0000	10.03
	37 N-Nitrosodipropylamine	70	4.442	4.442	(1.123)	28335	10.0000	9.587
	42 Nitrobenzene	77	4.597	4.597	(0.857)	40198	10.0000	9.845
	44 Isophorone	82	4.856	4.856	(0.905)	76804	10.0000	9.782
	45 2-Nitrophenol	139	4.960	4.960	(0.925)	23221	10.0000	9.585
	46 2,4-Dimethylphenol	107	5.012	5.012	(0.934)	42128	10.0000	9.787

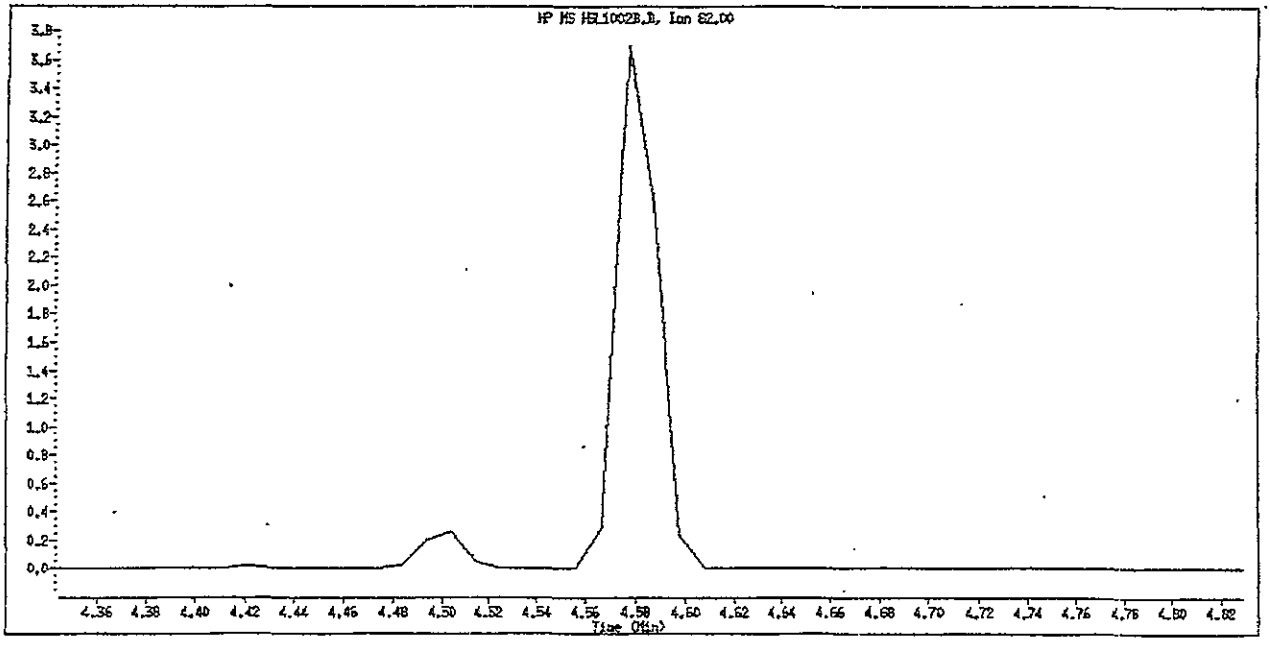
Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47 Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.956)	46230	10.0000	9.636
49 2,4-Dichlorophenol	162	5.229	5.229	(0.975)	32450	10.0000	9.744
50 Benzoic Acid	122	5.084	5.115	(0.948)	20056	10.0000	8.418
51 1,2,4-Trichlorobenzene	180	5.323	5.322	(0.992)	35544	10.0000	9.857
52 Naphthalene	128	5.395	5.395	(1.006)	138665	10.0000	10.18
54 4-Chloroaniline	127	5.488	5.488	(1.023)	52444	10.0000	9.825
57 Hexachlorobutadiene	225	5.613	5.613	(1.046)	17030	10.0000	9.650
60 4-Chloro-3-Methylphenol	107	6.069	6.069	(1.131)	35592	10.0000	9.570
63 2-Methylnaphthalene	142	6.203	6.203	(1.156)	83922	10.0000	9.811
66 Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	18919	10.0000	9.300
69 2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	20325	10.0000	9.344
70 2,4,5-Trichlorphenol	196	6.618	6.628	(0.886)	22419	10.0000	9.567
71 2-Chloronaphthalene	162	6.773	6.784	(0.907)	74574	10.0000	9.719
73 2-Nitroaniline	65	6.950	6.949	(0.931)	21647	10.0000	9.308
76 Dimethylphthalate	163	7.219	7.229	(0.967)	85330	10.0000	9.659
77 Acenaphthylene	152	7.281	7.281	(0.975)	130392	10.0000	9.758
79 2,6-Dinitrotoluene	165	7.291	7.302	(0.976)	18661	10.0000	9.067 (QM)
80 3-Nitroaniline	138	7.447	7.447	(0.997)	23598	10.0000	9.186 (q)
81 Acenaphthene	153	7.509	7.509	(1.006)	83474	10.0000	9.814
82 2,4-Dinitrophenol	184	7.571	7.572	(1.014)	7537	10.0000	10.11 (q)
83 Dibenzofuran	168	7.696	7.706	(1.031)	110503	10.0000	9.789
84 4-Nitrophenol	109	7.675	7.675	(1.028)	9643	10.0000	9.049 (Q)
86 2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	24530	10.0000	9.080
91 Fluorene	166	8.131	8.131	(1.089)	91225	10.0000	9.759
92 Diethylphthalate	149	8.100	8.100	(1.085)	88532	10.0000	9.788
93 4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	38113	10.0000	9.807
94 4-Nitroaniline	138	8.214	8.214	(1.100)	23002	10.0000	9.033
97 4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	11282	10.0000	11.10
98 N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	74860	11.7000	11.53
100 Azobenzene	77	8.349	8.348	(0.888)	82437	10.0000	9.784
101 4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	19823	10.0000	9.478
108 Hexachlorobenzene	284	8.981	8.981	(0.955)	23622	10.0000	10.11
110 Pentachlorophenol	266	9.240	9.240	(0.982)	10551	10.0000	10.90
114 Phenanthrene	178	9.437	9.437	(1.003)	134966	10.0000	9.995
115 Anthracene	178	9.499	9.499	(1.010)	130416	10.0000	9.667
118 Carbazole	167	9.768	9.768	(1.039)	120549	10.0000	9.782
120 Di-n-Butylphthalate	149	10.463	10.463	(1.112)	141693	10.0000	9.555
126 Fluoranthene	202	11.302	11.302	(1.202)	115262	10.0000	9.526
127 Benzidine	184	11.571	11.571	(0.840)	78774	10.0000	9.428
128 Pyrene	202	11.654	11.665	(0.846)	127577	10.0000	9.901
134 3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	66361	10.0000	8.997
136 Butylbenzylphthalate	149	12.991	12.991	(0.943)	62032	10.0000	9.605
138 Benzo(a)Anthracene	228	13.748	13.758	(0.998)	102788	10.0000	9.360
139 Chrysene	228	13.820	13.831	(1.003)	113552	10.0000	10.11
140 3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	38850	10.0000	9.379
141 bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.024)	83377	10.0000	9.372
142 Di-n-octylphthalate	149	15.157	15.167	(1.100)	126961	10.0000	8.528
144 Benzo(b)fluoranthene	252	15.572	15.582	(0.963)	84929	10.0000	8.954 (Q)
145 Benzo(k)fluoranthene	252	15.613	15.623	(0.966)	122065	10.0000	10.02 (q)
147 Benzo(e)pyrene	252	15.996	16.007	(0.990)	97140	10.0000	9.821
148 Benzo(a)pyrene	252	16.069	16.079	(0.994)	102327	10.0000	9.516
151 Indeno(1,2,3-cd)pyrene	276	17.789	17.800	(1.101)	76748	10.0000	8.824
152 Dibenzo(a,h)anthracene	278	17.841	17.841	(1.104)	88393	10.0000	9.097
153 Benzo(g,h,i)perylene	276	18.224	18.235	(1.128)	103135	10.0000	9.804

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				206994	10.0000	9.556 (A)

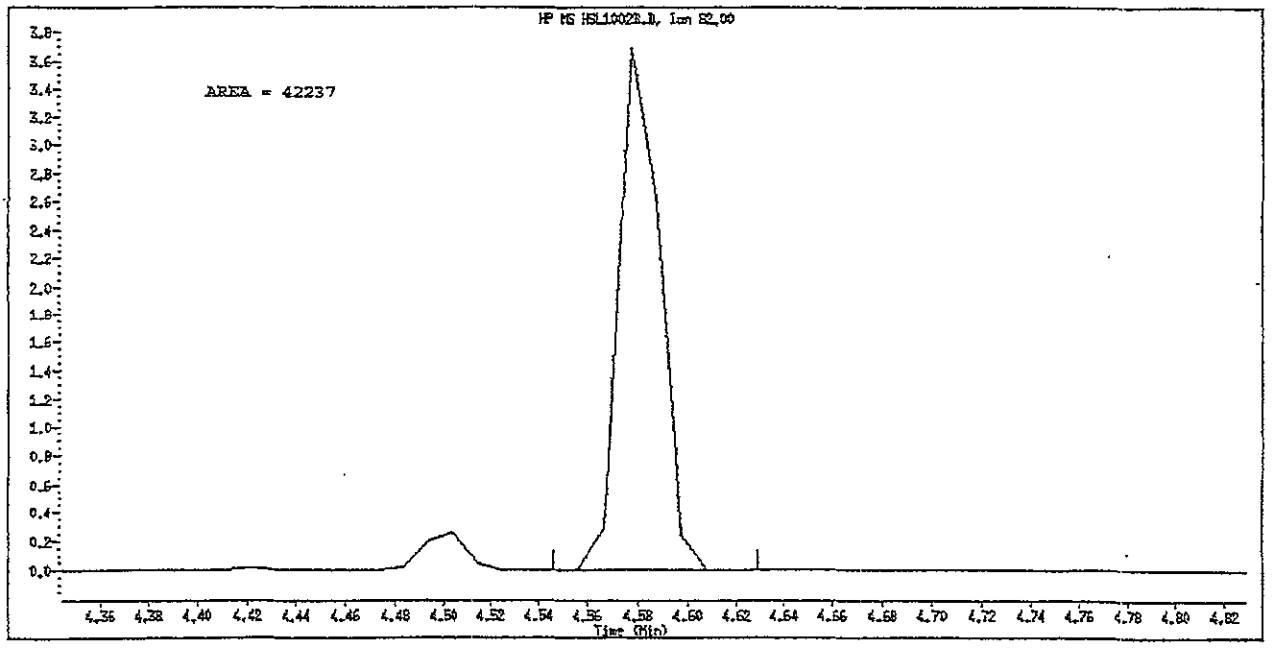
QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
M - Compound response manually integrated.
q - Qualifier signal exceeded ratio warning limit.

Data File Name: HSL1002B.D
Inj. Date and Time: 02-OCT-2010 12:53
Instrument ID: sv5.i
Client ID: 827DF.M
Compound Name: Nitrobenzene-d5
CAS #: 4165-60-0
Report Date: 10/03/2010



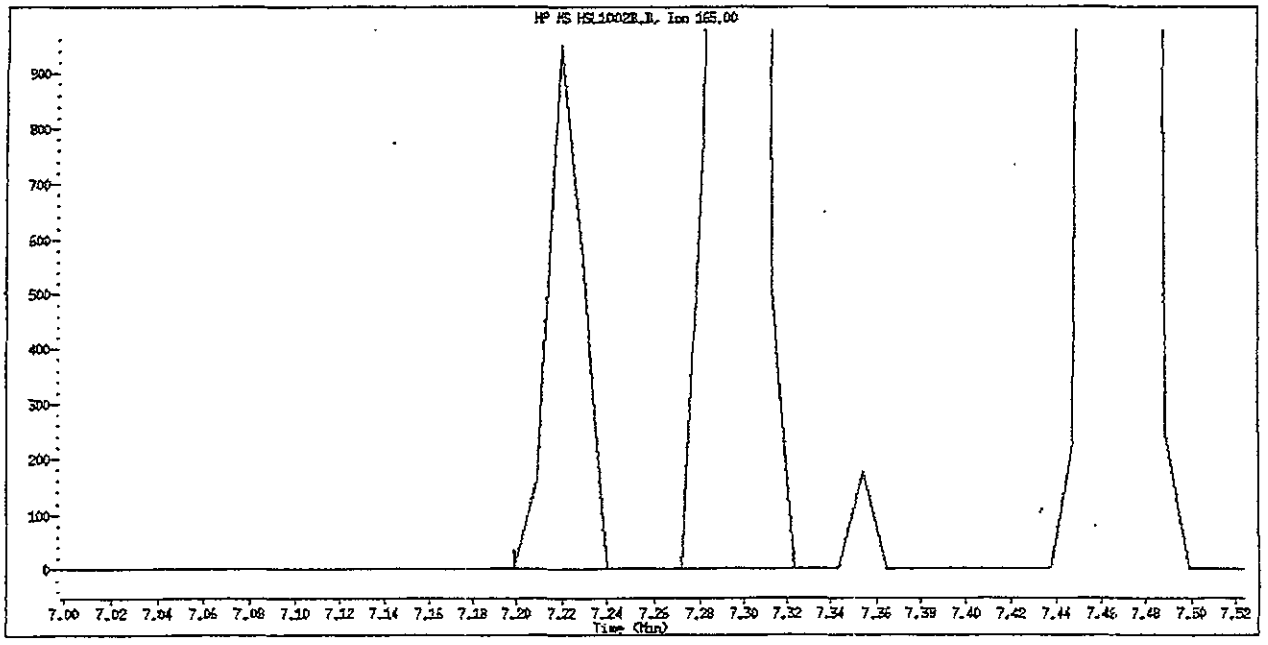
Original Integration



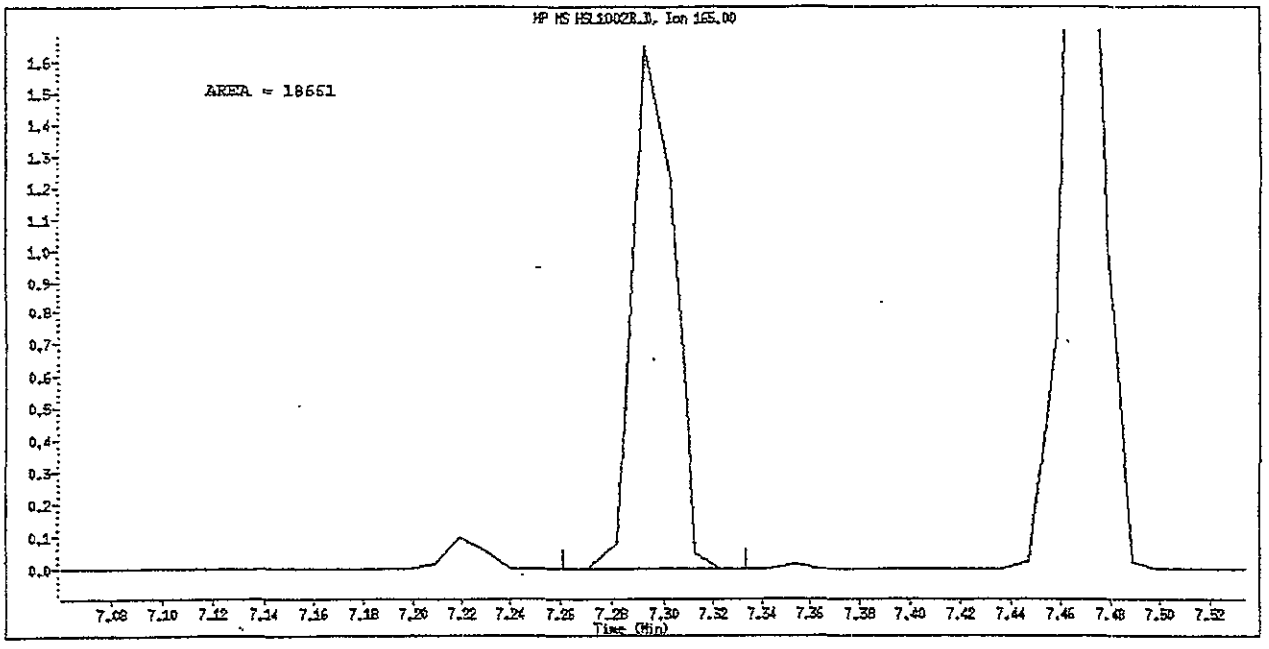
Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Peak Not Found

Data File Name: HSL1002B.D
Inj. Date and Time: 02-OCT-2010 12:53
Instrument ID: sv5.i
Client ID: 8270F.M
Compound Name: 2,6-Dinitrotoluene
CAS #: 606-20-2
Report Date: 10/03/2010



Original Integration



Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Poor Chromatography

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\100210.B\HSL1002B.D
Lab Smp Id: HSL_010 ug/ml CS-2 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 12:53
Operator : KT Inst ID: sv5.i
Smp Info : HSL_010 ug/ml CS-2;1;;2;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0308;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Meth Date : 02-Oct-2010 16:57 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 2 Calibration Sample, Level: 2
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
* 1	1,4-Dichlorobenzene-d4	152	3.955	3.955	(1.000)	116839	40.0000		(Q)
* 2	Naphthalene-d8	136	5.364	5.374	(1.000)	493196	40.0000		
* 3	Acenaphthene-d10	164	7.468	7.468	(1.000)	272639	40.0000		
* 4	Phenanthrene-d10	188	9.406	9.405	(1.000)	428440	40.0000		
* 5	Chrysene-d12	240	13.779	13.779	(1.000)	412260	40.0000		
* 6	Perylene-d12	264	16.162	16.162	(1.000)	419005	40.0000		
$ 7	2-Fluorophenol	112	2.732	2.732	(0.691)	38100	10.0000	8.835	
$ 8	Phenol-d5	99	3.613	3.613	(0.914)	48878	10.0000	8.913	
$ 9	2-Chlorophenol-d4	132	3.747	3.758	(0.948)	45430	10.0000	9.716	
$ 10	1,2-Dichlorobenzene-d4	152	4.151	4.162	(1.050)	28658	10.0000	9.820	
$ 11	Nitrobenzene-d5	82	Compound Not Detected.						
$ 12	2-Fluorobiphenyl	172	6.680	6.680	(0.895)	85886	10.0000	9.835	
$ 13	2,4,6-Tribromophenol	330	8.473	8.473	(1.135)	11265	10.0000	10.46	
$ 14	Terphenyl-d14	244	12.017	12.017	(0.872)	81026	10.0000	10.08	
15	N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	25783	10.0000	9.059	
16	Pyridine	79	1.737	1.726	(0.439)	40141	10.0000	8.464	
23	Aniline	93	3.654	3.654	(0.924)	63074	10.0000	9.172	(Q)
24	Phenol	94	3.623	3.623	(0.916)	57313	10.0000	9.827	(Q)
26	Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	40383	10.0000	9.114	
27	2-Chlorophenol	128	3.768	3.768	(0.953)	45449	10.0000	9.848	
28	1,3-Dichlorobenzene	146	3.913	3.923	(0.990)	49415	10.0000	9.589	
29	1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	52537	10.0000	10.19	
30	Benzyl Alcohol	108	4.120	4.120	(1.042)	30277	10.0000	9.547	
31	1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	47666	10.0000	9.755	
32	2-Methylphenol	108	4.255	4.255	(1.076)	40581	10.0000	9.413	
33	2,2'-oxybis(1-Chloropropane)	45	4.297	4.297	(1.086)	64869	10.0000	7.888	
34	4-Methylphenol	108	4.421	4.421	(1.118)	43497	10.0000	9.466	
36	Hexachloroethane	117	4.504	4.504	(1.139)	17770	10.0000	9.768	
37	N-Nitrosodipropylamine	70	4.442	4.442	(1.123)	28335	10.0000	8.809	
42	Nitrobenzene	77	4.597	4.597	(0.857)	40198	10.0000	9.246	
44	Isophorone	82	4.856	4.856	(0.905)	76804	10.0000	9.318	
45	2-Nitrophenol	139	4.960	4.960	(0.925)	23221	10.0000	9.784	
46	2,4-Dimethylphenol	107	5.012	5.012	(0.934)	42128	10.0000	9.559	

10-3-10

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.956)	46230	10.0000	9.421
49	2,4-Dichlorophenol	162	5.229	5.229	(0.975)	32450	10.0000	10.00
50	Benzoic Acid	122	5.084	5.115	(0.948)	20056	10.0000	8.164
51	1,2,4-Trichlorobenzene	180	5.323	5.322	(0.992)	35544	10.0000	10.11
52	Naphthalene	128	5.395	5.395	(1.006)	138665	10.0000	10.10
54	4-Chloroaniline	127	5.488	5.488	(1.023)	52444	10.0000	9.711
57	Hexachlorobutadiene	225	5.613	5.613	(1.046)	17030	10.0000	10.18
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.131)	35592	10.0000	9.536
63	2-Methylnaphthalene	142	6.203	6.203	(1.156)	83922	10.0000	10.02
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	18919	10.0000	9.098
69	2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	20325	10.0000	9.847
70	2,4,5-Trichlorophenol	196	6.628	6.628	(0.886)	22419	10.0000	9.889
71	2-Chloronaphthalene	162	6.773	6.784	(0.907)	74574	10.0000	9.756
73	2-Nitroaniline	65	6.950	6.949	(0.931)	21647	10.0000	8.456
76	Dimethylphthalate	163	7.219	7.229	(0.967)	85330	10.0000	9.665
77	Acenaphthylene	152	7.281	7.281	(0.975)	130392	10.0000	9.758
79	2,6-Dinitrotoluene	165	7.219	7.302	(0.967)	19698	10.0000	9.963 (Q)
80	3-Nitroaniline	138	7.447	7.447	(0.997)	23598	10.0000	9.002 (q)
81	Acenaphthene	153	7.509	7.509	(1.006)	83474	10.0000	9.804
82	2,4-Dinitrophenol	184	7.571	7.571	(1.014)	7537	10.0000	9.147 (q)
83	Dibenzofuran	168	7.696	7.706	(1.031)	110503	10.0000	9.824
84	4-Nitrophenol	109	7.675	7.675	(1.028)	9643	10.0000	8.425 (Q)
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	24530	10.0000	9.262
91	Fluorene	166	8.131	8.131	(1.089)	91225	10.0000	9.898
92	Diethylphthalate	149	8.100	8.100	(1.085)	88532	10.0000	9.594
93	4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	38113	10.0000	10.03
94	4-Nitroaniline	138	8.214	8.214	(1.100)	23002	10.0000	8.977
97	4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	11282	10.0000	11.76
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	74860	11.7000	11.21
100	Azobenzene	77	8.349	8.348	(0.888)	82437	10.0000	8.875
101	4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	19823	10.0000	9.575
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	23622	10.0000	10.56
110	Pentachlorophenol	266	9.240	9.240	(0.982)	10551	10.0000	7.861
114	Phenanthrene	178	9.437	9.437	(1.003)	134966	10.0000	10.10
115	Anthracene	178	9.498	9.499	(1.010)	130416	10.0000	9.697
118	Carbazole	167	9.768	9.768	(1.039)	120549	10.0000	9.637
120	Di-n-Butylphthalate	149	10.463	10.463	(1.112)	141693	10.0000	9.367
125	Fluoranthene	202	11.302	11.302	(1.202)	115262	10.0000	9.583
127	Benzidine	184	11.571	11.571	(0.840)	78774	10.0000	9.305
128	Pyrene	202	11.654	11.665	(0.846)	127577	10.0000	9.897
134	3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	66361	10.0000	9.134
136	Butylbenzylphthalate	149	12.991	12.991	(0.943)	62032	10.0000	9.418
138	Benzo(a)Anthracene	228	13.748	13.758	(0.998)	102788	10.0000	9.450
139	Chrysene	228	13.820	13.831	(1.003)	113552	10.0000	10.05
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	38850	10.0000	9.762
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.024)	83377	10.0000	9.194
142	Di-n-octylphthalate	149	15.157	15.167	(1.100)	126961	10.0000	8.756
144	Benzo(b)fluoranthene	252	15.572	15.582	(0.963)	84329	10.0000	8.548 (Q)
145	Benzo(k)fluoranthene	252	15.613	15.623	(0.966)	122065	10.0000	10.51 (q)
147	Benzo(e)pyrene	252	15.996	16.007	(0.990)	97140	10.0000	9.863
148	Benzo(a)pyrene	252	16.069	16.079	(0.994)	102327	10.0000	9.463
151	Indeno(1,2,3-cd)pyrene	276	17.789	17.800	(1.101)	76748	10.0000	8.022
152	Dibenzo(a,h)anthracene	278	17.841	17.841	(1.104)	88393	10.0000	9.016
153	Benzo(g,h,i)perylene	276	18.224	18.235	(1.128)	103135	10.0000	9.811

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				206994	10.0000	9.607(A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002B.D
Lab Smp Id: HSL_010 ug/ml CS-2
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\sv5\c\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0308;0;8270F.M

Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	116839	-4.72
2 Naphthalene-d8	530514	265257	1061028	493196	-7.03
3 Acenaphthene-d10	282538	141269	565076	272639	-3.50
4 Phenanthrene-d10	462722	231361	925444	428440	-7.41
5 Chrysene-d12	435850	217925	871700	412260	-5.41
6 Perylene-d12	422284	211142	844568	419005	-0.78

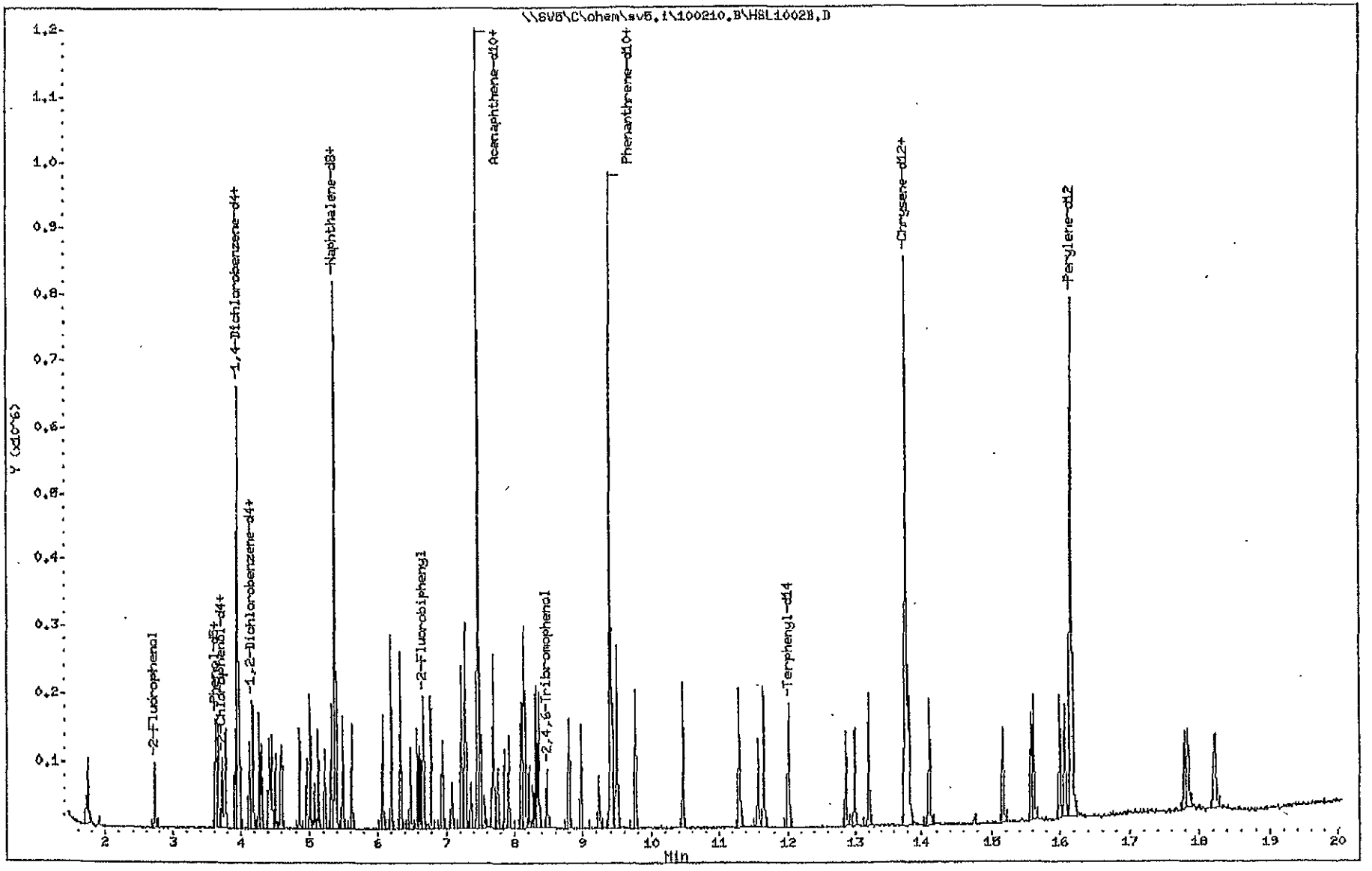
COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.96	3.46	4.46	3.96	0.00
2 Naphthalene-d8	5.37	4.87	5.87	5.36	-0.19
3 Acenaphthene-d10	7.47	6.97	7.97	7.47	0.00
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	0.00
5 Chrysene-d12	13.78	13.28	14.28	13.78	0.00
6 Perylene-d12	16.16	15.66	16.66	16.16	0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV5\C\chem\sv5.i\100210.B\HSL1002B.D
Date : 02-OCT-2010 12:53
Client ID: 8270F.H
Sample Info: HSL_010 ug/ml CS-2;1;;2;;;;;4

Instrument: sv5.i
Operator: KT
Column diameter: 2.00

Column phase:



TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002C.D
Lab Smp Id: HSL 020 ug/ml CS-3 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 13:18
Operator : KT Inst ID: sv5.i
Smp Info : HSL 020 ug/ml CS-3;1;;3;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0309;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:09 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 3 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SACP307UM

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
* 1	1,4-Dichlorobenzene-d4	152	3.954	3.955	(1.000)	145926	40.0000	(Q)
* 2	Naphthalene-d8	136	5.364	5.374	(1.000)	625682	40.0000	
* 3	Acenaphthene-d10	164	7.467	7.468	(1.000)	328608	40.0000	
* 4	Phenanthrene-d10	188	9.405	9.405	(1.000)	525834	40.0000	
* 5	Chrysene-d12	240	13.779	13.779	(1.000)	590727	40.0000	
* 6	Perylene-d12	264	16.162	16.162	(1.000)	619266	40.0000	
$ 7	2-Fluorophenol	112	2.732	2.732	(0.691)	100961	20.0000	19.63
$ 8	Phenol-d5	99	3.612	3.613	(0.914)	127066	20.0000	19.64
$ 9	2-Chlorophenol-d4	132	3.747	3.758	(0.948)	112302	20.0000	19.77
$ 10	1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	72837	20.0000	20.27(q)
$ 11	Nitrobenzene-d5	82	4.576	4.576	(0.853)	103440	20.0000	19.52
$ 12	2-Fluorobiphenyl	172	6.680	6.680	(0.895)	209764	20.0000	19.82
$ 13	2,4,6-Tribromophenol	330	8.473	8.473	(1.135)	28698	20.0000	20.10
$ 14	Terphenyl-d14	244	12.017	12.017	(0.872)	218324	20.0000	18.76
15	N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	66431	20.0000	19.76(q)
16	Pyridine	79	1.726	1.726	(0.437)	116339	20.0000	20.69(Q)
23	Aniline	93	3.654	3.654	(0.924)	160510	20.0000	19.50
24	Phenol	94	3.623	3.623	(0.916)	147994	20.0000	19.91
26	Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	101777	20.0000	19.53
27	2-Chlorophenol	128	3.768	3.768	(0.953)	114481	20.0000	20.07
28	1,3-Dichlorobenzene	146	3.913	3.923	(0.990)	122398	20.0000	19.70
29	1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	126965	20.0000	19.54
30	Benzyl Alcohol	108	4.120	4.120	(1.042)	72366	20.0000	18.87
31	1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	117073	20.0000	19.60
32	2-Methylphenol	108	4.255	4.255	(1.076)	101499	20.0000	19.45
33	2,2'-oxybis(1-Chloropropane)	45	4.296	4.297	(1.086)	166596	20.0000	20.08
34	4-Methylphenol	108	4.421	4.421	(1.118)	106723	20.0000	19.26
36	Hexachloroethane	117	4.504	4.504	(1.139)	44196	20.0000	19.98
37	N-Nitrosodipropylamine	70	4.441	4.442	(1.123)	73913	20.0000	20.02
42	Nitrobenzene	77	4.597	4.597	(0.857)	101809	20.0000	19.65
44	Isophorone	82	4.856	4.856	(0.905)	191333	20.0000	19.21
45	2-Nitrophenol	139	4.960	4.960	(0.925)	58938	20.0000	19.18
46	2,4-Dimethylphenol	107	5.011	5.012	(0.934)	107325	20.0000	19.65

10-3-10

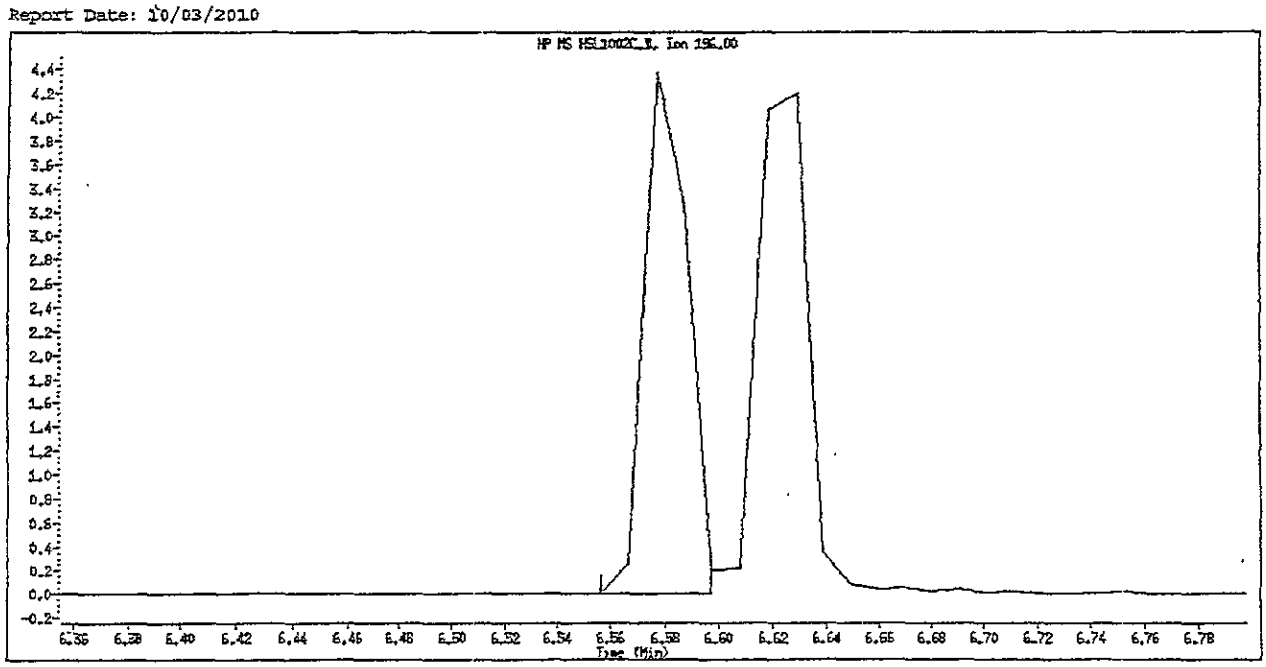
Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47 Bis(2-chloroethoxy)methane	93	5.125	5.126	(0.956)	120646	20.0000	19.82
49 2,4-Dichlorophenol	162	5.229	5.229	(0.975)	84525	20.0000	20.01
50 Benzoic Acid	122	5.094	5.115	(0.950)	54506	20.0000	18.03
51 1,2,4-Trichlorobenzene	180	5.322	5.322	(0.992)	89082	20.0000	19.47
52 Naphthalene	128	5.395	5.395	(1.006)	336100	20.0000	19.46
54 4-Chloroaniline	127	5.488	5.488	(1.023)	135348	20.0000	19.99
57 Hexachlorobutadiene	225	5.613	5.613	(1.046)	45138	20.0000	20.16
60 4-Chloro-3-Methylphenol	107	6.068	6.069	(1.131)	90970	20.0000	19.28
63 2-Methylnaphthalene	142	6.203	6.203	(1.156)	212981	20.0000	19.62
66 Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	47478	20.0000	19.36
69 2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	49658	20.0000	18.94 (Q)
70 2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	55529	20.0000	19.66 (QM)
71 2-Chloronaphthalene	162	6.784	6.784	(0.908)	180754	20.0000	19.54
73 2-Nitroaniline	65	6.949	6.949	(0.931)	54872	20.0000	19.58
76 Dimethylphthalate	163	7.219	7.229	(0.967)	213272	20.0000	20.03
77 Acenaphthylene	152	7.281	7.281	(0.975)	315165	20.0000	19.57
79 2,6-Dinitrotoluene	165	7.291	7.302	(0.976)	49111	20.0000	19.80 (QM)
80 3-Nitroaniline	138	7.447	7.447	(0.997)	59114	20.0000	19.09
81 Acenaphthene	153	7.509	7.509	(1.006)	208228	20.0000	20.31
82 2,4-Dinitrophenol	184	7.571	7.572	(1.014)	23799	20.0000	19.52
83 Dibenzofuran	168	7.695	7.706	(1.031)	271431	20.0000	19.95
84 4-Nitrophenol	109	7.675	7.675	(1.028)	25164	20.0000	19.59 (Q)
86 2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	63223	20.0000	19.42
91 Fluorene	166	8.131	8.131	(1.089)	220647	20.0000	19.58
92 Diethylphthalate	149	8.100	8.100	(1.085)	216140	20.0000	19.83
93 4-Chlorophenyl-phenylether	204	8.151	8.152	(1.092)	93458	20.0000	19.95
94 4-Nitroaniline	138	8.214	8.214	(1.100)	61333	20.0000	19.98
97 4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	32982	20.0000	20.44
98 N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	186206	23.4000	23.36
100 Azobenzene	77	8.348	8.348	(0.888)	203290	20.0000	19.66
101 4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	50693	20.0000	19.75
108 Hexachlorobenzene	284	8.980	8.981	(0.955)	54528	20.0000	19.02
110 Pentachlorophenol	266	9.240	9.240	(0.982)	30451	20.0000	20.33
114 Phenanthrene	178	9.436	9.437	(1.003)	329718	20.0000	19.89
115 Anthracene	178	9.499	9.499	(1.010)	326558	20.0000	19.72
118 Carbazole	167	9.768	9.768	(1.038)	298921	20.0000	19.76
120 Di-n-Butylphthalate	149	10.462	10.463	(1.112)	358075	20.0000	19.58
126 Fluoranthene	202	11.302	11.302	(1.202)	308182	20.0000	20.75
127 Benzidine	184	11.571	11.571	(0.840)	222260	20.0000	18.56
128 Pyrene	202	11.665	11.665	(0.847)	345805	20.0000	18.73
134 3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	198960	20.0000	18.82
136 Butylbenzylphthalate	149	12.991	12.991	(0.943)	174685	20.0000	18.88
138 Benzo(a)Anthracene	228	13.758	13.758	(0.998)	304948	20.0000	19.38
139 Chrysene	228	13.820	13.831	(1.003)	314030	20.0000	19.51
140 3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	115458	20.0000	19.45
141 bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.024)	248201	20.0000	19.47
142 Di-n-octylphthalate	149	15.157	15.167	(1.100)	400592	20.0000	19.66
144 Benzo(b)fluoranthene	252	15.582	15.582	(0.964)	256213	20.0000	18.28 (Q)
145 Benzo(k)fluoranthene	252	15.613	15.623	(0.966)	371629	20.0000	20.65 (Q)
147 Benzo(e)pyrene	252	15.996	16.007	(0.990)	281015	20.0000	19.22
148 Benzo(a)pyrene	252	16.069	16.079	(0.994)	307781	20.0000	19.37
151 Indeno(1,2,3-cd)pyrene	276	17.789	17.800	(1.101)	228110	20.0000	17.74
152 Dibenzo(a,h)anthracene	278	17.841	17.841	(1.104)	270172	20.0000	18.81
153 Benzo(g,h,i)perylene	276	18.224	18.235	(1.128)	301520	20.0000	19.39

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				627842	20.0000	19.61 (A)

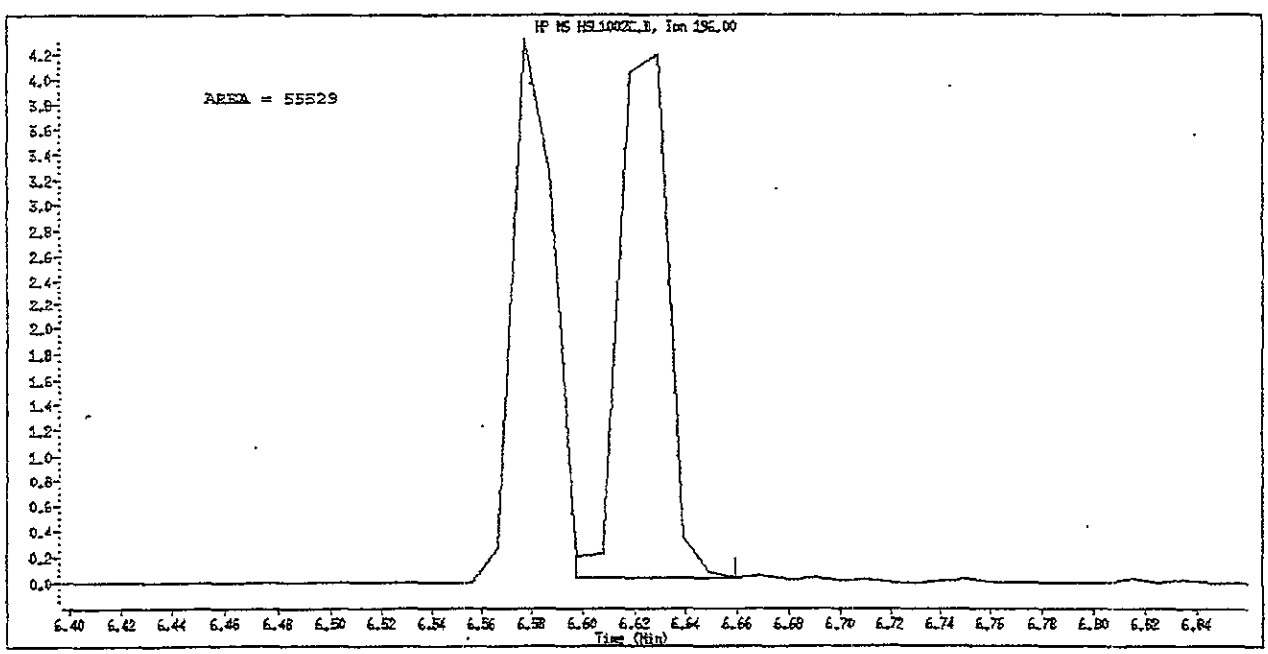
QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
M - Compound response manually integrated.
q - Qualifier signal exceeded ratio warning limit.

Data File Name: HSL1002C.D
Inj. Date and Time: 02-OCT-2010 13:18
Instrument ID: sv5.i
Client ID: 8270F.M
Compound Name: 2,4,5-Trichlorphenol
CAS #: 95-95-4
Report Date: 10/03/2010



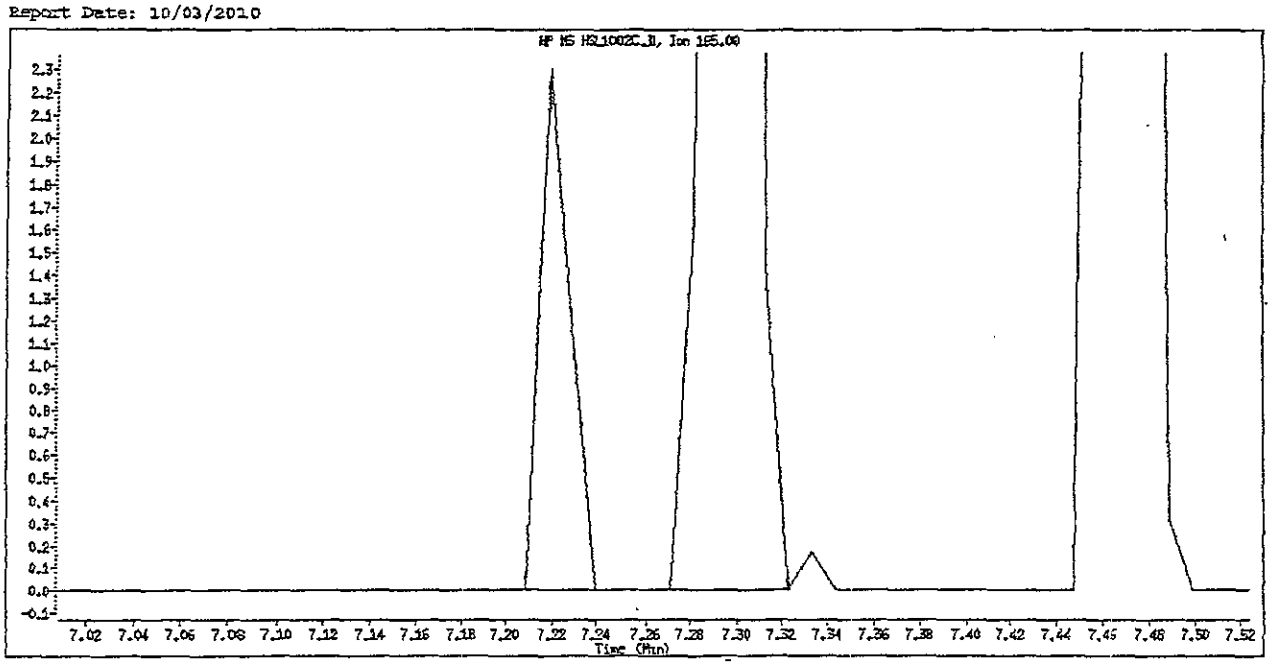
Original Integration



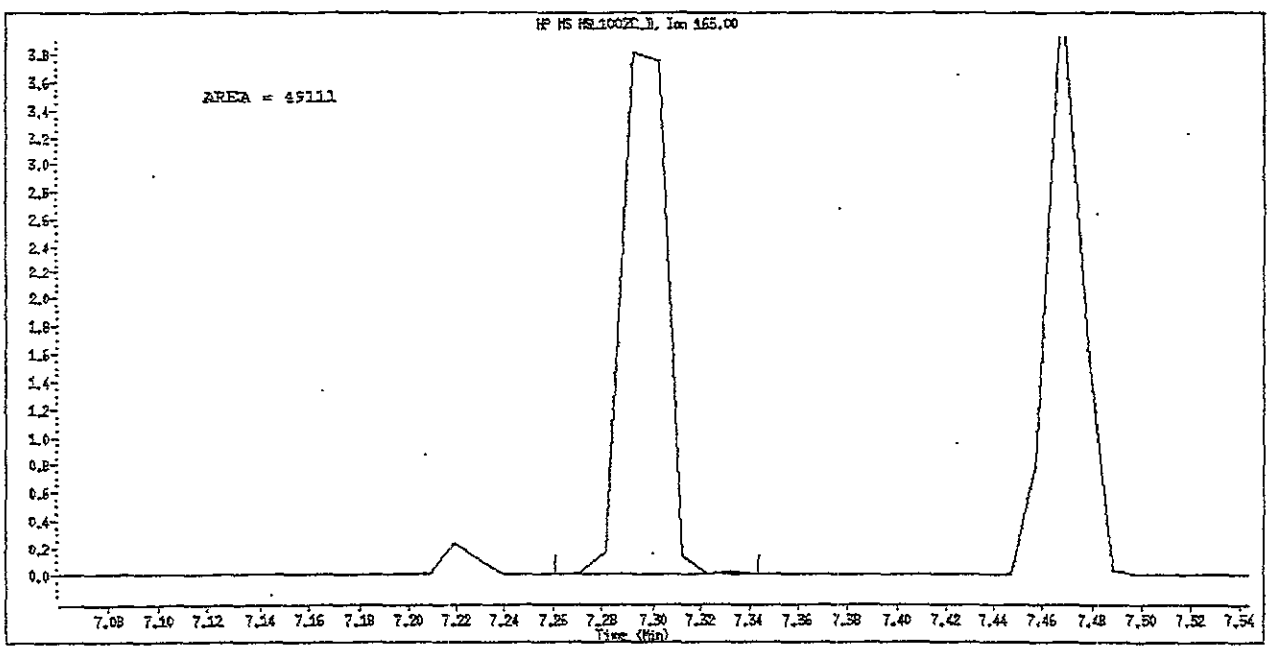
Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Wrong Peak

Data File Name: BSL1002C.D
Inj. Date and Time: 02-OCT-2010 13:18
Instrument ID: svS.i
Client ID: 8270F.M
Compound Name: 2,6-Dinitrotoluene
CAS #: 606-20-2
Report Date: 10/03/2010



Original Integration



Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Poor Chromatography

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\100210.B\HSL1002C.D
Lab Smp Id: HSL_020 ug/ml CS-3 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 13:18
Operator : KT Inst ID: sv5.i
Smp Info : HSL_020 ug/ml CS-3;1;;3;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0309;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Meth Date : 02-Oct-2010 16:57 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 3 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
*	1 1,4-Dichlorobenzene-d4	152	3.954	3.955	(1.000)	145926	40.0000		(Q)
*	2 Naphthalene-d8	136	5.364	5.374	(1.000)	625682	40.0000		
*	3 Acenaphthene-d10	164	7.467	7.468	(1.000)	328608	40.0000		
*	4 Phenanthrene-d10	188	9.405	9.405	(1.000)	525834	40.0000		
*	5 Chrysene-d12	240	13.779	13.779	(1.000)	590727	40.0000		
*	6 Perylene-d12	264	16.162	16.162	(1.000)	619266	40.0000		
$	7 2-Fluorophenol	112	2.732	2.732	(0.691)	100961	20.0000	18.75	
$	8 Phenol-d5	99	3.612	3.613	(0.914)	127066	20.0000	18.55	
$	9 2-Chlorophenol-d4	132	3.747	3.758	(0.948)	112302	20.0000	19.23	
$	10 1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	72837	20.0000	19.98	(q)
$	11 Nitrobenzene-d5	82	4.576	4.576	(0.853)	103440	20.0000	18.64	
$	12 2-Fluorobiphenyl	172	6.680	6.680	(0.895)	209764	20.0000	19.93	
$	13 2,4,6-Tribromophenol	330	8.473	8.473	(1.135)	28698	20.0000	22.12	
$	14 Terphenyl-d14	244	12.017	12.017	(0.872)	218324	20.0000	18.95	
	15 N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	66431	20.0000	18.69	
	16 Pyridine	79	1.726	1.726	(0.437)	116339	20.0000	19.64	
	23 Aniline	93	3.654	3.654	(0.924)	160510	20.0000	18.69	
	24 Phenol	94	3.623	3.623	(0.916)	147994	20.0000	20.32	
	26 Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	101777	20.0000	18.39	
	27 2-Chlorophenol	128	3.768	3.768	(0.953)	114481	20.0000	19.86	
	28 1,3-Dichlorobenzene	146	3.913	3.923	(0.990)	122398	20.0000	19.22	
	29 1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	126965	20.0000	19.72	
	30 Benzyl Alcohol	108	4.120	4.120	(1.042)	72366	20.0000	18.27	
	31 1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	117073	20.0000	19.18	
	32 2-Methylphenol	108	4.255	4.255	(1.076)	101499	20.0000	18.85	
	33 2,2'-oxybis(1-Chloropropane)	45	4.296	4.297	(1.086)	166596	20.0000	16.22	
	34 4-Methylphenol	108	4.421	4.421	(1.118)	106723	20.0000	18.60	
	36 Hexachloroethane	117	4.504	4.504	(1.139)	44196	20.0000	19.45	
	37 N-Nitrosodipropylamine	70	4.441	4.442	(1.123)	73913	20.0000	18.40	
	42 Nitrobenzene	77	4.597	4.597	(0.857)	101809	20.0000	18.46	
	44 Isophorone	82	4.856	4.856	(0.905)	191333	20.0000	18.30	
	45 2-Nitrophenol	139	4.960	4.960	(0.925)	58938	20.0000	19.57	
	46 2,4-Dimethyphenol	107	5.011	5.012	(0.934)	107325	20.0000	19.20	

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47 Bis(2-chloroethoxy)methane	93	5.125	5.126	(0.956)	120646	20.0000	19.38
49 2,4-Dichlorophenol	162	5.229	5.229	(0.975)	84525	20.0000	20.54
50 Benzoic Acid	122	5.094	5.115	(0.950)	54506	20.0000	17.49
51 1,2,4-Trichlorobenzene	180	5.322	5.322	(0.992)	89082	20.0000	19.97
52 Naphthalene	128	5.395	5.395	(1.006)	336100	20.0000	19.30
54 4-Chloroaniline	127	5.488	5.488	(1.023)	135348	20.0000	19.76
57 Hexachlorobutadiene	225	5.613	5.613	(1.046)	45138	20.0000	21.26
60 4-Chloro-3-Methylphenol	107	6.068	6.069	(1.131)	90970	20.0000	19.21
63 2-Methylnaphthalene	142	6.203	6.203	(1.156)	212981	20.0000	20.04
66 Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	47478	20.0000	18.94
69 2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	49658	20.0000	19.96 (Q)
70 2,4,5-Trichlorphenol	196	6.576	6.628	(0.881)	49658	20.0000	18.17 (Q)
71 2-Chloronaphthalene	162	6.784	6.784	(0.908)	180754	20.0000	19.62
73 2-Nitroaniline	65	6.949	6.949	(0.931)	54872	20.0000	17.78
76 Dimethylphthalate	163	7.219	7.229	(0.967)	213272	20.0000	20.04
77 Acenaphthylene	152	7.281	7.281	(0.975)	315165	20.0000	19.57
79 2,6-Dinitrotoluene	165	7.218	7.302	(0.967)	51125	20.0000	21.45 (Q)
80 3-Nitroaniline	138	7.447	7.447	(0.997)	59114	20.0000	18.71
81 Acenaphthene	153	7.509	7.509	(1.006)	208228	20.0000	20.29
82 2,4-Dinitrophenol	184	7.571	7.571	(1.014)	23799	20.0000	19.22
83 Dibenzofuran	168	7.695	7.706	(1.031)	271431	20.0000	20.02
84 4-Nitrophenol	109	7.675	7.675	(1.028)	25164	20.0000	18.24 (Q)
86 2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	63223	20.0000	19.81
91 Fluorene	166	8.131	8.131	(1.089)	220647	20.0000	19.86
92 Diethylphthalate	149	8.100	8.100	(1.085)	216140	20.0000	19.43
93 4-Chlorophenyl-phenylether	204	8.151	8.152	(1.092)	93468	20.0000	20.41
94 4-Nitroaniline	138	8.214	8.214	(1.100)	61333	20.0000	19.86
97 4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	32982	20.0000	20.90
98 N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	186206	23.4000	22.72
100 Azobenzene	77	8.348	8.348	(0.888)	203290	20.0000	17.83
101 4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	50693	20.0000	19.95
108 Hexachlorobenzene	284	8.980	8.981	(0.955)	54528	20.0000	19.87
110 Pentachlorophenol	266	9.240	9.240	(0.982)	30451	20.0000	18.48
114 Phenanthrene	178	9.436	9.437	(1.003)	329718	20.0000	20.10
115 Anthracene	178	9.499	9.499	(1.010)	326558	20.0000	19.78
118 Carbazole	167	9.768	9.768	(1.039)	298921	20.0000	19.47
120 Di-n-Butylphthalate	149	10.462	10.463	(1.112)	358075	20.0000	19.29
126 Fluoranthene	202	11.302	11.302	(1.202)	308182	20.0000	20.88
127 Benzidine	184	11.571	11.571	(0.840)	222260	20.0000	18.32
128 Pyrene	202	11.665	11.665	(0.847)	345805	20.0000	18.72
134 3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	198960	20.0000	19.11
136 Butylbenzylphthalate	149	12.991	12.991	(0.943)	174685	20.0000	18.51
138 Benzo(a)Anthracene	228	13.758	13.758	(0.998)	304968	20.0000	19.57
139 Chrysene	228	13.820	13.831	(1.003)	314030	20.0000	19.39
140 3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	115458	20.0000	20.25
141 bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.024)	248201	20.0000	19.10
142 Di-n-octylphthalate	149	15.157	15.167	(1.100)	400592	20.0000	19.28
144 Benzo(b)fluoranthene	252	15.582	15.582	(0.964)	256213	20.0000	17.45 (Q)
145 Benzo(k)fluoranthene	252	15.613	15.623	(0.966)	371629	20.0000	21.66 (q)
147 Benzo(e)pyrene	252	15.996	16.007	(0.990)	281015	20.0000	19.30
148 Benzo(a)pyrene	252	16.069	16.079	(0.994)	307781	20.0000	19.26
151 Indeno(1,2,3-cd)pyrene	276	17.789	17.800	(1.101)	228110	20.0000	16.13
152 Dibenzo(a,h)anthracene	278	17.841	17.841	(1.104)	270172	20.0000	18.64
153 Benzo(g,h,i)perylene	276	18.224	18.235	(1.128)	301520	20.0000	19.41

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				627842	20.0000	19.72 (A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002C.D
Lab Smp Id: HSL_020 ug/ml CS-3
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\sv5\c\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0309;0;8270F.M

Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	145926	19.00
2 Naphthalene-d8	530514	265257	1061028	625682	17.94
3 Acenaphthene-d10	282538	141269	565076	328608	16.31
4 Phenanthrene-d10	462722	231361	925444	525834	13.64
5 Chrysene-d12	435850	217925	871700	590727	35.53
6 Perylene-d12	422284	211142	844568	619266	46.65

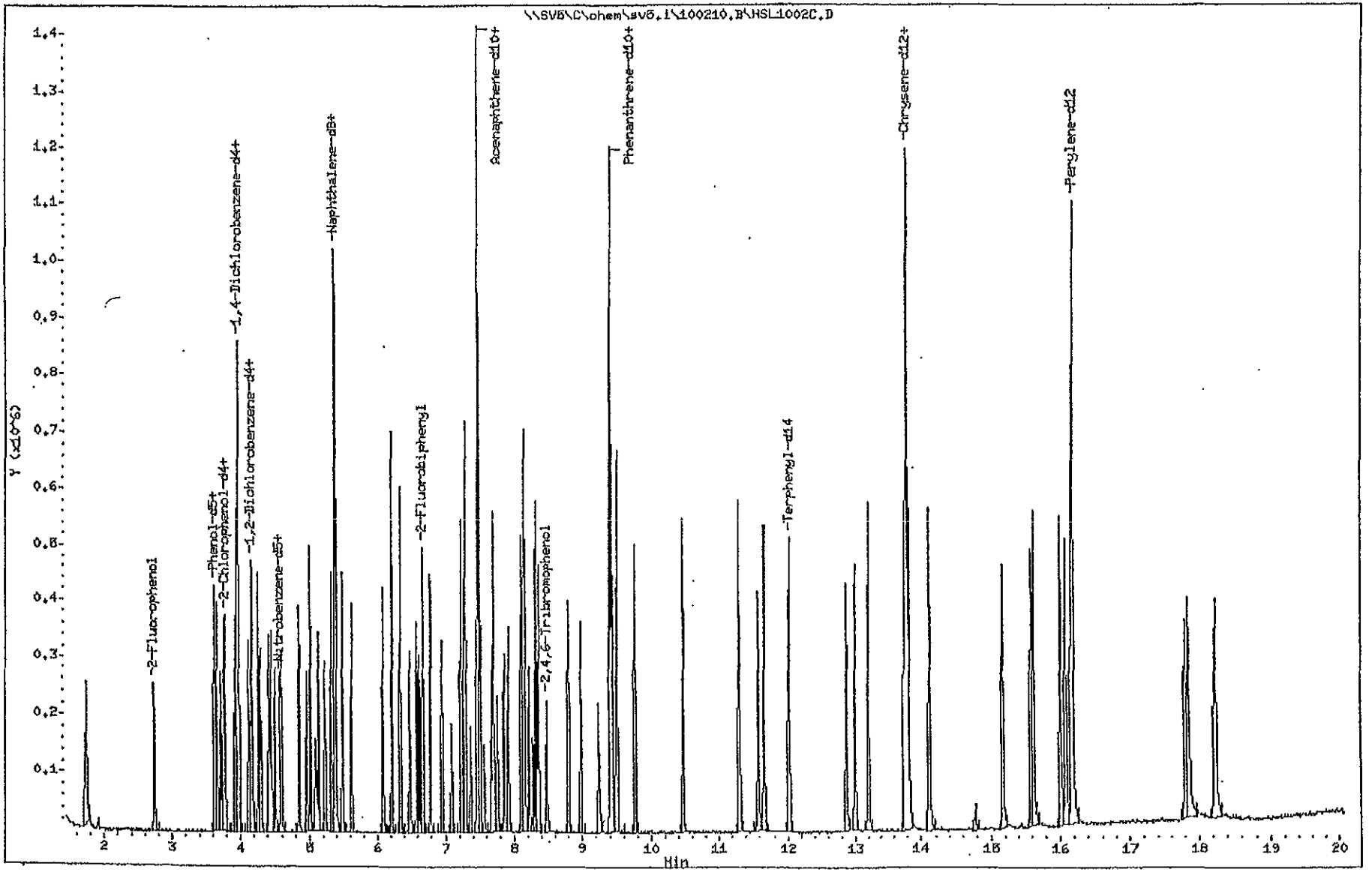
COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.96	3.46	4.46	3.95	-0.00
2 Naphthalene-d8	5.37	4.87	5.87	5.36	-0.20
3 Acenaphthene-d10	7.47	6.97	7.97	7.47	-0.00
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	-0.00
5 Chrysene-d12	13.78	13.28	14.28	13.78	-0.00
6 Perylene-d12	16.16	15.66	16.66	16.16	-0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV5\C\chem\sv5.i\100210.B\HSL1002C.D
Date : 02-OCT-2010 13:18
Client ID: 8270F,H
Sample Info: HSL_020 ug/ml CS-3:1::3:::::4

Instrument: sv5.i
Operator: KT
Column diameter: 2.00

Column phase:



TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002D.D
Lab Smp Id: HSL_050 ug/ml CS-4 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 13:44
Operator : KT Inst ID: sv5.i
Smp Info : HSL_050 ug/ml CS-4;1;;4;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0310;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:09 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 4 Calibration Sample, Level: 4
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SACP307UM

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
*	1 1,4-Dichlorobenzene-d4	152	3.955	3.955	(1.000)	122625	40.0000	
*	2 Naphthalene-d8	136	5.374	5.374	(1.000)	530514	40.0000	
*	3 Acenaphthene-d10	164	7.468	7.468	(1.000)	282538	40.0000	
*	4 Phenanthrene-d10	188	9.405	9.405	(1.000)	462722	40.0000	
*	5 Chrysene-d12	240	13.779	13.779	(1.000)	435850	40.0000	
*	6 Perylene-d12	264	16.162	16.162	(1.000)	422284	40.0000	
$	7 2-Fluorophenol	112	2.732	2.732	(0.691)	220986	50.0000	51.13
$	8 Phenol-d5	99	3.613	3.613	(0.914)	274382	50.0000	50.48
$	9 2-Chlorophenol-d4	132	3.758	3.758	(0.950)	244352	50.0000	51.19
$	10 1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	151616	50.0000	50.20
$	11 Nitrobenzene-d5	82	4.576	4.576	(0.852)	226162	50.0000	50.33
$	12 2-Fluorobiphenyl	172	6.680	6.680	(0.895)	473978	50.0000	52.08
$	13 2,4,6-Tribromophenol	330	8.473	8.473	(1.135)	63321	50.0000	51.57
$	14 Terphenyl-d14	244	12.017	12.017	(0.872)	438253	50.0000	51.05
	15 N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	140972	50.0000	49.90 (M)
	16 Pyridine	79	1.726	1.726	(0.437)	240053	50.0000	50.81 (M)
	23 Aniline	93	3.654	3.654	(0.924)	346504	50.0000	50.08
	24 Phenol	94	3.623	3.623	(0.916)	311820	50.0000	49.93
	26 Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	220455	50.0000	50.34
	27 2-Chlorophenol	128	3.768	3.768	(0.953)	242442	50.0000	50.57
	28 1,3-Dichlorobenzene	146	3.923	3.923	(0.992)	265384	50.0000	50.82
	29 1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	271151	50.0000	49.66
	30 Benzyl Alcohol	108	4.120	4.120	(1.042)	160914	50.0000	49.94
	31 1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	257606	50.0000	51.32
	32 2-Methylphenol	108	4.255	4.255	(1.076)	218610	50.0000	49.86
	33 2,2'-oxybis(1-Chloropropane)	45	4.297	4.297	(1.086)	349371	50.0000	50.12
	34 4-Methylphenol	108	4.421	4.421	(1.118)	233354	50.0000	50.11
	36 Hexachloroethane	117	4.504	4.504	(1.139)	94106	50.0000	50.62
	37 N-Nitrosodipropylamine	70	4.442	4.442	(1.123)	156914	50.0000	50.59
	42 Nitrobenzene	77	4.597	4.597	(0.855)	219387	50.0000	49.95
	44 Isophorone	82	4.856	4.856	(0.904)	420061	50.0000	49.74
	45 2-Nitrophenol	139	4.960	4.960	(0.923)	132771	50.0000	50.95
	46 2,4-Dimethylphenol	107	5.012	5.012	(0.933)	231517	50.0000	50.00

10-3-10

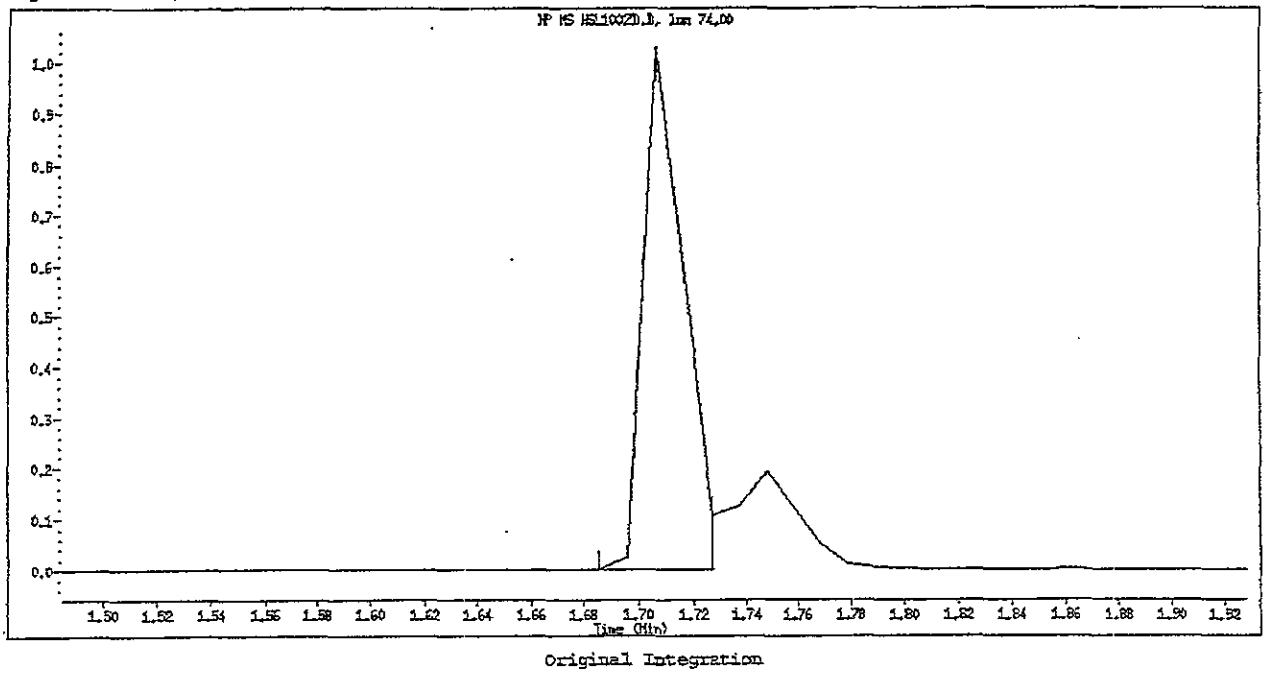
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.954)	253648	50.0000	49.15
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	179296	50.0000	50.05
50	Benzoic Acid	122	5.115	5.115	(0.952)	128366	50.0000	50.08
51	1,2,4-Trichlorobenzene	180	5.322	5.322	(0.990)	197265	50.0000	50.86
52	Naphthalene	128	5.395	5.395	(1.004)	726980	50.0000	49.49
54	4-Chloroaniline	127	5.488	5.488	(1.021)	291184	50.0000	50.72
57	Hexachlorobutadiene	225	5.613	5.613	(1.044)	95592	50.0000	50.36
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	205388	50.0000	51.34
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	464646	50.0000	50.50
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	106908	50.0000	49.76
69	2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	113001	50.0000	50.13
70	2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	128196	50.0000	52.79
71	2-Chloronaphthalene	162	6.784	6.784	(0.908)	403257	50.0000	50.72
73	2-Nitroaniline	65	6.949	6.949	(0.931)	124335	50.0000	51.59
76	Dimethylphthalate	163	7.229	7.229	(0.968)	475258	50.0000	51.91
77	Acenaphthylene	152	7.281	7.281	(0.975)	712158	50.0000	51.43
79	2,6-Dinitrotoluene	165	7.302	7.302	(0.978)	110261	50.0000	51.69
80	3-Nitroaniline	138	7.447	7.447	(0.997)	141396	50.0000	53.11
81	Acenaphthene	153	7.509	7.509	(1.006)	448691	50.0000	50.90
82	2,4-Dinitrophenol	184	7.571	7.572	(1.014)	58864	50.0000	47.37
83	Dibenzofuran	168	7.706	7.706	(1.032)	598735	50.0000	51.18
84	4-Nitrophenol	109	7.675	7.675	(1.028)	56777	50.0000	51.41
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	148875	50.0000	53.18
91	Fluorene	166	8.131	8.131	(1.089)	496097	50.0000	51.01
92	Diethylphthalate	149	8.100	8.100	(1.085)	487067	50.0000	51.96
93	4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	209308	50.0000	51.97
94	4-Nitroaniline	138	8.214	8.214	(1.100)	135397	50.0000	51.31
97	4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	76137	50.0000	46.58
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	409656	58.6000	58.41
100	Azobenzene	77	8.348	8.348	(0.888)	459950	50.0000	50.55
101	4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	115283	50.0000	51.04
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	124963	50.0000	49.54
110	Pentachlorophenol	266	9.240	9.240	(0.982)	67882	50.0000	45.48
114	Phenanthrene	178	9.437	9.437	(1.003)	718164	50.0000	49.24
115	Anthracene	178	9.499	9.499	(1.010)	728681	50.0000	50.01
118	Carbazole	167	9.768	9.768	(1.039)	660885	50.0000	49.65
120	Di-n-Butylphthalate	149	10.463	10.463	(1.112)	799142	50.0000	49.90
126	Fluoranthene	202	11.302	11.302	(1.202)	639252	50.0000	48.92
127	Benzidine	184	11.571	11.571	(0.840)	450332	50.0000	50.98
128	Pyrene	202	11.665	11.665	(0.847)	701084	50.0000	51.46
134	3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	385489	50.0000	49.44
136	Butylbenzylphthalate	149	12.991	12.991	(0.943)	340978	50.0000	49.94
138	Benzo(a)Anthracene	228	13.758	13.758	(0.998)	565271	50.0000	49.03
139	Chrysene	228	13.831	13.831	(1.004)	597685	50.0000	50.33
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	217413	50.0000	49.65
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.024)	464144	50.0000	49.35
142	Di-n-octylphthalate	149	15.167	15.167	(1.101)	732406	50.0000	48.72
144	Benzo(b)fluoranthene	252	15.582	15.582	(0.964)	527487	50.0000	55.18
145	Benzo(k)fluoranthene	252	15.623	15.623	(0.967)	580084	50.0000	47.27
147	Benzo(e)pyrene	252	16.007	16.007	(0.990)	506622	50.0000	50.82
148	Benzo(a)pyrene	252	16.079	16.079	(0.995)	542578	50.0000	50.06
151	Indeno(1,2,3-cd)pyrene	276	17.800	17.800	(1.101)	447085	50.0000	51.00(M)
152	Dibenzo(a,h)anthracene	278	17.841	17.841	(1.104)	486893	50.0000	49.72
153	Benzo(g,h,i)perylene	276	18.235	18.235	(1.128)	527720	50.0000	49.77

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				1107571	50.0000	50.74 (A)

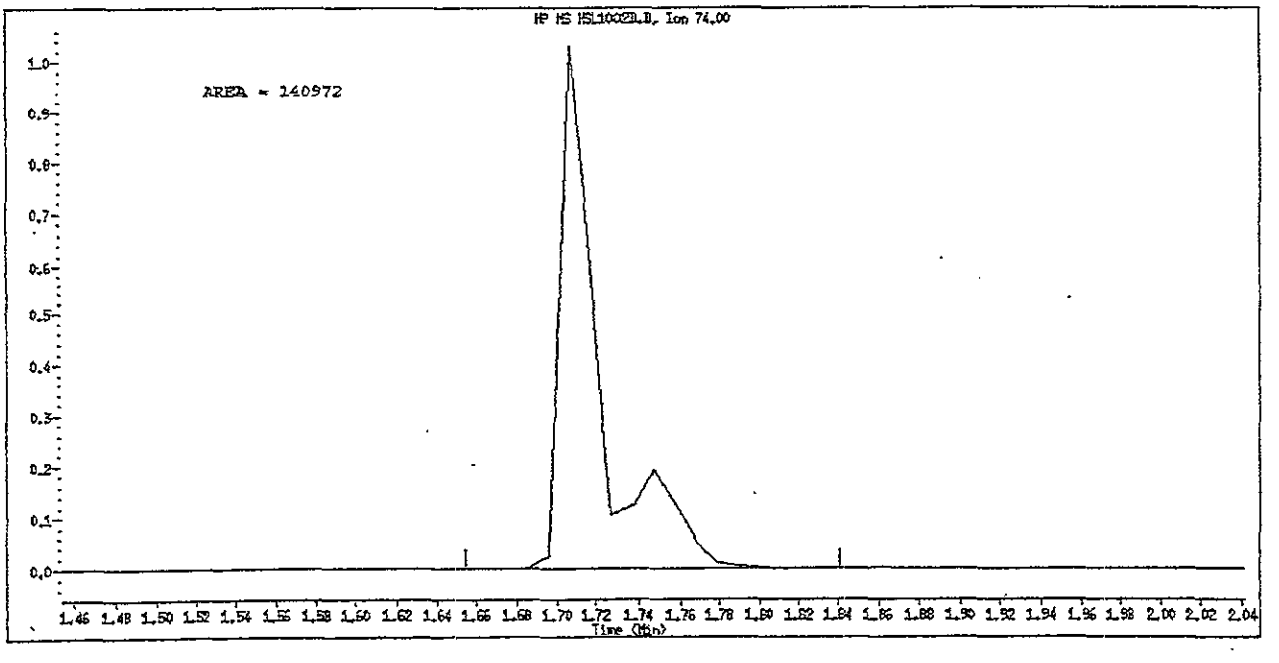
QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
M - Compound response manually integrated.

Data File Name: ESL1002D.D
Inj. Date and Time: 02-OCT-2010 13:44
Instrument ID: sv5.i
Client ID: 8270F.M
Compound Name: N-Nitrosodimethylamine
CAS #: 62-75-9
Report Date: 10/03/2010



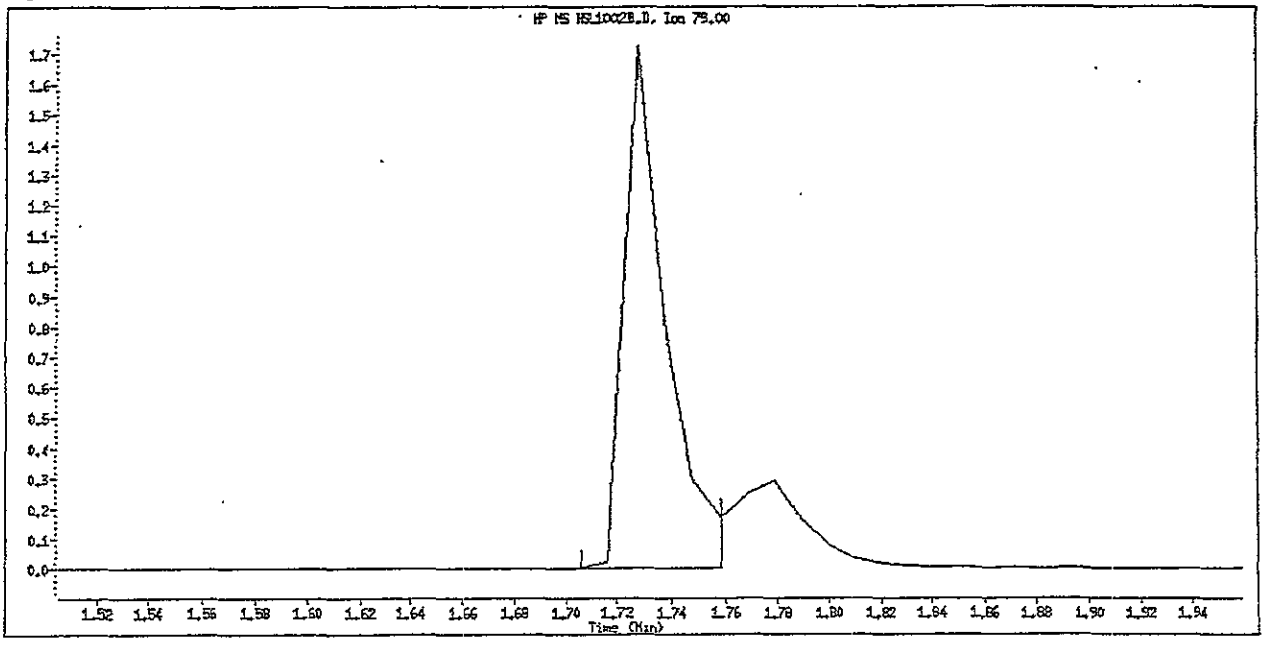
Original Integration



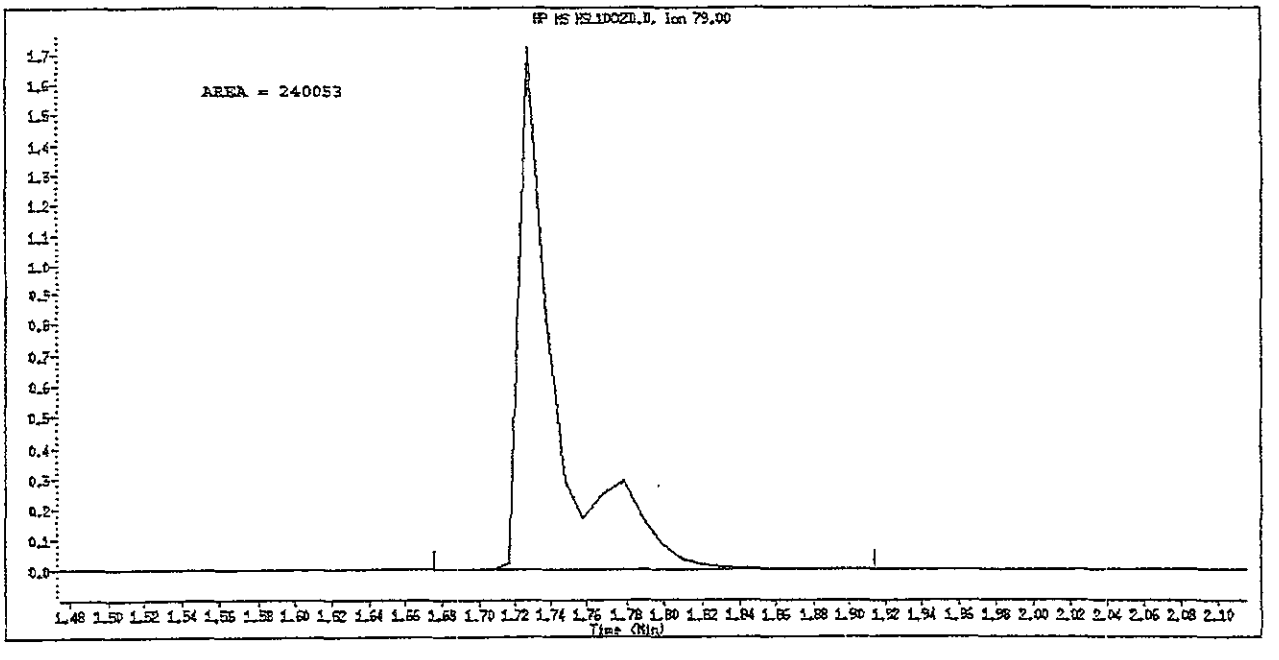
Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Poor Chromatography

Data File Name: HSL1002D.D
Inj. Date and Time: 02-OCT-2010 13:44
Instrument ID: sv5.i
Client ID: 8270F.M
Compound Name: Pyridine
CAS #: 110-86-1
Report Date: 10/03/2010



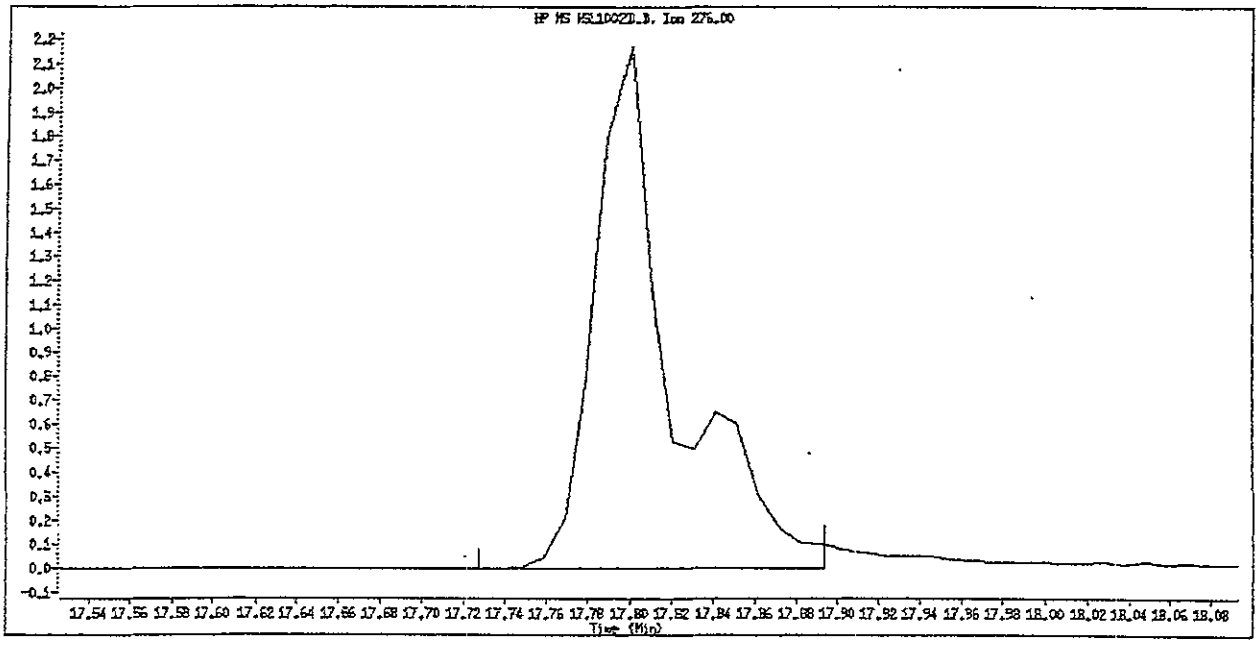
Original Integration



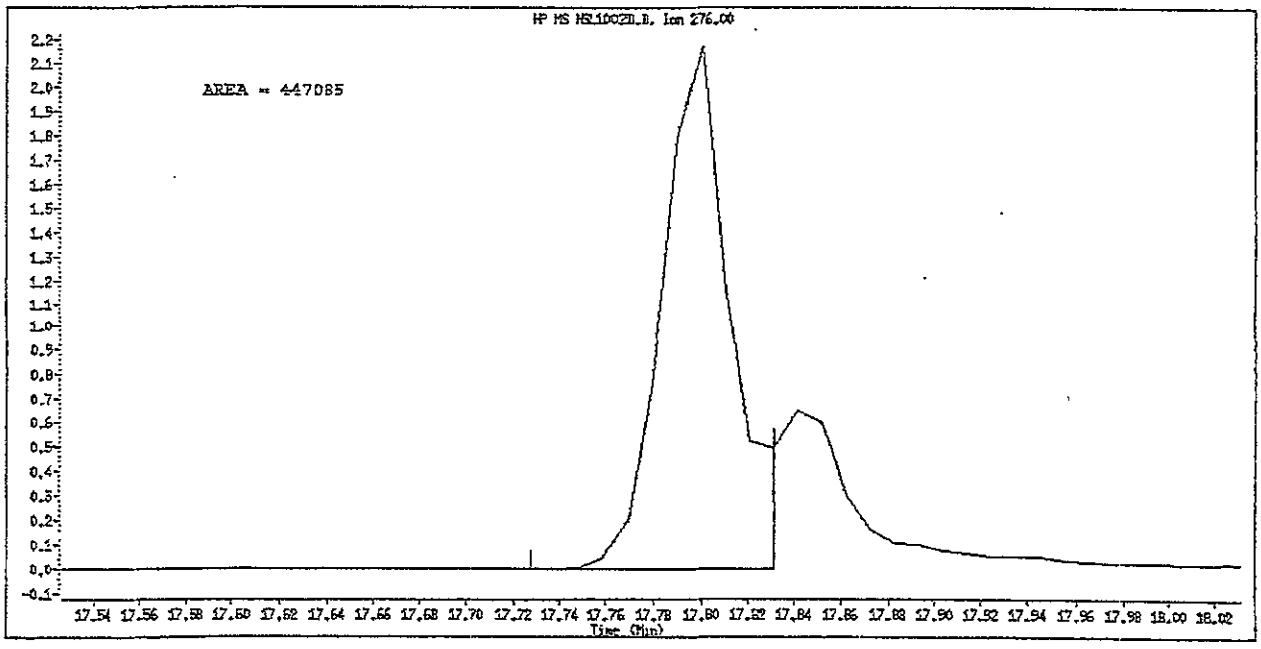
Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Poor Chromatography

Data File Name: ESL1002D.D
Inj. Date and Time: 02-OCT-2010 13:44
Instrument ID: sv5.i
Client ID: E270F.M
Compound Name: Indeno(1,2,3-cd)pyrene
CAS #: 193-39-5
Report Date: 10/03/2010



Original Integration



Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Poor Chromatography

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\100210.B\HSL1002D.D
Lab Smp Id: HSL_050 ug/ml CS-4 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 13:44
Operator : KT Inst ID: sv5.i
Smp Info : HSL_050 ug/ml CS-4;1;;4;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0310;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Meth Date : 02-Oct-2010 16:57 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 4 Calibration Sample, Level: 4
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
* 1	1,4-Dichlorobenzene-d4	152	3.955	3.955	(1.000)	122625	40.0000	
* 2	Naphthalene-d8	136	5.374	5.374	(1.000)	530514	40.0000	
* 3	Acenaphthene-d10	164	7.468	7.468	(1.000)	282538	40.0000	
* 4	Phenanthrene-d10	188	9.405	9.405	(1.000)	462722	40.0000	
* 5	Chrysene-d12	240	13.779	13.779	(1.000)	435850	40.0000	
* 6	Perylene-d12	264	16.162	16.162	(1.000)	422284	40.0000	
$ 7	2-Fluorophenol	112	2.732	2.732	(0.691)	220986	50.0000	48.83
$ 8	Phenol-d5	99	3.613	3.613	(0.914)	274382	50.0000	47.67
$ 9	2-Chlorophenol-d4	132	3.758	3.758	(0.950)	264352	50.0000	49.80
$ 10	1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	151616	50.0000	49.50
$ 11	Nitrobenzene-d5	82	4.576	4.576	(0.852)	226162	50.0000	48.07
$ 12	2-Fluorobiphenyl	172	6.680	6.680	(0.895)	473578	50.0000	52.38
$ 13	2,4,6-Tribromophenol	330	8.473	8.473	(1.135)	63311	50.0000	56.75
$ 14	Terphenyl-d14	244	12.017	12.017	(0.872)	436253	50.0000	51.56
15	N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	105836	50.0000	35.43
16	Pyridine	79	1.726	1.726	(0.437)	182664	50.0000	36.70
23	Aniline	93	3.654	3.654	(0.924)	346504	50.0000	48.01
24	Phenol	94	3.623	3.623	(0.916)	311820	50.0000	50.94
26	Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	220455	50.0000	47.40
27	2-Chlorophenol	128	3.768	3.768	(0.953)	242442	50.0000	50.05
28	1,3-Dichlorobenzene	146	3.923	3.923	(0.992)	265384	50.0000	49.58
29	1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	271151	50.0000	50.11
30	Benzyl Alcohol	108	4.120	4.120	(1.042)	160914	50.0000	48.35
31	1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	257606	50.0000	50.23
32	2-Methylphenol	108	4.255	4.255	(1.076)	218610	50.0000	48.31
33	2,2'-oxybis(1-Chloropropane)	45	4.297	4.297	(1.086)	349371	50.0000	40.48
34	4-Methylphenol	108	4.421	4.421	(1.118)	233354	50.0000	48.39
36	Hexachloroethane	117	4.504	4.504	(1.139)	94106	50.0000	49.29
37	N-Nitrosodinpropylamine	70	4.442	4.442	(1.123)	156914	50.0000	46.48
42	Nitrobenzene	77	4.597	4.597	(0.855)	219387	50.0000	46.91
44	Isophorone	82	4.856	4.856	(0.904)	420061	50.0000	47.38
45	2-Nitrophenol	139	4.960	4.960	(0.923)	132771	50.0000	52.00
46	2,4-Dimethylphenol	107	5.012	5.012	(0.933)	231517	50.0000	48.84

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.954)	253648	50.0000	48.05
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	179296	50.0000	51.39
50	Benzoic Acid	122	5.115	5.115	(0.952)	128366	50.0000	48.58
51	1,2,4-Trichlorobenzene	180	5.322	5.322	(0.990)	197265	50.0000	52.15
52	Naphthalene	128	5.395	5.395	(1.004)	724980	50.0000	49.10
54	4-Chloroaniline	127	5.488	5.488	(1.021)	291184	50.0000	50.12
57	Hexachlorobutadiene	225	5.613	5.613	(1.044)	95592	50.0000	53.11
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	205388	50.0000	51.16
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	464646	50.0000	51.57
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	104908	50.0000	48.68
69	2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	113001	50.0000	52.83
70	2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	128196	50.0000	54.56
71	2-Chloronaphthalene	162	6.784	6.784	(0.908)	403257	50.0000	50.91
73	2-Nitroaniline	65	6.949	6.949	(0.931)	124335	50.0000	46.87
76	Dimethylphthalate	163	7.229	7.229	(0.968)	475258	50.0000	51.95
77	Acenaphthylene	152	7.281	7.281	(0.975)	712158	50.0000	51.43
79	2,6-Dinitrotoluene	165	7.302	7.302	(0.978)	110261	50.0000	53.82
80	3-Nitroaniline	138	7.447	7.447	(0.997)	141396	50.0000	52.05
81	Acenaphthene	153	7.509	7.509	(1.006)	448691	50.0000	50.85
82	2,4-Dinitrophenol	184	7.571	7.571	(1.014)	58864	50.0000	48.70
83	Dibenzofuran	168	7.706	7.706	(1.032)	598735	50.0000	51.36
84	4-Nitrophenol	109	7.675	7.675	(1.028)	56777	50.0000	47.87
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	148875	50.0000	54.24
91	Fluorene	166	8.131	8.131	(1.089)	494097	50.0000	51.73
92	Diethylphthalate	149	8.100	8.100	(1.085)	487067	50.0000	50.93
93	4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	209308	50.0000	53.15
94	4-Nitroaniline	138	8.214	8.214	(1.100)	135397	50.0000	50.99
97	4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	76137	50.0000	46.45
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	409666	58.6000	56.82
100	Azobenzene	77	8.348	8.348	(0.888)	459960	50.0000	45.85
101	4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	115283	50.0000	51.56
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	124963	50.0000	51.74
110	Pentachlorophenol	266	9.240	9.240	(0.982)	67882	50.0000	46.83
114	Phenanthrene	178	9.437	9.437	(1.003)	718164	50.0000	49.76
115	Anthracene	178	9.499	9.499	(1.010)	728681	50.0000	50.17
118	Carbazole	167	9.768	9.768	(1.039)	660885	50.0000	48.92
120	Di-n-Butylphthalate	149	10.463	10.463	(1.112)	799142	50.0000	48.91
126	Fluoranthene	202	11.302	11.302	(1.202)	639252	50.0000	49.21
127	Benzidine	184	11.571	11.571	(0.840)	450332	50.0000	50.32
128	Pyrene	202	11.665	11.665	(0.847)	701084	50.0000	51.44
134	3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	385489	50.0000	50.19
136	Butylbenzylphthalate	149	12.991	12.991	(0.943)	340978	50.0000	48.97
138	Benzo(a)Anthracene	228	13.758	13.758	(0.998)	569271	50.0000	49.51
139	Chrysene	228	13.831	13.831	(1.004)	597685	50.0000	50.03
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	217413	50.0000	51.67
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.024)	464144	50.0000	48.41
142	Di-n-octylphthalate	149	15.167	15.167	(1.101)	732406	50.0000	47.78
144	Benzo(b)fluoranthene	252	15.582	15.582	(0.964)	527487	50.0000	52.68
145	Benzo(k)fluoranthene	252	15.623	15.623	(0.967)	580084	50.0000	49.57
147	Benzo(e)pyrene	252	16.007	16.007	(0.990)	506622	50.0000	51.04
148	Benzo(a)pyrene	252	16.079	16.079	(0.995)	542578	50.0000	49.78
151	Indeno(1,2,3-cd)pyrene	276	17.800	17.800	(1.101)	564014	50.0000	58.49
152	Dibenzo(a,h)anthracene	278	17.841	17.841	(1.104)	486893	50.0000	49.27
153	Benzo(g,h,i)perylene	276	18.235	18.235	(1.128)	527720	50.0000	49.81

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				1107571	50.0000	51.00 (A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.

Data File: \\sv5\c\chem\sv5.i\100210.B\HSL1002D.D Page 1
Report Date: 03-Oct-2010 11:14

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002D.D
Lab Smp Id: HSL_050 ug/ml CS-4
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\sv5\c\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0310;0;8270F.M

Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	122625	0.00
2 Naphthalene-d8	530514	265257	1061028	530514	0.00
3 Acenaphthene-d10	282538	141269	565076	282538	0.00
4 Phenanthrene-d10	462722	231361	925444	462722	0.00
5 Chrysene-d12	435850	217925	871700	435850	0.00
6 Perylene-d12	422284	211142	844568	422284	0.00

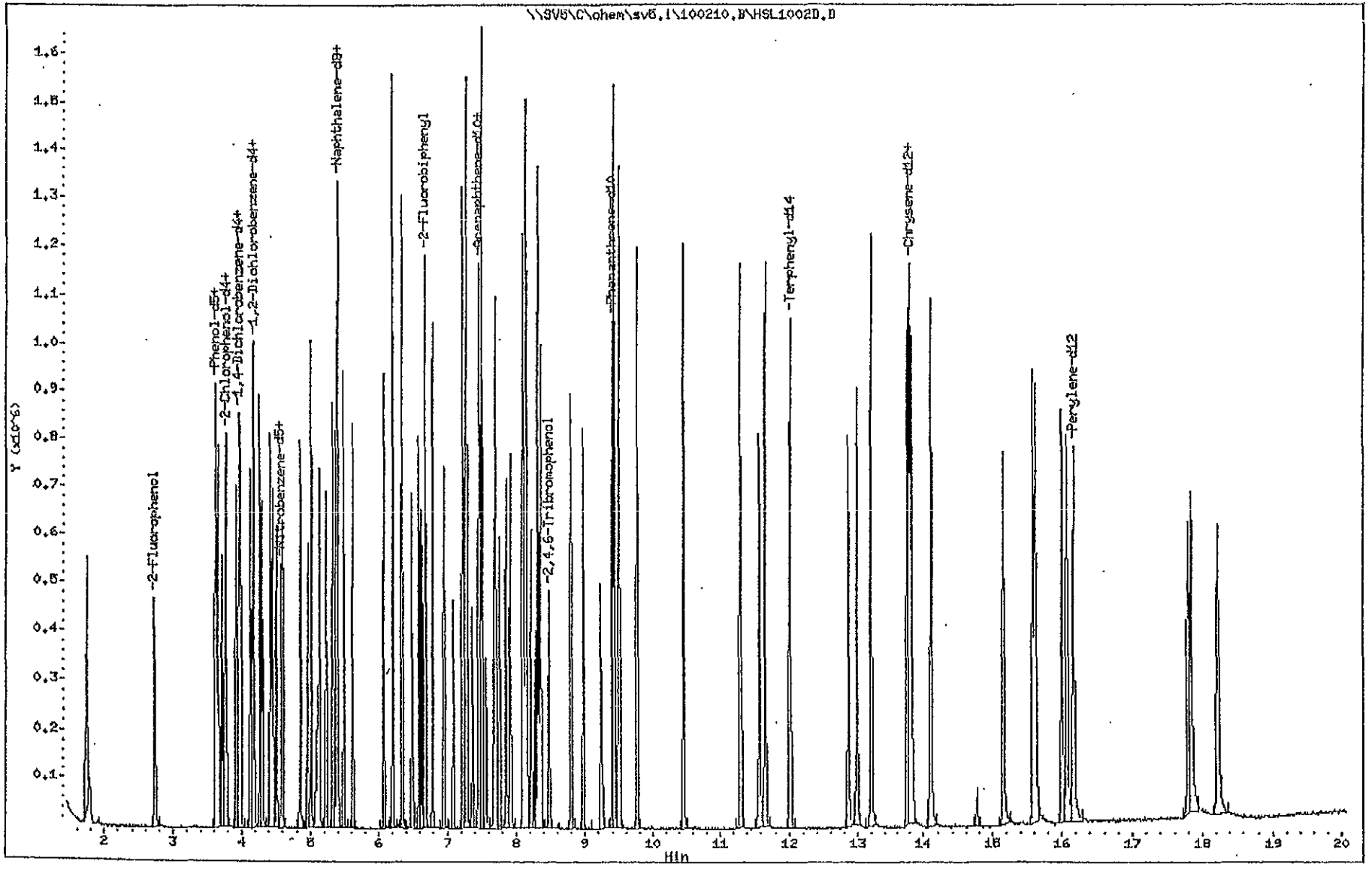
COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.96	3.46	4.46	3.96	0.00
2 Naphthalene-d8	5.37	4.87	5.87	5.37	0.00
3 Acenaphthene-d10	7.47	6.97	7.97	7.47	0.00
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	0.00
5 Chrysene-d12	13.78	13.28	14.28	13.78	0.00
6 Perylene-d12	16.16	15.66	16.66	16.16	0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV5\C\chem\sv5.i\100210.B\HSL1002D.D
Date : 02-OCT-2010 13:44
Client ID: 8270F,M
Sample Info: HSL_050 ug/ml CS-4;1;;4;;;;4

Instrument: sv5.i
Operator: KT
Column diameter: 2.00

Column phase:



TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002E.D
Lab Smp Id: HSL_080 ug/ml CS-5 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 14:09
Operator : KT Inst ID: sv5.i
Smp Info : HSL_080 ug/ml CS-5;1;;5;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0311;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:09 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 5 Calibration Sample, Level: 5
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SACP307UM

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
*	1 1,4-Dichlorobenzene-d4	152	3.954	3.955	(1.000)	126989	40.0000		(q)
*	2 Naphthalene-d8	136	5.374	5.374	(1.000)	553454	40.0000		
*	3 Acenaphthene-d10	164	7.468	7.468	(1.000)	300315	40.0000		
*	4 Phenanthrene-d10	188	9.405	9.405	(1.000)	477777	40.0000		
*	5 Chrysene-d12	240	13.789	13.779	(1.000)	486126	40.0000		
*	6 Perylene-d12	264	16.162	16.162	(1.000)	482782	40.0000		
$	7 2-Fluorophenol	112	2.742	2.732	(0.693)	364547	80.0000	81.44	
$	8 Phenol-d5	99	3.612	3.613	(0.914)	459352	80.0000	81.61	
$	9 2-Chlorophenol-d4	132	3.758	3.758	(0.950)	399981	80.0000	80.92	
$	10 1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	252754	80.0000	80.82	
$	11 Nitrobenzene-d5	82	4.587	4.576	(0.853)	371989	80.0000	79.35	
$	12 2-Fluorobiphenyl	172	6.680	6.680	(0.895)	755916	80.0000	78.14	
$	13 2,4,6-Tribromophenol	330	8.483	8.473	(1.136)	107063	80.0000	82.04	
$	14 Terphenyl-d14	244	12.017	12.017	(0.871)	758812	80.0000	79.25	
	15 N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	236570	80.0000	80.86	(q)
	16 Pyridine	79	1.726	1.726	(0.437)	386806	80.0000	79.06	(Q)
	23 Aniline	93	3.654	3.654	(0.924)	583513	80.0000	81.44	(Q)
	24 Phenol	94	3.623	3.623	(0.916)	524930	80.0000	81.16	(Q)
	26 Bis(2-chloroethyl)ether	93	3.716	3.715	(0.940)	362044	80.0000	79.83	
	27 2-Chlorophenol	128	3.768	3.768	(0.953)	398210	80.0000	80.21	
	28 1,3-Dichlorobenzene	146	3.923	3.923	(0.992)	428311	80.0000	79.20	
	29 1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	452588	80.0000	80.04	
	30 Benzyl Alcohol	108	4.120	4.120	(1.042)	273768	80.0000	82.05	
	31 1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	415025	80.0000	79.84	
	32 2-Methylphenol	108	4.255	4.255	(1.076)	369704	80.0000	81.43	
	33 2,2'-oxybis(1-Chloropropane)	45	4.296	4.297	(1.086)	576575	80.0000	79.88	
	34 4-Methylphenol	108	4.421	4.421	(1.118)	387704	80.0000	80.39	
	36 Hexachloroethane	117	4.504	4.504	(1.139)	153472	80.0000	79.72	
	37 N-Nitrosodipropylamine	70	4.442	4.442	(1.123)	265916	80.0000	82.78	
	42 Nitrobenzene	77	4.597	4.597	(0.855)	369479	80.0000	80.64	
	44 Isophorone	82	4.856	4.856	(0.904)	704520	80.0000	79.96	
	45 2-Nitrophenol	139	4.960	4.960	(0.923)	221628	80.0000	81.52	
	46 2,4-Dimethylphenol	107	5.011	5.012	(0.933)	385073	80.0000	79.72	

10-3-10

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.125	5.126	(0.954)	426158	80.0000	79.16
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	301897	80.0000	80.78
50	Benzoic Acid	122	5.125	5.115	(0.954)	232711	80.0000	87.04
51	1,2,4-Trichlorobenzene	180	5.322	5.322	(0.990)	323096	80.0000	79.84
52	Naphthalene	128	5.395	5.395	(1.004)	1216155	80.0000	79.58
54	4-Chloroaniline	127	5.488	5.488	(1.021)	484619	80.0000	80.91
57	Hexachlorobutadiene	225	5.613	5.613	(1.044)	159233	80.0000	80.41
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	335335	80.0000	80.35
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	781029	80.0000	81.36
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	181608	80.0000	81.05
69	2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	194036	80.0000	80.98
70	2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	211635	80.0000	81.99
71	2-Chloronaphthalene	162	6.784	6.784	(0.908)	668023	80.0000	79.04
73	2-Nitroaniline	65	6.949	6.949	(0.931)	209144	80.0000	81.65
76	Dimethylphthalate	163	7.229	7.229	(0.968)	787815	80.0000	80.96
77	Acenaphthylene	152	7.281	7.281	(0.975)	1190475	80.0000	80.88
79	2,6-Dinitrotoluene	165	7.302	7.302	(0.978)	187961	80.0000	82.91
80	3-Nitroaniline	138	7.457	7.447	(0.999)	232287	80.0000	82.09
81	Acenaphthene	153	7.509	7.509	(1.006)	727612	80.0000	77.66
82	2,4-Dinitrophenol	184	7.571	7.572	(1.014)	110384	80.0000	78.64
83	Dibenzofuran	168	7.706	7.706	(1.032)	991740	80.0000	79.76(Q)
84	4-Nitrophenol	109	7.675	7.675	(1.028)	102888	80.0000	87.65(Q)
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	246471	80.0000	82.83
91	Fluorene	166	8.131	8.131	(1.089)	834271	80.0000	81.03
92	Diethylphthalate	149	8.100	8.100	(1.085)	792071	80.0000	79.50
93	4-Chlorophenyl-phenylether	204	8.151	8.152	(1.092)	340608	80.0000	79.56
94	4-Nitroaniline	138	8.224	8.214	(1.101)	235541	80.0000	83.97
97	4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	134784	80.0000	76.76
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	695826	93.7000	96.08
100	Azobenzene	77	8.348	8.348	(0.888)	765053	80.0000	81.43
101	4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	187352	80.0000	80.33
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	207655	80.0000	79.72
110	Pentachlorophenol	266	9.240	9.240	(0.982)	126397	80.0000	78.86
114	Phenanthrene	178	9.437	9.437	(1.003)	1188468	80.0000	78.92
115	Anthracene	178	9.509	9.499	(1.011)	1218608	80.0000	81.00
118	Carbazole	167	9.768	9.768	(1.039)	1118637	80.0000	81.39
120	Di-n-Butylphthalate	149	10.462	10.463	(1.112)	1351860	80.0000	81.75
126	Fluoranthene	202	11.302	11.302	(1.202)	1107116	80.0000	82.05
127	Benzidine	184	11.571	11.571	(0.839)	799205	80.0000	81.12
128	Pyrene	202	11.665	11.665	(0.846)	1221015	80.0000	80.36
134	3,3'-dimethylbenzidine	212	12.867	12.867	(0.933)	715866	80.0000	82.31
136	Butylbenzylphthalate	149	12.991	12.991	(0.942)	598812	80.0000	78.63
138	Benzo(a)Anthracene	228	13.758	13.758	(0.998)	1034950	80.0000	79.92
139	Chrysene	228	13.830	13.831	(1.003)	1040163	80.0000	78.52
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.001)	392335	80.0000	80.33
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.023)	820296	80.0000	78.20
142	Di-n-octylphthalate	149	15.167	15.167	(1.100)	1354893	80.0000	80.80
144	Benzo(b)fluoranthene	252	15.582	15.582	(0.964)	920884	80.0000	84.26(Q)
145	Benzo(k)fluoranthene	252	15.623	15.623	(0.967)	1102899	80.0000	78.61(Q)
147	Benzo(e)pyrene	252	16.007	16.007	(0.990)	936566	80.0000	82.18
148	Benzo(a)pyrene	252	16.079	16.079	(0.995)	1039045	80.0000	83.86
151	Indeno(1,2,3-cd)pyrene	276	17.799	17.800	(1.101)	811625	80.0000	80.99
152	Dibenzo(a,h)anthracene	278	17.851	17.841	(1.105)	926841	80.0000	82.79
153	Benzo(g,h,i)perylene	276	18.235	18.235	(1.128)	982275	80.0000	81.04

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				2023783	80.0000	81.09 (A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\100210.B\HSL1002E.D
Lab Smp Id: HSL 080 ug/ml CS-5 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 14:09
Operator : KT Inst ID: sv5.i
Smp Info : HSL 080 ug/ml CS-5;1;;5;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0311;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Meth Date : 02-Oct-2010 16:57 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 5 Calibration Sample, Level: 5
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
* 1	1,4-Dichlorobenzene-d4	152	3.954	3.955	(1.000)	126989	40.0000		(q)
* 2	Naphthalene-d8	136	5.374	5.374	(1.000)	553454	40.0000		
* 3	Acenaphthene-d10	164	7.468	7.468	(1.000)	300315	40.0000		
* 4	Phenanthrene-d10	188	9.405	9.405	(1.000)	477777	40.0000		
* 5	Chrysene-d12	240	13.789	13.779	(1.000)	486126	40.0000		
* 6	Perylene-d12	264	16.162	16.162	(1.000)	482782	40.0000		
$ 7	2-Fluorophenol	112	2.742	2.732	(0.693)	364547	80.0000	77.78	
$ 8	Phenol-d5	99	3.612	3.613	(0.914)	459352	80.0000	77.07	
$ 9	2-Chlorophenol-d4	132	3.758	3.758	(0.950)	399981	80.0000	78.71	
$ 10	1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	252754	80.0000	79.68	
$ 11	Nitrobenzene-d5	82	4.587	4.576	(0.853)	371989	80.0000	75.79	
$ 12	2-Fluorobiphenyl	172	6.680	6.680	(0.895)	755916	80.0000	78.58	
$ 13	2,4,6-Tribromophenol	330	8.483	8.473	(1.136)	107063	80.0000	90.29	
$ 14	Terphenyl-d14	244	12.017	12.017	(0.871)	758812	80.0000	80.04	
15	N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	236570	80.0000	76.48	
16	Pyridine	79	1.726	1.726	(0.437)	386806	80.0000	75.04	
23	Aniline	93	3.654	3.654	(0.924)	583513	80.0000	78.07	(Q)
24	Phenol	94	3.623	3.623	(0.916)	524930	80.0000	82.81	(Q)
26	Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	362044	80.0000	75.18	
27	2-Chlorophenol	128	3.768	3.768	(0.953)	398210	80.0000	79.39	
28	1,3-Dichlorobenzene	146	3.923	3.923	(0.992)	428311	80.0000	77.27	
29	1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	452588	80.0000	80.76	
30	Benzyl Alcohol	108	4.120	4.120	(1.042)	273768	80.0000	79.43	
31	1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	415025	80.0000	78.14	
32	2-Methylphenol	108	4.255	4.255	(1.076)	369704	80.0000	78.90	
33	2,2'-oxybis(1-Chloropropane)	45	4.296	4.297	(1.086)	576575	80.0000	64.50	
34	4-Methylphenol	108	4.421	4.421	(1.118)	387704	80.0000	77.63	
36	Hexachloroethane	117	4.504	4.504	(1.139)	153472	80.0000	77.62	
37	N-Nitrosodinpropylamine	70	4.442	4.442	(1.123)	265916	80.0000	76.06	
42	Nitrobenzene	77	4.597	4.597	(0.855)	369479	80.0000	75.74	
44	Isophorone	82	4.856	4.856	(0.904)	704520	80.0000	76.17	
45	2-Nitrophenol	139	4.960	4.960	(0.923)	221628	80.0000	83.21	
46	2,4-Dimethyphenol	107	5.011	5.012	(0.933)	385073	80.0000	77.86	

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.125	5.126	(0.954)	426158	80.0000	77.39
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	301897	80.0000	82.94
50	Benzoic Acid	122	5.125	5.115	(0.954)	232711	80.0000	84.41
51	1,2,4-Trichlorobenzene	180	5.322	5.322	(0.990)	323096	80.0000	81.88
52	Naphthalene	128	5.395	5.395	(1.004)	1216155	80.0000	78.94
54	4-Chloroaniline	127	5.488	5.488	(1.021)	484619	80.0000	79.97
57	Hexachlorobutadiene	225	5.613	5.613	(1.044)	159233	80.0000	84.81
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	335335	80.0000	80.06
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	781029	80.0000	83.09
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	181608	80.0000	79.29
69	2,4,6-Trichlorophenol	196	6.576	6.576	(0.881)	194036	80.0000	85.34
70	2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	211635	80.0000	84.74
71	2-Chloronaphthalene	162	6.784	6.784	(0.908)	668023	80.0000	79.34
73	2-Nitroaniline	65	6.949	6.949	(0.931)	209144	80.0000	74.17
76	Dimethylphthalate	163	7.229	7.229	(0.968)	787815	80.0000	81.01
77	Acenaphthylene	152	7.281	7.281	(0.975)	1190475	80.0000	80.88
79	2,6-Dinitrotoluene	165	7.302	7.302	(0.978)	187961	80.0000	86.31
80	3-Nitroaniline	138	7.457	7.447	(0.999)	232287	80.0000	80.44
81	Acenaphthene	153	7.509	7.509	(1.006)	727612	80.0000	77.58
82	2,4-Dinitrophenol	184	7.571	7.571	(1.014)	110384	80.0000	81.10
83	Dibenzofuran	168	7.706	7.706	(1.032)	991740	80.0000	80.04(Q)
84	4-Nitrophenol	109	7.675	7.675	(1.028)	102888	80.0000	81.61(Q)
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	246471	80.0000	84.49
91	Fluorene	166	8.131	8.131	(1.089)	834271	80.0000	82.18
92	Diethylphthalate	149	8.100	8.100	(1.085)	792071	80.0000	77.92
93	4-Chlorophenyl-phenylether	204	8.151	8.152	(1.092)	340608	80.0000	81.38
94	4-Nitroaniline	138	8.224	8.214	(1.101)	235541	80.0000	83.45
97	4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	134784	80.0000	75.96
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	695826	93.7000	93.46
100	Azobenzene	77	8.348	8.348	(0.888)	765053	80.0000	73.86
101	4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	187352	80.0000	81.15
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	207655	80.0000	83.28
110	Pentachlorophenol	266	9.240	9.240	(0.982)	126397	80.0000	84.45
114	Phenanthrene	178	9.437	9.437	(1.003)	1188468	80.0000	79.75
115	Anthracene	178	9.509	9.499	(1.011)	1218608	80.0000	81.25
118	Carbazole	167	9.768	9.768	(1.039)	1118637	80.0000	80.19
120	Di-n-Butylphthalate	149	10.462	10.463	(1.112)	1351860	80.0000	80.14
126	Fluoranthene	202	11.302	11.302	(1.202)	1107116	80.0000	82.54
127	Benzidine	184	11.571	11.571	(0.839)	799205	80.0000	80.06
128	Pyrene	202	11.665	11.665	(0.846)	1221015	80.0000	80.33
134	3,3'-dimethylbenzidine	212	12.867	12.867	(0.933)	715866	80.0000	83.56
136	Butylbenzylphthalate	149	12.991	12.991	(0.942)	598812	80.0000	77.10
138	Benzo(a)Anthracene	228	13.758	13.758	(0.998)	1034950	80.0000	80.70
139	Chrysene	228	13.830	13.831	(1.003)	1040163	80.0000	78.06
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.001)	392335	80.0000	83.60
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.023)	820296	80.0000	76.71
142	Di-n-octylphthalate	149	15.167	15.167	(1.100)	1354893	80.0000	79.24
144	Benzo(b)fluoranthene	252	15.582	15.582	(0.964)	920884	80.0000	80.44(Q)
145	Benzo(k)fluoranthene	252	15.623	15.623	(0.967)	1102899	80.0000	82.44(Q)
147	Benzo(e)pyrene	252	16.007	16.007	(0.990)	936566	80.0000	82.53
148	Benzo(a)pyrene	252	16.079	16.079	(0.995)	1039045	80.0000	83.39
151	Indeno(1,2,3-cd)pyrene	276	17.799	17.800	(1.101)	811625	80.0000	73.62
152	Dibenzo(a,h)anthracene	278	17.851	17.841	(1.105)	926841	80.0000	82.04
153	Benzo(g,h,i)perylene	276	18.235	18.235	(1.128)	982275	80.0000	81.10

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	CW-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				2023783	80.0000	81.52(A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002E.D
Lab Smp Id: HSL_080 ug/ml CS-5
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\sv5\c\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0311;0;8270F.M

Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	126989	3.56
2 Naphthalene-d8	530514	265257	1061028	553454	4.32
3 Acenaphthene-d10	282538	141269	565076	300315	6.29
4 Phenanthrene-d10	462722	231361	925444	477777	3.25
5 Chrysene-d12	435850	217925	871700	486126	11.54
6 Perylene-d12	422284	211142	844568	482782	14.33

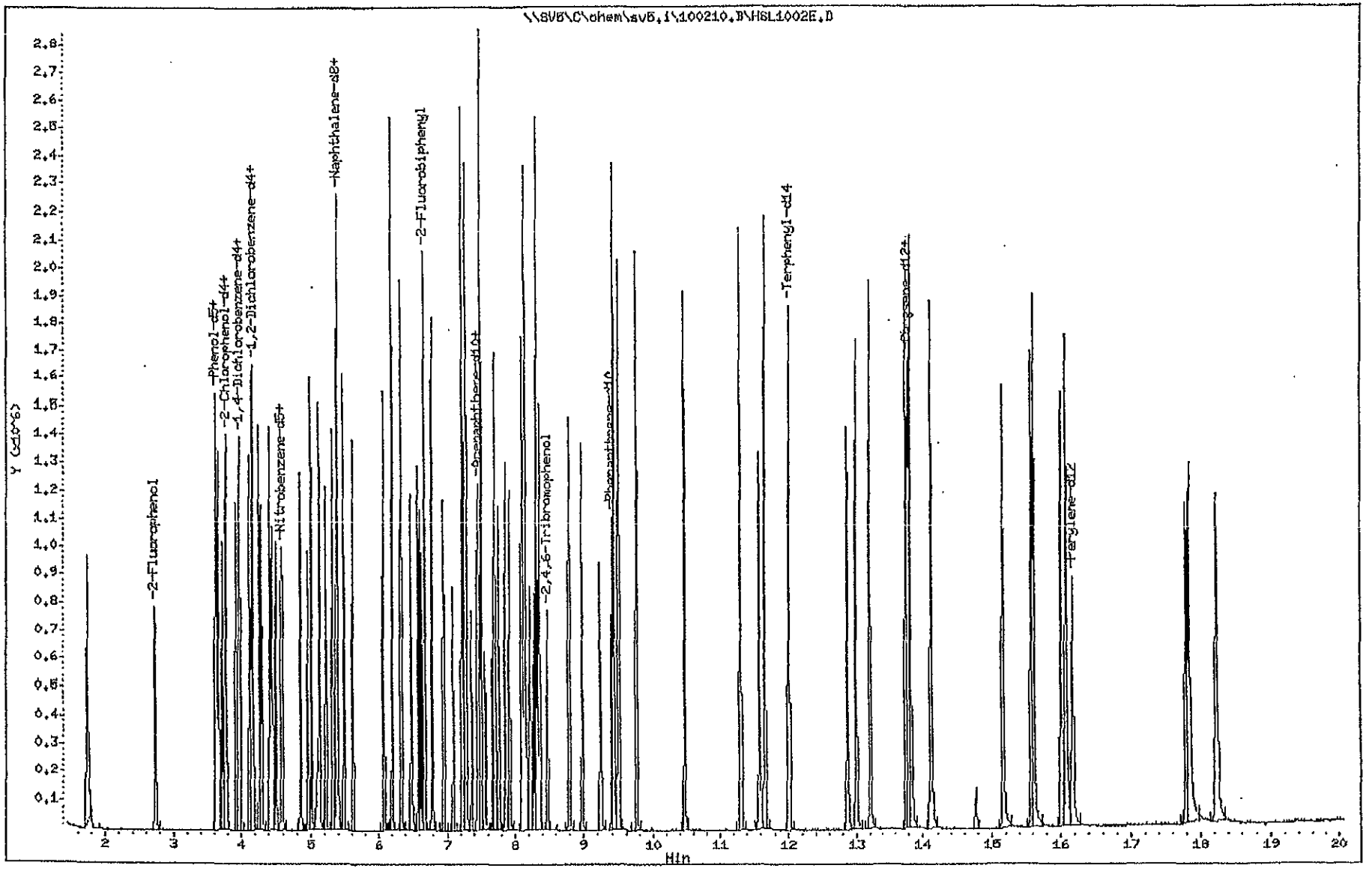
COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.96	3.46	4.46	3.95	-0.00
2 Naphthalene-d8	5.37	4.87	5.87	5.37	-0.00
3 Acenaphthene-d10	7.47	6.97	7.97	7.47	-0.00
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	-0.00
5 Chrysene-d12	13.78	13.28	14.28	13.79	0.07
6 Perylene-d12	16.16	15.66	16.66	16.16	-0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV8\C\chem\sv8.i\100210.B\HSL1002E.D
Date : 02-OCT-2010 14:09
Client ID: 8270F.H
Sample Info: HSL_080 ug/ml CS-5;1;;5;;;;;4

Instrument: sv8.i
Operator: KT
Column diameter: 2.00

Column phase:



TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002F.D
Lab Smp Id: HSL_120 ug/ml CS-6 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 14:35
Operator : KT Inst ID: sv5.i
Smp Info : HSL_120 ug/ml CS-6;1;;6;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0312;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:09 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 6 Calibration Sample, Level: 6
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SACP307UM

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
* 1	1,4-Dichlorobenzene-d4	152	3.955	3.955	(1.000)	137751	40.0000		(Q)
* 2	Naphthalene-d8	136	5.374	5.374	(1.000)	591665	40.0000		
* 3	Acenaphthene-d10	164	7.468	7.468	(1.000)	322596	40.0000		
* 4	Phenanthrene-d10	188	9.406	9.405	(1.000)	515607	40.0000		
* 5	Chrysene-d12	240	13.789	13.779	(1.000)	509570	40.0000		
* 6	Perylene-d12	264	16.173	16.162	(1.000)	539588	40.0000		
$ 7	2-Fluorophenol	112	2.732	2.732	(0.691)	588026	120.000	121.1	
$ 8	Phenol-d5	99	3.613	3.613	(0.914)	759824	120.000	124.4	
$ 9	2-Chlorophenol-d4	132	3.758	3.758	(0.950)	652805	120.000	121.7	
$ 10	1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	407247	120.000	120.0	
$ 11	Nitrobenzene-d5	82	4.587	4.576	(0.853)	623501	120.000	124.4	
$ 12	2-Fluorobiphenyl	172	6.680	6.680	(0.895)	1255441	120.000	120.8	
$ 13	2,4,6-Tribromophenol	330	8.483	8.473	(1.136)	179055	120.000	127.7	
$ 14	Terphenyl-d14	244	12.017	12.017	(0.871)	1251844	120.000	124.7	
15	N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	388111	120.000	122.3	(q)
16	Pyridine	79	1.727	1.726	(0.437)	633334	120.000	119.3	(Q)
23	Aniline	93	3.654	3.654	(0.924)	964533	120.000	124.1	(Q)
24	Phenol	94	3.623	3.623	(0.916)	851671	120.000	121.4	(Q)
26	Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	596323	120.000	121.2	
27	2-Chlorophenol	128	3.768	3.768	(0.953)	653244	120.000	121.3	
28	1,3-Dichlorobenzene	146	3.924	3.923	(0.992)	712032	120.000	121.4	
29	1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	740915	120.000	120.8	
30	Benzyl Alcohol	108	4.120	4.120	(1.042)	450249	120.000	124.4	
31	1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	679448	120.000	120.5	
32	2-Methylphenol	108	4.255	4.255	(1.076)	603987	120.000	122.6	
33	2,2'-oxybis(1-Chloropropane)	45	4.297	4.297	(1.086)	941514	120.000	120.2	
34	4-Methylphenol	108	4.421	4.421	(1.118)	664202	120.000	123.1	
36	Hexachloroethane	117	4.504	4.504	(1.139)	245394	120.000	117.5	
37	N-Nitrosodipropylamine	70	4.452	4.442	(1.126)	628242	120.000	122.9	
42	Nitrobenzene	77	4.607	4.597	(0.857)	593736	120.000	121.2	
44	Isophorone	82	4.867	4.856	(0.906)	1179801	120.000	125.2	
45	2-Nitrophenol	139	4.960	4.960	(0.923)	367467	120.000	126.4	
46	2,4-Dimethylphenol	107	5.012	5.012	(0.933)	638328	120.000	123.6	

10-3-10

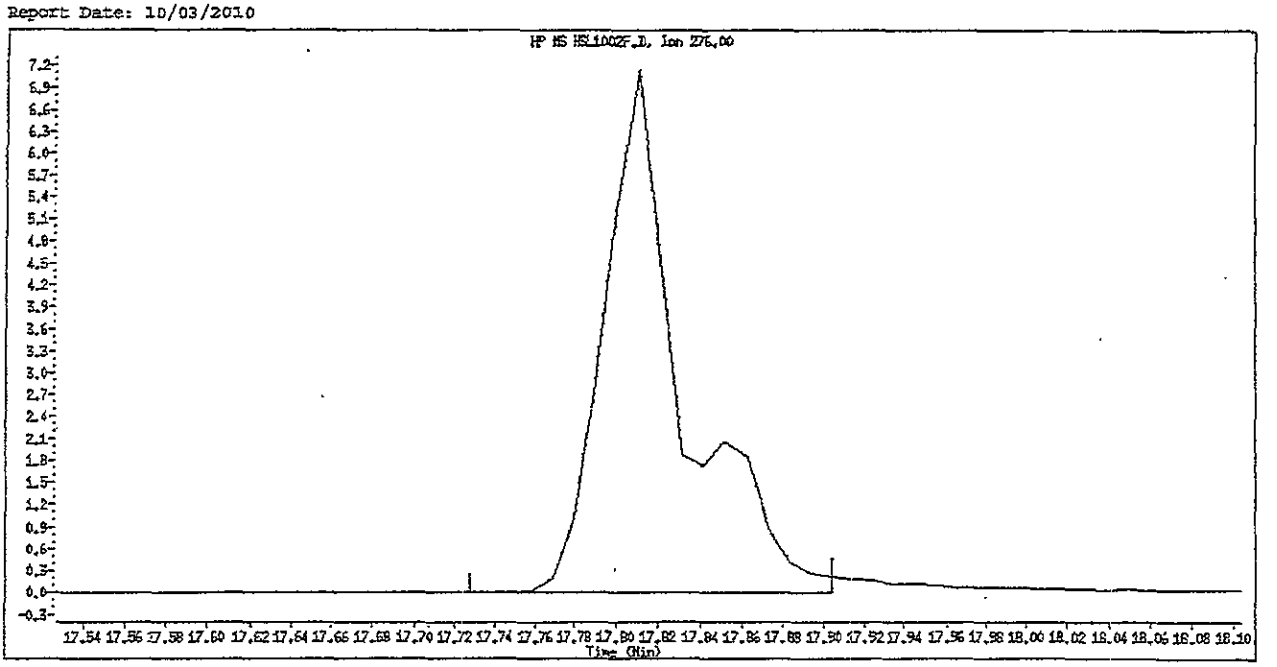
	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
47	Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.954)	707504	120.000	122.8	
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	500185	120.000	125.2	
50	Benzoic Acid	122	5.146	5.115	(0.958)	395333	120.000	138.3	
51	1,2,4-Trichlorobenzene	180	5.333	5.322	(0.992)	531764	120.000	122.9	
52	Naphthalene	128	5.395	5.395	(1.004)	2020315	120.000	123.7	
54	4-Chloroaniline	127	5.488	5.488	(1.021)	797064	120.000	124.5	
57	Hexachlorobutadiene	225	5.613	5.613	(1.044)	255231	120.000	120.6	
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	563840	120.000	126.4	
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	1263302	120.000	123.1	
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	312226	120.000	129.7	
69	2,4,6-Trichlorophenol	196	6.587	6.576	(0.882)	331223	120.000	128.7	
70	2,4,5-Trichlorophenol	196	6.628	6.628	(0.888)	343374	120.000	123.8	
71	2-Chloronaphthalene	162	6.784	6.784	(0.908)	1107604	120.000	122.0	
73	2-Nitroaniline	65	6.950	6.949	(0.931)	346408	120.000	125.9	
76	Dimethylphthalate	163	7.229	7.229	(0.968)	1286101	120.000	123.0	
77	Acenaphthylene	152	7.281	7.281	(0.975)	1933504	120.000	122.3	
79	2,6-Dinitrotoluene	165	7.302	7.302	(0.978)	311050	120.000	127.7	
80	3-Nitroaniline	138	7.457	7.447	(0.999)	382849	120.000	125.9	
81	Acenaphthene	153	7.509	7.509	(1.006)	1207616	120.000	120.0	
82	2,4-Dinitrophenol	184	7.582	7.572	(1.015)	199007	120.000	124.7	
83	Dibenzofuran	168	7.706	7.706	(1.032)	1630240	120.000	122.0	(q)
84	4-Nitrophenol	109	7.675	7.675	(1.028)	161169	120.000	127.8	(Q)
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	409418	120.000	128.1	
91	Fluorene	166	8.131	8.131	(1.089)	1333949	120.000	120.6	
92	Diethylphthalate	149	8.110	8.100	(1.086)	1329206	120.000	124.2	
93	4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	558370	120.000	121.4	
94	4-Nitroaniline	138	8.224	8.214	(1.101)	378421	120.000	125.6	
97	4,6-Dinitro-2-methylphenol	198	8.286	8.276	(0.881)	236477	120.000	122.1	
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	1123239	141.000	143.7	
100	Azobenzene	77	8.359	8.348	(0.889)	1266722	120.000	124.9	
101	4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	318358	120.000	126.5	
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	335728	120.000	119.4	
110	Pentachlorophenol	266	9.240	9.240	(0.982)	215360	120.000	122.2	
114	Phenanthrene	178	9.437	9.437	(1.003)	1942962	120.000	119.6	
115	Anthracene	178	9.509	9.499	(1.011)	2014183	120.000	124.0	
118	Carbazole	167	9.768	9.768	(1.039)	1828217	120.000	123.3	
120	Di-n-Butylphthalate	149	10.463	10.463	(1.112)	2225048	120.000	124.7	
126	Fluoranthene	202	11.302	11.302	(1.202)	1829791	120.000	125.6	
127	Benzidine	184	11.582	11.571	(0.840)	1320429	120.000	127.8	
128	Pyrene	202	11.665	11.665	(0.846)	1963825	120.000	123.3	
134	3,3'-dimethylbenzidine	212	12.877	12.867	(0.934)	1214012	120.000	133.2	
136	Butylbenzylphthalate	149	12.991	12.991	(0.942)	997218	120.000	124.9	
138	Benzo(a)Anthracene	228	13.758	13.758	(0.998)	1694281	120.000	124.8	
139	Chrysene	228	13.831	13.831	(1.003)	1715841	120.000	123.6	
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.001)	653016	120.000	127.5	
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.023)	1368794	120.000	124.5	
142	Di-n-octylphthalate	149	15.167	15.167	(1.100)	2256614	120.000	128.4	
144	Benzo(b)fluoranthene	252	15.592	15.582	(0.964)	1475217	120.000	120.8	(Q)
145	Benzo(k)fluoranthene	252	15.623	15.623	(0.966)	1935987	120.000	123.5	(q)
147	Benzo(e)pyrene	252	16.007	16.007	(0.990)	1569049	120.000	123.2	
148	Benzo(a)pyrene	252	16.079	16.079	(0.994)	1720343	120.000	124.2	
151	Indeno(1,2,3-cd)pyrene	276	17.810	17.800	(1.101)	1517263	120.000	135.5	(M)
152	Dibenzo(a,h)anthracene	278	17.851	17.841	(1.104)	1534040	120.000	130.6	
153	Benzo(g,h,i)perylene	276	18.245	18.235	(1.128)	1706123	120.000	125.9	

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				3411204	120.000	122.3 (A)

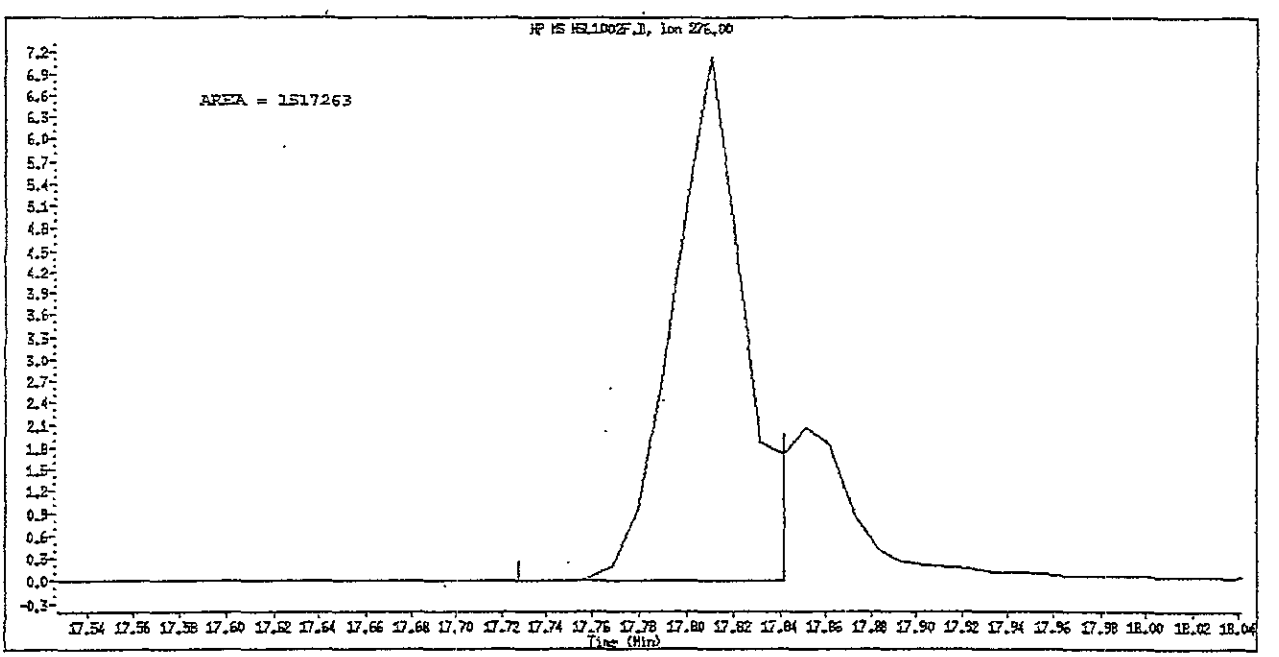
QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
M - Compound response manually integrated.
q - Qualifier signal exceeded ratio warning limit.

Data File Name: HSL1002F.D
Inj. Date and Time: 02-OCT-2010 14:35
Instrument ID: sv5.i
Client ID: 827DF.M
Compound Name: Indeno(1,2,3-cd)pyrene
CAS #: 193-39-5
Report Date: 10/03/2010



Original Integration



Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Poor Chromatography

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\100210.B\HSL1002F.D
Lab Smp Id: HSL 120 ug/ml CS-6 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 14:35
Operator : KT Inst ID: sv5.i
Smp Info : HSL_120 ug/ml CS-6;1;;6;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0312;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Meth Date : 02-Oct-2010 16:57 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 6 Calibration Sample, Level: 6
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)	
*	1 1,4-Dichlorobenzene-d4	152	3.955	3.955	(1.000)	137751	40.0000		(Q)
*	2 Naphthalene-d8	136	5.374	5.374	(1.000)	591665	40.0000		
*	3 Acenaphthene-d10	164	7.468	7.468	(1.000)	322596	40.0000		
*	4 Phenanthrene-d10	188	9.406	9.405	(1.000)	515607	40.0000		
*	5 Chrysene-d12	240	13.789	13.779	(1.000)	509570	40.0000		
*	6 Perylene-d12	264	16.173	16.162	(1.000)	539588	40.0000		
$	7 2-Fluorophenol	112	2.732	2.732	(0.691)	588028	120.000	115.7	
$	8 Phenol-d5	99	3.613	3.613	(0.914)	759824	120.000	117.5	
$	9 2-Chlorophenol-d4	132	3.758	3.758	(0.950)	652805	120.000	118.4	
$	10 1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	407247	120.000	118.4	
$	11 Nitrobenzene-d5	82	4.587	4.576	(0.853)	623501	120.000	118.8	
$	12 2-Fluorobiphenyl	172	6.680	6.680	(0.895)	1255441	120.000	121.5	
$	13 2,4,6-Tribromophenol	330	8.483	8.473	(1.136)	179055	120.000	140.6	
$	14 Terphenyl-d14	244	12.017	12.017	(0.871)	1251844	120.000	126.0	
	15 N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	388111	120.000	115.7	
	16 Pyridine	79	1.727	1.726	(0.437)	633334	120.000	113.3	
	23 Aniline	93	3.654	3.654	(0.924)	964533	120.000	119.0	(Q)
	24 Phenol	94	3.623	3.623	(0.916)	851671	120.000	123.6	(Q)
	26 Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	596323	120.000	114.2	
	27 2-Chlorophenol	128	3.768	3.768	(0.953)	653244	120.000	120.0	
	28 1,3-Dichlorobenzene	146	3.924	3.923	(0.992)	712032	120.000	118.4	
	29 1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	740915	120.000	121.9	
	30 Benzyl Alcohol	108	4.120	4.120	(1.042)	450249	120.000	120.4	
	31 1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	679448	120.000	117.9	
	32 2-Methylphenol	108	4.255	4.255	(1.076)	603987	120.000	118.8	
	33 2,2'-oxybis(1-Chloropropane)	45	4.297	4.297	(1.086)	941514	120.000	97.10	
	34 4-Methylphenol	108	4.421	4.421	(1.118)	644202	120.000	118.9	
	36 Hexachloroethane	117	4.504	4.504	(1.139)	245394	120.000	114.4	
	37 N-Nitrosodipropylamine	70	4.452	4.442	(1.126)	428242	120.000	112.9	
	42 Nitrobenzene	77	4.607	4.597	(0.857)	593736	120.000	113.8	
	44 Isophorone	82	4.867	4.856	(0.906)	1179801	120.000	119.3	
	45 2-Nitrophenol	139	4.960	4.960	(0.923)	367467	120.000	129.0	
	46 2,4-Dimethyphenol	107	5.012	5.012	(0.933)	638328	120.000	120.7	

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.954)	707504	120.000	120.2
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	500185	120.000	128.5
50	Benzoic Acid	122	5.146	5.115	(0.958)	395333	120.000	134.1
51	1,2,4-Trichlorobenzene	180	5.333	5.322	(0.992)	531764	120.000	126.0
52	Naphthalene	128	5.395	5.395	(1.004)	2020315	120.000	122.7
54	4-Chloroaniline	127	5.488	5.488	(1.021)	797064	120.000	123.0
57	Hexachlorobutadiene	225	5.613	5.613	(1.044)	255231	120.000	127.2
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	563840	120.000	125.9
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	1263302	120.000	125.7
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	312226	120.000	126.9
69	2,4,6-Trichlorophenol	196	6.587	6.576	(0.882)	331223	120.000	135.6
70	2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	343374	120.000	128.0
71	2-Chloronaphthalene	162	6.784	6.784	(0.908)	1107604	120.000	122.5
73	2-Nitroaniline	65	6.950	6.949	(0.931)	346408	120.000	114.4
76	Dimethylphthalate	163	7.229	7.229	(0.968)	1286101	120.000	123.1
77	Acenaphthylene	152	7.281	7.281	(0.975)	1933504	120.000	122.3
79	2,6-Dinitrotoluene	165	7.302	7.302	(0.978)	311050	120.000	133.0
80	3-Nitroaniline	138	7.457	7.447	(0.999)	382849	120.000	123.4
81	Acenaphthene	153	7.509	7.509	(1.006)	1207616	120.000	119.9
82	2,4-Dinitrophenol	184	7.582	7.571	(1.015)	199007	120.000	127.2
83	Dibenzofuran	168	7.706	7.706	(1.032)	1630240	120.000	122.5 (q)
84	4-Nitrophenol	109	7.675	7.675	(1.028)	161169	120.000	119.0 (Q)
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	409418	120.000	130.6
91	Fluorene	166	8.131	8.131	(1.089)	1333949	120.000	122.3
92	Diethylphthalate	149	8.110	8.100	(1.086)	1329206	120.000	121.7
93	4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	558370	120.000	124.2
94	4-Nitroaniline	138	8.224	8.214	(1.101)	378421	120.000	124.8
97	4,6-Dinitro-2-methylphenol	198	8.286	8.276	(0.881)	236477	120.000	120.3
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	1123239	141.000	139.8
100	Azobenzene	77	8.359	8.348	(0.889)	1266722	120.000	113.3
101	4-Bromophenyl-phenylether	248	8.794	8.794	(0.935)	318358	120.000	127.8
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	335728	120.000	124.8
110	Pentachlorophenol	266	9.240	9.240	(0.982)	215360	120.000	133.3
114	Phenanthrene	178	9.437	9.437	(1.003)	1942962	120.000	120.8
115	Anthracene	178	9.509	9.499	(1.011)	2014183	120.000	124.4
118	Carbazole	167	9.768	9.768	(1.039)	1828217	120.000	121.4
120	Di-n-Butylphthalate	149	10.463	10.463	(1.112)	2225048	120.000	122.2
126	Fluoranthene	202	11.302	11.302	(1.202)	1829791	120.000	126.4
127	Benzidine	184	11.582	11.571	(0.840)	1320429	120.000	126.2
128	Pyrene	202	11.665	11.665	(0.846)	1963825	120.000	123.2
134	3,3'-dimethylbenzidine	212	12.877	12.867	(0.934)	1214012	120.000	135.2
136	Butylbenzylphthalate	149	12.991	12.991	(0.942)	997218	120.000	122.5
138	Benzo(a)Anthracene	228	13.758	13.758	(0.998)	1694281	120.000	126.0
139	Chrysene	228	13.831	13.831	(1.003)	1715841	120.000	122.8
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.001)	653016	120.000	132.7
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.023)	1368794	120.000	122.1
142	Di-n-octylphthalate	149	15.167	15.167	(1.100)	2256614	120.000	125.9
144	Benzo(b)fluoranthene	252	15.592	15.582	(0.964)	1475217	120.000	115.3 (Q)
145	Benzo(k)fluoranthene	252	15.623	15.623	(0.966)	1935987	120.000	129.5 (q)
147	Benzo(e)pyrene	252	16.007	16.007	(0.990)	1569049	120.000	123.7
148	Benzo(a)pyrene	252	16.079	16.079	(0.994)	1720343	120.000	123.5
151	Indeno(1,2,3-cd)pyrene	276	17.810	17.800	(1.101)	1867193	120.000	151.5
152	Dibenzo(a,h)anthracene	278	17.851	17.841	(1.104)	1634040	120.000	129.4
153	Benzo(g,h,i)perylene	276	18.245	18.235	(1.128)	1706123	120.000	126.0

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				3411204	120.000	122.9 (A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002F.D
Lab Smp Id: HSL 120 ug/ml CS-6
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:
Method File: \\sv5\c\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0312;0;8270F.M

Test Mode:
Use Initial Calibration Level 4.

		AREA LIMIT			
COMPOUND	STANDARD	LOWER	UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	137751	12.34
2 Naphthalene-d8	530514	265257	1061028	591665	11.53
3 Acenaphthene-d10	282538	141269	565076	322596	14.18
4 Phenanthrene-d10	462722	231361	925444	515607	11.43
5 Chrysene-d12	435850	217925	871700	509570	16.91
6 Perylene-d12	422284	211142	844568	539588	27.78

		RT LIMIT			
COMPOUND	STANDARD	LOWER	UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.96	3.46	4.46	3.96	0.00
2 Naphthalene-d8	5.37	4.87	5.87	5.37	0.00
3 Acenaphthene-d10	7.47	6.97	7.97	7.47	0.00
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	0.00
5 Chrysene-d12	13.78	13.28	14.28	13.79	0.08
6 Perylene-d12	16.16	15.66	16.66	16.17	0.06

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV5\C\chem\sv5.i\100210.B\HSL1002F.D

Page 5

Date : 02-OCT-2010 14:35

Client ID: 8270F,M

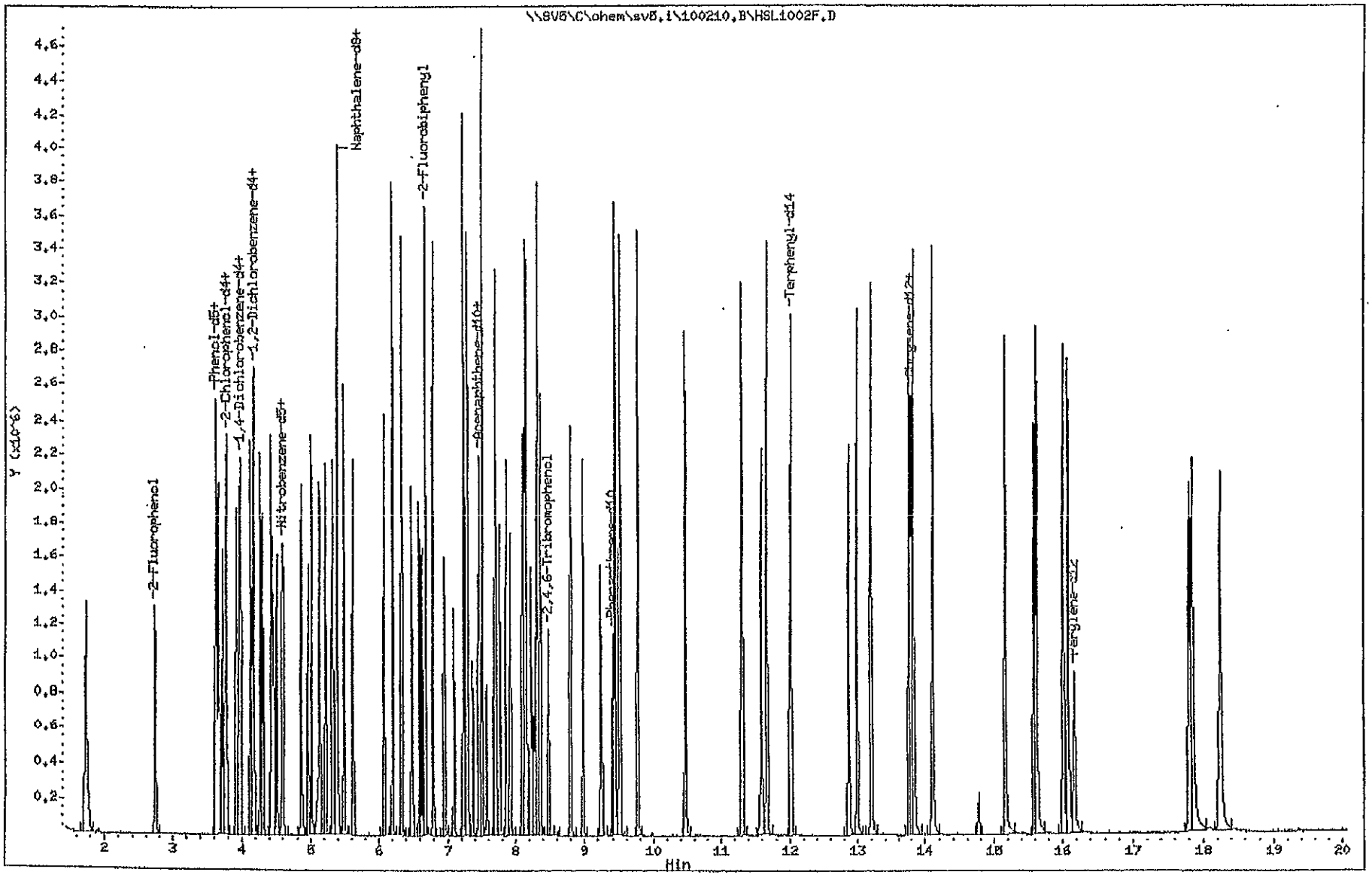
Instrument: sv5.i

Sample Info: HSL_120 ug/ml CS-5;1;;;5;;;;;4

Operator: KT

Column phase:

Column diameter: 2.00



TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002G.D
Lab Smp Id: HSL 160 ug/ml CS-7 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 15:00
Operator : KT Inst ID: sv5.i
Smp Info : HSL 160 ug/ml CS-7;1;;7;;;;4
Misc Info : 3;;0;1 8270STD.SUB;10MSSV0313;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:09 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 7 Calibration Sample, Level: 7
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SACP307UM

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
* 1	1,4-Dichlorobenzene-d4	152	3.954	3.955	(1.000)	141009	40.0000	(Q)
* 2	Naphthalene-d8	136	5.374	5.374	(1.000)	622461	40.0000	
* 3	Acenaphthene-d10	164	7.478	7.468	(1.000)	328259	40.0000	
* 4	Phenanthrene-d10	188	9.405	9.405	(1.000)	532284	40.0000	
* 5	Chrysene-d12	240	13.789	13.779	(1.000)	539557	40.0000	
* 6	Perylene-d12	264	16.172	16.162	(1.000)	560436	40.0000	
$ 7	2-Fluorophenol	112	2.732	2.732	(0.691)	810154	160.000	163.0 (A)
$ 8	Phenol-d5	99	3.623	3.613	(0.916)	1035724	160.000	165.7 (A)
$ 9	2-Chlorophenol-d4	132	3.757	3.758	(0.950)	890073	160.000	162.2 (A)
$ 10	1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	557810	160.000	160.6 (A)
$ 11	Nitrobenzene-d5	82	4.587	4.576	(0.853)	845795	160.000	160.4 (A)
$ 12	2-Fluorobiphenyl	172	6.680	6.680	(0.893)	1707074	160.000	161.4 (A)
$ 13	2,4,6-Tribromophenol	330	8.483	8.473	(1.134)	241468	160.000	169.3 (A)
$ 14	Terphenyl-d14	244	12.017	12.017	(0.871)	1728892	160.000	162.7 (A)
15	N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	529253	160.000	162.9 (Aq)
16	Pyridine	79	1.726	1.726	(0.437)	860850	160.000	158.4 (Q)
23	Aniline	93	3.654	3.654	(0.924)	1318620	160.000	165.8 (AQ)
24	Phenol	94	3.633	3.623	(0.919)	1166090	160.000	162.4 (AQ)
26	Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	813702	160.000	161.6 (A)
27	2-Chlorophenol	128	3.768	3.768	(0.953)	885754	160.000	160.7 (A)
28	1,3-Dichlorobenzene	146	3.923	3.923	(0.992)	972719	160.000	162.0 (A)
29	1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	1023408	160.000	163.0 (A)
30	Benzyl Alcohol	108	4.120	4.120	(1.042)	617653	160.000	166.7 (A)
31	1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	928919	160.000	160.9 (A)
32	2-Methylphenol	108	4.265	4.255	(1.079)	834149	160.000	165.4 (A)
33	2,2'-oxybis(1-Chloropropane)	45	4.296	4.297	(1.086)	1290345	160.000	161.0 (A)
34	4-Methylphenol	108	4.421	4.421	(1.118)	895481	160.000	167.2 (A)
36	Hexachloroethane	117	4.504	4.504	(1.139)	343605	160.000	160.7 (A)
37	N-Nitrosodipropylamine	70	4.452	4.442	(1.126)	590870	160.000	165.6 (A)
42	Nitrobenzene	77	4.607	4.597	(0.857)	844093	160.000	163.8 (A)
44	Isophorone	82	4.866	4.856	(0.906)	1628636	160.000	164.4 (A)
45	2-Nitrophenol	139	4.960	4.960	(0.923)	510613	160.000	167.0 (A)
46	2,4-Dimethyphenol	107	5.022	5.012	(0.934)	890994	160.000	164.0 (A)

10-3-10

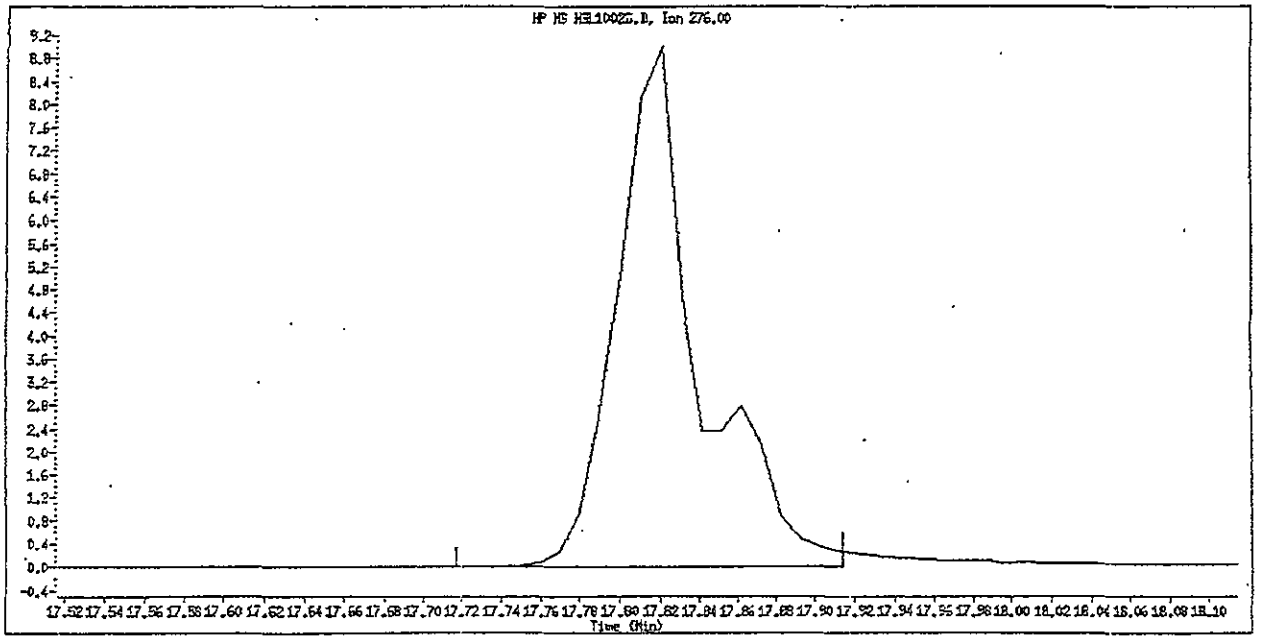
Compounds	QUANT SIG MASS	RT	REF RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47 Bis(2-chloroethoxy)methane	93	5.136	5.126	(0.956)	959710	160.000	158.5
49 2,4-Dichlorophenol	162	5.229	5.229	(0.973)	692405	160.000	164.7 (A)
50 Benzoic Acid	122	5.167	5.115	(0.961)	552251	160.000	183.6 (A)
51 1,2,4-Trichlorobenzene	180	5.333	5.322	(0.992)	724320	160.000	159.2
52 Naphthalene	128	5.395	5.395	(1.004)	2744968	160.000	159.7
54 4-Chloroaniline	127	5.488	5.488	(1.021)	1092223	160.000	162.1 (A)
57 Hexachlorobutadiene	225	5.612	5.613	(1.044)	360358	160.000	161.8 (A)
60 4-Chloro-3-Methylphenol	107	6.068	6.069	(1.129)	767831	160.000	163.6 (A)
63 2-Methylnaphthalene	142	6.203	6.203	(1.154)	1723402	160.000	159.6
66 Hexachlorocyclopentadiene	237	6.483	6.483	(0.867)	435738	160.000	177.9 (A)
69 2,4,6-Trichlorophenol	196	6.587	6.576	(0.881)	441685	160.000	168.6 (A)
70 2,4,5-Trichlorphenol	196	6.628	6.628	(0.886)	474468	160.000	168.2 (A)
71 2-Chloronaphthalene	162	6.783	6.784	(0.907)	1511253	160.000	163.6 (A)
73 2-Nitroaniline	65	6.960	6.949	(0.931)	476342	160.000	170.1 (A)
76 Dimethylphthalate	163	7.229	7.229	(0.967)	1710061	160.000	160.8 (A)
77 Acenaphthylene	152	7.291	7.281	(0.975)	2665048	160.000	165.6 (A)
79 2,6-Dinitrotoluene	165	7.302	7.302	(0.976)	408436	160.000	164.8 (A)
80 3-Nitroaniline	138	7.457	7.447	(0.997)	520002	160.000	168.1 (A)
81 Acenaphthene	153	7.509	7.509	(1.004)	1647377	160.000	160.9 (A)
82 2,4-Dinitrophenol	184	7.581	7.572	(1.014)	265655	160.000	157.7
83 Dibenzofuran	168	7.706	7.706	(1.030)	2246304	160.000	165.3 (A)
84 4-Nitrophenol	109	7.685	7.675	(1.028)	228516	160.000	178.1 (Aq)
86 2,4-Dinitrotoluene	165	7.778	7.768	(1.040)	566055	160.000	174.0 (A)
91 Fluorene	166	8.141	8.131	(1.089)	1846653	160.000	164.1 (A)
92 Diethylphthalate	149	8.110	8.100	(1.085)	1813127	160.000	166.5 (A)
93 4-Chlorophenyl-phenylether	204	8.151	8.152	(1.090)	757562	160.000	161.9 (A)
94 4-Nitroaniline	138	8.224	8.214	(1.100)	531151	160.000	173.2 (A)
97 4,6-Dinitro-2-methylphenol	198	8.286	8.276	(0.881)	324244	160.000	160.7 (A)
98 N-Nitrosodiphenylamine	169	8.328	8.317	(0.885)	1542041	187.000	191.1 (A)
100 Azobenzene	77	8.359	8.348	(0.889)	1646477	160.000	157.3
101 4-Bromophenyl-phenylether	248	8.804	8.794	(0.936)	421894	160.000	162.4 (A)
108 Hexachlorobenzene	284	8.980	8.981	(0.955)	465305	160.000	160.3 (A)
110 Pentachlorophenol	266	9.250	9.240	(0.983)	293184	160.000	159.9
114 Phenanthrene	178	9.447	9.437	(1.004)	2695719	160.000	160.7 (A)
115 Anthracene	178	9.509	9.499	(1.011)	2703105	160.000	161.3 (A)
118 Carbazole	167	9.768	9.768	(1.039)	2479487	160.000	161.9 (A)
120 Di-n-Butylphthalate	149	10.473	10.463	(1.113)	3164666	160.000	171.8 (A)
126 Fluoranthene	202	11.312	11.302	(1.203)	2500453	160.000	166.3 (A)
127 Benzidine	184	11.582	11.571	(0.840)	1864289	160.000	170.5 (A)
128 Pyrene	202	11.664	11.665	(0.846)	2714930	160.000	161.0 (A)
134 3,3'-dimethylbenzidine	212	12.877	12.867	(0.934)	1724989	160.000	178.7 (A)
136 Butylbenzylphthalate	149	12.991	12.991	(0.942)	1401117	160.000	165.8 (A)
138 Benzo(a)Anthracene	228	13.768	13.758	(0.998)	2393908	160.000	166.6 (A)
139 Chrysene	228	13.841	13.831	(1.004)	2422526	160.000	164.5 (A)
140 3,3'-Dichlorobenzidine	252	13.810	13.799	(1.002)	915413	160.000	168.9 (A)
141 bis(2-ethylhexyl) Phthalate	149	14.110	14.110	(1.023)	1906885	160.000	163.8 (A)
142 Di-n-octylphthalate	149	15.167	15.167	(1.100)	2253965	160.000	174.8 (A)
144 Benzo(b)fluoranthene	252	15.592	15.582	(0.964)	2299398	160.000	181.2 (AQ)
145 Benzo(k)fluoranthene	252	15.634	15.623	(0.967)	2475925	160.000	152.0 (q)
147 Benzo(e)pyrene	252	16.017	16.007	(0.990)	2178628	160.000	164.7 (A)
148 Benzo(a)pyrene	252	16.089	16.079	(0.995)	2387962	160.000	166.0 (A)
151 Indeno(1,2,3-cd)pyrene	276	17.820	17.800	(1.102)	2196805	160.000	188.8 (AM)
152 Dibenzo(a,h)anthracene	278	17.862	17.841	(1.104)	2250528	160.000	173.2 (A)
153 Benzo(g,h,i)perylene	276	18.255	18.235	(1.129)	2332007	160.000	165.7 (A)

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				4775333	160.000	164.8 (A)

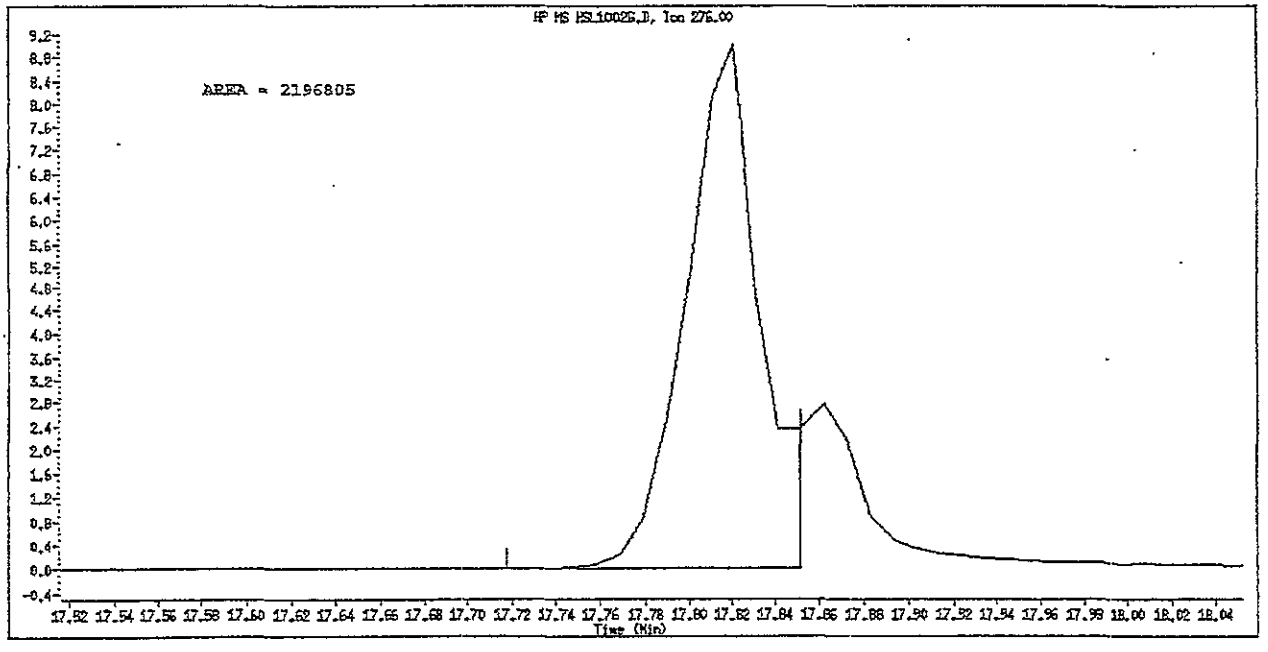
QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
M - Compound response manually integrated.
q - Qualifier signal exceeded ratio warning limit.

Data File Name: ESL1002G.D
Inj. Date and Time: 02-OCT-2010 15:00
Instrument ID: svS.i
Client ID: 8270F.M
Compound Name: Indeno(1,2,3-cd)pyrene
CAS #: 193-39-5
Report Date: 10/03/2010



Original Integration



Manual Integration

Manually Integrated By: truongk
Manual Integration Reason: Poor Chromatography

TestAmerica West Sacramento

Method 8270C

Data file : \\SV5\C\chem\sv5.i\100210.B\HSL1002G.D
Lab Smp Id: HSL 160 ug/ml CS-7 Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 15:00
Operator : KT Inst ID: sv5.i
Smp Info : HSL 160 ug/ml CS-7;1;;7;;;;4
Misc Info : 3;;0;1 8270STD.SUB;10MSSV0313;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Meth Date : 02-Oct-2010 16:57 onishim Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 7 Calibration Sample, Level: 7
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SV5

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
* 1	1,4-Dichlorobenzene-d4	152	3.954	3.955	(1.000)	141009	40.0000	(Q)
* 2	Naphthalene-d8	136	5.374	5.374	(1.000)	622461	40.0000	
* 3	Acenaphthene-d10	164	7.478	7.468	(1.000)	328259	40.0000	
* 4	Phenanthrene-d10	188	9.405	9.405	(1.000)	532284	40.0000	
* 5	Chrysene-d12	240	13.789	13.779	(1.000)	539557	40.0000	
* 6	Perylene-d12	264	16.172	16.162	(1.000)	560436	40.0000	
$ 7	2-Fluorophenol	112	2.732	2.732	(0.691)	810154	160.000	155.7
$ 8	Phenol-d5	99	3.623	3.613	(0.916)	1035724	160.000	156.5
$ 9	2-Chlorophenol-d4	132	3.757	3.758	(0.950)	890073	160.000	157.7
$ 10	1,2-Dichlorobenzene-d4	152	4.162	4.162	(1.052)	557810	160.000	158.4
$ 11	Nitrobenzene-d5	82	4.587	4.576	(0.853)	845796	160.000	153.2
$ 12	2-Fluorobiphenyl	172	6.660	6.680	(0.893)	1707074	160.000	162.4 (A)
$ 13	2,4,6-Tribromophenol	330	8.483	8.473	(1.134)	241468	160.000	186.3 (A)
$ 14	Terphenyl-d14	244	12.017	12.017	(0.871)	1728892	160.000	164.3 (A)
15	N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	529253	160.000	154.1
16	Pyridine	79	1.726	1.726	(0.437)	860850	160.000	150.4
23	Aniline	93	3.654	3.654	(0.924)	1318620	160.000	158.9 (Q)
24	Phenol	94	3.633	3.623	(0.919)	1166090	160.000	165.7 (AQ)
26	Bis(2-chloroethyl)ether	93	3.716	3.716	(0.940)	813702	160.000	152.2
27	2-Chlorophenol	128	3.768	3.768	(0.953)	885754	160.000	159.0
28	1,3-Dichlorobenzene	146	3.923	3.923	(0.992)	972719	160.000	158.0
29	1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	1023408	160.000	164.5 (A)
30	Benzyl Alcohol	108	4.120	4.120	(1.042)	617653	160.000	161.4 (A)
31	1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	928919	160.000	157.5
32	2-Methylphenol	108	4.265	4.255	(1.079)	834149	160.000	160.3 (A)
33	2,2'-oxybis(1-Chloropropane)	45	4.296	4.297	(1.086)	1290345	160.000	130.0
34	4-Methylphenol	108	4.421	4.421	(1.118)	895481	160.000	161.5 (A)
36	Hexachloroethane	117	4.504	4.504	(1.139)	343605	160.000	156.5
37	N-Nitrosodipropylamine	70	4.452	4.442	(1.126)	590870	160.000	152.2
42	Nitrobenzene	77	4.607	4.597	(0.857)	844093	160.000	153.8
44	Isophorone	82	4.866	4.856	(0.906)	1628636	160.000	156.6
45	2-Nitrophenol	139	4.960	4.960	(0.923)	510613	160.000	170.5 (A)
46	2,4-Dimethylphenol	107	5.022	5.012	(0.934)	890994	160.000	160.2 (A)

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	AMOUNTS ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.136	5.126	(0.956)	959710	160.000	155.0
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	692405	160.000	169.1(A)
50	Benzoic Acid	122	5.167	5.115	(0.961)	552251	160.000	178.1(A)
51	1,2,4-Trichlorobenzene	180	5.333	5.322	(0.992)	724320	160.000	163.2(A)
52	Naphthalene	128	5.395	5.395	(1.004)	2744968	160.000	158.4
54	4-Chloroaniline	127	5.488	5.488	(1.021)	1092223	160.000	160.2(A)
57	Hexachlorobutadiene	225	5.612	5.613	(1.044)	360358	160.000	170.6(A)
60	4-Chloro-3-Methylphenol	107	6.068	6.069	(1.129)	767831	160.000	163.0(A)
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	1723402	160.000	163.0(A)
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.867)	435738	160.000	174.0(A)
69	2,4,6-Trichlorophenol	196	6.587	6.576	(0.881)	441685	160.000	177.7(A)
70	2,4,5-Trichlorphenol	196	6.628	6.628	(0.886)	474468	160.000	173.8(A)
71	2-Chloronaphthalene	162	6.783	6.784	(0.907)	1511253	160.000	164.2(A)
73	2-Nitroaniline	65	6.960	6.949	(0.931)	476342	160.000	154.5
76	Dimethylphthalate	163	7.229	7.229	(0.967)	1710061	160.000	160.9(A)
77	Acenaphthylene	152	7.291	7.281	(0.975)	2665048	160.000	165.6(A)
79	2,6-Dinitrotoluene	165	7.302	7.302	(0.976)	408436	160.000	171.6(A)
80	3-Nitroaniline	138	7.457	7.447	(0.997)	520002	160.000	164.8(A)
81	Acenaphthene	153	7.509	7.509	(1.004)	1647377	160.000	160.7(A)
82	2,4-Dinitrophenol	184	7.581	7.571	(1.014)	265655	160.000	158.9
83	Dibenzofuran	168	7.706	7.706	(1.030)	2246304	160.000	165.8(A)
84	4-Nitrophenol	109	7.685	7.675	(1.028)	228516	160.000	165.8(Aq)
86	2,4-Dinitrotoluene	165	7.778	7.768	(1.040)	566055	160.000	177.5(A)
91	Fluorene	166	8.141	8.131	(1.089)	1846653	160.000	166.4(A)
92	Diethylphthalate	149	8.110	8.100	(1.085)	1813127	160.000	163.2(A)
93	4-Chlorophenyl-phenylether	204	8.151	8.152	(1.090)	757562	160.000	165.6(A)
94	4-Nitroaniline	138	8.224	8.214	(1.100)	531151	160.000	172.2(A)
97	4,6-Dinitro-2-methylphenol	198	8.286	8.276	(0.881)	324264	160.000	158.0
98	N-Nitrosodiphenylamine	169	8.328	8.317	(0.885)	1542041	187.000	185.9(A)
100	Azobenzene	77	8.359	8.348	(0.889)	1646477	160.000	142.7
101	4-Bromophenyl-phenylether	248	8.804	8.794	(0.936)	421894	160.000	164.0(A)
108	Hexachlorobenzene	284	8.980	8.981	(0.955)	465305	160.000	167.5(A)
110	Pentachlorophenol	266	9.250	9.240	(0.983)	293184	160.000	175.8(A)
114	Phenanthrene	178	9.447	9.437	(1.004)	2695719	160.000	162.4(A)
115	Anthracene	178	9.509	9.499	(1.011)	2703105	160.000	161.8(A)
118	Carbazole	167	9.768	9.758	(1.039)	2479487	160.000	159.5
120	Di-n-Butylphthalate	149	10.473	10.463	(1.113)	3164666	160.000	168.4(A)
126	Fluoranthene	202	11.312	11.302	(1.203)	2500453	160.000	167.3(A)
127	Benzidine	184	11.582	11.571	(0.840)	1864289	160.000	168.3(A)
128	Pyrene	202	11.664	11.665	(0.846)	2714930	160.000	160.9(A)
134	3,3'-dimethylbenzidine	212	12.877	12.867	(0.934)	1724989	160.000	181.4(A)
136	Butylbenzylphthalate	149	12.991	12.991	(0.942)	1401117	160.000	162.5(A)
138	Benzo(a)Anthracene	228	13.768	13.758	(0.998)	2393908	160.000	168.2(A)
139	Chrysene	228	13.841	13.831	(1.004)	2422526	160.000	163.8(A)
140	3,3'-Dichlorobenzidine	252	13.810	13.799	(1.002)	915413	160.000	175.7(A)
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.023)	1906885	160.000	160.7(A)
142	Di-n-octylphthalate	149	15.167	15.167	(1.100)	3253965	160.000	171.5(A)
144	Benzo(b)fluoranthene	252	15.592	15.582	(0.964)	2299398	160.000	173.0(AQ)
145	Benzo(k)fluoranthene	252	15.634	15.623	(0.967)	2475935	160.000	159.4(q)
147	Benzo(e)pyrene	252	16.017	16.007	(0.990)	2178528	160.000	165.4(A)
148	Benzo(a)pyrene	252	16.089	16.079	(0.995)	2387962	160.000	165.1(A)
151	Indeno(1,2,3-cd)pyrene	276	17.820	17.800	(1.102)	2617878	160.000	204.6(A)
152	Dibenzo(a,h)anthracene	278	17.862	17.841	(1.104)	2250528	160.000	171.6(A)
153	Benzo(g,h,i)perylene	276	18.255	18.235	(1.129)	2332087	160.000	165.9(A)

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				4775333	160.000	165.7(A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.
Q - Qualifier signal failed the ratio test.
q - Qualifier signal exceeded ratio warning limit.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002G.D
Lab Smp Id: HSL_160 ug/ml CS-7
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\sv5\c\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0313;0;8270F.M

Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	141009	14.99
2 Naphthalene-d8	530514	265257	1061028	622461	17.33
3 Acenaphthene-d10	282538	141269	565076	328259	16.18
4 Phenanthrene-d10	462722	231361	925444	532284	15.03
5 Chrysene-d12	435850	217925	871700	539557	23.79
6 Perylene-d12	422284	211142	844568	560436	32.72

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.96	3.46	4.46	3.95	-0.00
2 Naphthalene-d8	5.37	4.87	5.87	5.37	-0.00
3 Acenaphthene-d10	7.47	6.97	7.97	7.48	0.14
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	-0.00
5 Chrysene-d12	13.78	13.28	14.28	13.79	0.07
6 Perylene-d12	16.16	15.66	16.66	16.17	0.06

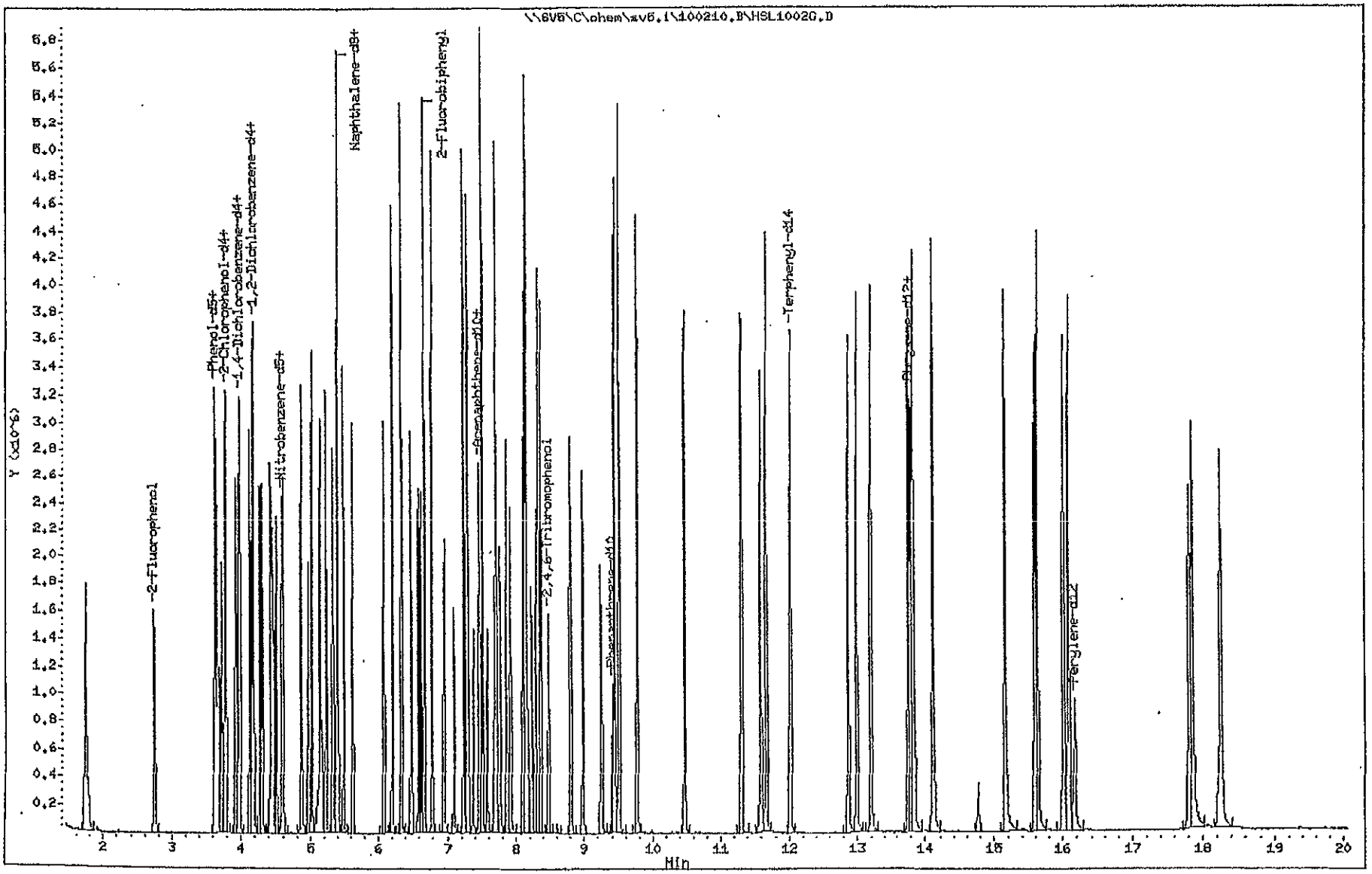
AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV5\C\ohem\sv5.i\100210.B\HSL1002G.D
Date : 02-OCT-2010 15:00
Client ID: 8270F.M
Sample Info: HSL_160 ug/ml CS-7;1;;7;;;;4

Instrument: sv5.i

Operator: KT
Column diameter: 2.00

Column phase:



TestAmerica West Sacramento

CONTINUING CALIBRATION COMPOUNDS

Instrument ID: sv5.i Injection Date: 02-OCT-2010 16:11
Lab File ID: HSL1002H.D Init. Cal. Date(s): 17-AUG-2010 02-OCT-2010
Analysis Type: Init. Cal. Times: 17:32 15:00
Lab Sample ID: HSL_050 ug/ml ICV Quant Type: ISTD
Method: \\sv5\c\chem\sv5.i\100210.B\8270f.m

COMPOUND	RRF / AMOUNT	RF50	CCAL RRF50	MIN RRF	%D / %DRIFT	MAX %D / %DRIFT	CURVE TYPE
$ 7 2-Fluorophenol	1.40992	1.41047	1.41047	0.010	0.03876	50.00000	Averaged
$ 8 Phenol-d5	1.77296	1.74907	1.74907	0.010	-1.34746	50.00000	Averaged
$ 9 2-Chlorophenol-d4	1.55698	1.55303	1.55303	0.010	-0.25385	50.00000	Averaged
$ 10 1,2-Dichlorobenzene-d4	0.98513	0.98502	0.98502	0.010	-0.01093	50.00000	Averaged
$ 11 Nitrobenzene-d5	0.33879	0.32706	0.32706	0.010	-3.46219	50.00000	Averaged
$ 12 2-Fluorobiphenyl	1.28852	1.25302	1.25302	0.010	-2.75502	50.00000	Averaged
$ 13 2,4,6-Tribromophenol	0.17381	0.17822	0.17822	0.010	2.53174	50.00000	Averaged
$ 14 Terphenyl-d14	0.78789	0.74054	0.74054	0.010	-6.00962	50.00000	Averaged
15 N-Nitrosodimethylamine	0.92154	0.91645	0.91645	0.010	-0.55265	50.00000	Averaged
16 Pyridine	1.54111	1.49084	1.49084	0.010	-3.26208	50.00000	Averaged
23 Aniline	2.25673	1.90520	1.90520	0.010	-15.57680	50.00000	Averaged
24 Phenol	2.03729	2.01343	2.01343	0.010	-1.17106	20.00000	Averaged
26 Bis(2-chloroethyl)ether	1.42859	1.41690	1.41690	0.010	-0.81844	50.00000	Averaged
27 2-Chlorophenol	1.56381	1.57626	1.57626	0.010	0.79611	50.00000	Averaged
28 1,3-Dichlorobenzene	1.70337	1.74104	1.74104	0.010	2.21094	50.00000	Averaged
29 1,4-Dichlorobenzene	1.78118	1.77637	1.77637	0.010	-0.26978	20.00000	Averaged
30 Benzyl Alcohol	1.05101	1.07153	1.07153	0.010	1.95228	50.00000	Averaged
31 1,2-Dichlorobenzene	1.63746	1.64144	1.64144	0.010	0.24267	50.00000	Averaged
32 2-Methylphenol	1.43012	1.41817	1.41817	0.010	-0.83592	50.00000	Averaged
33 2,2'-oxybis(1-Chloropropane	2.27365	2.14153	2.14153	0.010	-5.81096	50.00000	Averaged
34 4-Methylphenol	1.51904	1.42403	1.42403	0.010	-6.25452	50.00000	Averaged
36 Hexachloroethane	0.60636	0.62081	0.62081	0.010	2.38271	50.00000	Averaged
37 N-Nitrosodinpropylamine	1.01180	0.99863	0.99863	0.050	-1.30217	50.00000	Averaged
42 Nitrobenzene	0.33116	0.32452	0.32452	0.010	-2.00546	50.00000	Averaged
44 Isophorone	0.63679	0.62370	0.62370	0.010	-2.05513	50.00000	Averaged
45 2-Nitrophenol	0.19648	0.20090	0.20090	0.010	2.25050	20.00000	Averaged
46 2,4-Dimethyphenol	0.34911	0.33078	0.33078	0.010	-5.25153	50.00000	Averaged
47 Bis(2-chloroethoxy)methane	0.38908	0.37434	0.37434	0.010	-3.78942	50.00000	Averaged
49 2,4-Dichlorophenol	0.27010	0.26945	0.26945	0.010	-0.23923	20.00000	Averaged
50 Benzoic Acid	0.19324	0.20284	0.20284	0.010	4.96710	50.00000	Averaged
51 1,2,4-Trichlorobenzene	0.29246	0.28203	0.28203	0.010	-3.56320	50.00000	Averaged
52 Naphthalene	1.10443	1.07116	1.07116	0.010	-3.01217	50.00000	Averaged
54 4-Chloroaniline	0.43288	0.40664	0.40664	0.010	-6.06033	50.00000	Averaged
57 Hexachlorobutadiene	0.14313	0.14742	0.14742	0.010	2.99976	20.00000	Averaged
60 4-Chloro-3-Methylphenol	0.30164	0.29442	0.29442	0.010	-2.39317	20.00000	Averaged
63 2-Methylnaphthalene	0.69378	0.71003	0.71003	0.010	2.34296	50.00000	Averaged
66 Hexachlorocyclopentadiene	0.29846	0.32228	0.32228	0.050	7.98199	50.00000	Averaged
69 2,4,6-Trichlorophenol	0.31913	0.32462	0.32462	0.010	1.71977	20.00000	Averaged
70 2,4,5-Trichlorphenol	0.34380	0.34503	0.34503	0.010	0.35814	50.00000	Averaged
71 2-Chloronaphthalene	1.12571	1.09768	1.09768	0.010	-2.48963	50.00000	Averaged
73 2-Nitroaniline	0.34119	0.32550	0.32550	0.010	-4.59608	50.00000	Averaged
76 Dimethylphthalate	1.29606	1.28355	1.28355	0.010	-0.96554	50.00000	Averaged

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TestAmerica West Sacramento

CONTINUING CALIBRATION COMPOUNDS

Instrument ID: sv5.i Injection Date: 02-OCT-2010 16:11
Lab File ID: HSL1002H.D Init. Cal. Date(s): 17-AUG-2010 02-OCT-2010
Analysis Type: Init. Cal. Times: 17:32 15:00
Lab Sample ID: HSL_050 ug/ml ICV Quant Type: ISTD
Method: \\sv5\c\chem\sv5.i\100210.B\8270f.m

COMPOUND	RRF / AMOUNT	RF50	CCAL RRF50	MIN RRF	%D / %DRIFT	MAX %D / %DRIFT	CURVE TYPE
77 Acenaphthylene	1.96037	1.90194	1.90194	0.010	-2.98044	50.00000	Averaged
79 2,6-Dinitrotoluene	0.30197	0.30334	0.30334	0.010	0.45457	50.00000	Averaged
80 3-Nitroaniline	0.37691	0.37836	0.37836	0.010	0.38563	50.00000	Averaged
81 Acenaphthene	1.24787	1.19989	1.19989	0.010	-3.84461	20.00000	Averaged
82 2,4-Dinitrophenol	50.00000	48.07731	0.16950	0.050	-3.84537	0.000e+000	Quadratic
83 Dibenzofuran	1.65612	1.64309	1.64309	0.010	-0.78683	50.00000	Averaged
84 4-Nitrophenol	0.15634	0.16205	0.16205	0.050	3.65012	50.00000	Averaged
86 2,4-Dinitrotoluene	0.39633	0.40639	0.40639	0.010	2.53669	50.00000	Averaged
91 Fluorene	1.37139	1.36209	1.36209	0.010	-0.67828	50.00000	Averaged
92 Diethylphthalate	1.32699	1.28445	1.28445	0.010	-3.20581	50.00000	Averaged
93 4-Chlorophenyl-phenylether	0.57019	0.56986	0.56986	0.010	-0.05862	50.00000	Averaged
94 4-Nitroaniline	0.37361	0.40608	0.40608	0.010	8.68956	50.00000	Averaged
97 4,6-Dinitro-2-methylphenol	50.00000	48.62001	0.13800	0.010	-2.75999	0.000e+000	Linear
98 N-Nitrosodiphenylamine	0.60628	0.49086	0.49086	0.010	-19.03836	20.00000	Averaged
100 Azobenzene	0.78650	0.77322	0.77322	0.010	-1.70096	50.00000	Averaged
101 4-Bromophenyl-phenylether	0.19527	0.19536	0.19536	0.010	0.04546	50.00000	Averaged
108 Hexachlorobenzene	0.21807	0.22026	0.22026	0.010	1.00466	50.00000	Averaged
110 Pentachlorophenol	50.00000	50.72441	0.13218	0.010	1.44881	0.000e+000	Linear
114 Phenanthrene	1.26074	1.20864	1.20864	0.010	-4.13307	50.00000	Averaged
115 Anthracene	1.25955	1.22825	1.22825	0.010	-2.48429	50.00000	Averaged
118 Carbazole	1.15061	1.15083	1.15083	0.010	0.01942	50.00000	Averaged
120 Di-n-Butylphthalate	1.38442	1.39149	1.39149	0.010	0.51078	50.00000	Averaged
126 Fluoranthene	1.12969	1.19302	1.19302	0.010	5.60642	20.00000	Averaged
127 Benzidine	0.81067	0.30175	0.30175	0.010	-62.77740	50.00000	Averaged
128 Pyrene	1.25025	1.13023	1.13023	0.010	-9.59978	50.00000	Averaged
134 3,3'-dimethylbenzidine	0.71564	0.26880	0.26880	0.010	-62.43954	50.00000	Averaged
136 Butylbenzylphthalate	0.62663	0.58836	0.58836	0.010	-6.10747	50.00000	Averaged
138 Benzo(a)Anthracene	1.06548	0.99285	0.99285	0.010	-6.81596	50.00000	Averaged
139 Chrysene	1.08994	1.04703	1.04703	0.010	-3.93621	50.00000	Averaged
140 3,3'-Dichlorobenzidine	0.40189	0.37691	0.37691	0.010	-6.21534	50.00000	Averaged
141 bis(2-ethylhexyl)Phthalate	0.86316	0.80149	0.80149	0.010	-7.14468	50.00000	Averaged
142 Di-n-octylphthalate	1.37975	1.27404	1.27404	0.010	-7.66156	20.00000	Averaged
144 Benzo(b)fluoranthene	0.90549	0.90498	0.90498	0.010	-0.05663	50.00000	Averaged
145 Benzo(k)fluoranthene	1.16236	1.22175	1.22175	0.010	5.10982	50.00000	Averaged
147 Benzo(e)pyrene	0.94425	0.98421	0.98421	0.010	4.23177	50.00000	Averaged
148 Benzo(a)pyrene	1.02655	0.95393	0.95393	0.010	-7.07365	20.00000	Averaged
151 Indeno(1,2,3-cd)pyrene	0.83029	0.81846	0.81846	0.010	-1.42489	50.00000	Averaged
152 Dibenzo(a,h)anthracene	0.92758	0.99090	0.99090	0.010	6.82730	50.00000	Averaged
153 Benzo(g,h,i)perylene	1.00427	1.08674	1.08674	0.010	8.21177	50.00000	Averaged
M 162 benzo b,k Fluoranthene Tota	2.06785	2.12673	2.12673	0.010	2.84748	50.00000	Averaged

See NT

See RD

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TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002H.D
Lab Smp Id: HSL_050 ug/ml ICV Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 16:11
Operator : KT Inst ID: sv5.i
Smp Info : HSL_050 ug/ml ICV;2;;4;;;;4
Misc Info : 3;;0;1_8270STD.SUB;10MSSV0314;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:20 sv5.i Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 8 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: 1_8270STD.SUB
Target Version: 4.14
Processing Host: SACP307DM

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
*	1 1,4-Dichlorobenzene-d4	152	3.954	3.954	(1.000)	98364	40.0000	
*	2 Naphthalene-d8	136	5.374	5.374	(1.000)	431655	40.0000	
*	3 Acenaphthene-d10	164	7.468	7.468	(1.000)	236662	40.0000	
*	4 Phenanthrene-d10	188	9.405	9.405	(1.000)	380734	40.0000	
*	5 Chrysene-d12	240	13.789	13.789	(1.000)	421719	40.0000	
*	6 Perylene-d12	264	16.173	16.173	(1.000)	419419	40.0000	
$	7 2-Fluorophenol	112	2.732	2.732	(0.691)	173424	50.0000	50.02
$	8 Phenol-d5	99	3.613	3.613	(0.914)	215057	50.0000	49.33
$	9 2-Chlorophenol-d4	132	3.747	3.747	(0.948)	190953	50.0000	49.87
$	10 1,2-Dichlorobenzene-d4	152	4.151	4.151	(1.050)	121113	50.0000	49.99
$	11 Nitrobenzene-d5	82	4.576	4.576	(0.852)	176474	50.0000	48.27
$	12 2-Fluorobiphenyl	172	6.680	6.680	(0.895)	370679	50.0000	48.62
$	13 2,4,6-Tribromophenol	330	8.483	8.483	(1.136)	52721	50.0000	51.26
$	14 Terphenyl-d14	244	12.017	12.017	(0.871)	390377	50.0000	47.00
	15 N-Nitrosodimethylamine	74	1.706	1.706	(0.431)	112682	50.0000	49.72(Q)
	16 Pyridine	79	1.726	1.726	(0.437)	183306	50.0000	48.37
	23 Aniline	93	3.654	3.654	(0.924)	234254	50.0000	42.21
	24 Phenol	94	3.623	3.623	(0.916)	247561	50.0000	49.41(Q)
	26 Bis(2-chloroethyl) ether	93	3.716	3.716	(0.940)	174215	50.0000	49.59
	27 2-Chlorophenol	128	3.768	3.768	(0.953)	193809	50.0000	50.40
	28 1,3-Dichlorobenzene	146	3.913	3.913	(0.990)	214069	50.0000	51.10
	29 1,4-Dichlorobenzene	146	3.975	3.975	(1.005)	218414	50.0000	49.86
	30 Benzyl Alcohol	108	4.120	4.120	(1.042)	131750	50.0000	50.98
	31 1,2-Dichlorobenzene	146	4.172	4.172	(1.055)	201823	50.0000	50.12
	32 2-Methylphenol	108	4.255	4.255	(1.076)	174371	50.0000	49.58
	33 2,2'-oxybis(1-Chloropropane)	45	4.296	4.296	(1.086)	263312	50.0000	47.09
	34 4-Methylphenol	108	4.410	4.410	(1.115)	175092	50.0000	46.87
	36 Hexachloroethane	117	4.504	4.504	(1.139)	76332	50.0000	51.19
	37 N-Nitrosodipropylamine	70	4.442	4.442	(1.123)	122786	50.0000	49.35
	42 Nitrobenzene	77	4.597	4.597	(0.855)	175102	50.0000	49.00
	44 Isophorone	82	4.856	4.856	(0.904)	336530	50.0000	48.97
	45 2-Nitrophenol	139	4.960	4.960	(0.923)	108399	50.0000	51.12
	46 2,4-Dimethyphenol	107	5.012	5.012	(0.933)	178479	50.0000	47.37

	Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
47	Bis(2-chloroethoxy)methane	93	5.126	5.126	(0.954)	201982	50.0000	48.10
49	2,4-Dichlorophenol	162	5.229	5.229	(0.973)	145389	50.0000	49.88
50	Benzoic Acid	122	5.115	5.115	(0.952)	109446	50.0000	52.48
51	1,2,4-Trichlorobenzene	180	5.322	5.322	(0.990)	152177	50.0000	48.22
52	Naphthalene	128	5.395	5.395	(1.004)	577964	50.0000	48.49
54	4-Chloroaniline	127	5.488	5.488	(1.021)	219411	50.0000	46.97
57	Hexachlorobutadiene	225	5.613	5.613	(1.044)	79543	50.0000	51.50
60	4-Chloro-3-Methylphenol	107	6.069	6.069	(1.129)	158858	50.0000	48.80
63	2-Methylnaphthalene	142	6.203	6.203	(1.154)	383110	50.0000	51.17
66	Hexachlorocyclopentadiene	237	6.483	6.483	(0.868)	95339	50.0000	53.99
69	2,4,6-Trichlorophenol	196	6.587	6.587	(0.882)	96032	50.0000	50.86
70	2,4,5-Trichlorphenol	196	6.628	6.628	(0.888)	102070	50.0000	50.18
71	2-Chloronaphthalene	162	6.784	6.784	(0.908)	324725	50.0000	48.76
73	2-Nitroaniline	65	6.949	6.949	(0.931)	96293	50.0000	47.70
76	Dimethylphthalate	163	7.229	7.229	(0.968)	379709	50.0000	49.52
77	Acenaphthylene	152	7.281	7.281	(0.975)	562646	50.0000	48.51
79	2,6-Dinitrotoluene	165	7.302	7.302	(0.978)	89736	50.0000	50.23
80	3-Nitroaniline	138	7.457	7.457	(0.999)	111929	50.0000	50.19
81	Acenaphthene	153	7.509	7.509	(1.006)	354961	50.0000	48.08
82	2,4-Dinitrophenol	184	7.582	7.582	(1.015)	50142	50.0000	48.08
83	Dibenzofuran	168	7.706	7.706	(1.032)	486071	50.0000	49.61
84	4-Nitrophenol	109	7.675	7.675	(1.028)	47938	50.0000	51.82(Q)
86	2,4-Dinitrotoluene	165	7.768	7.768	(1.040)	120220	50.0000	51.27
91	Fluorene	166	8.131	8.131	(1.089)	402944	50.0000	49.66
92	Diethylphthalate	149	8.100	8.100	(1.085)	379976	50.0000	48.40
93	4-Chlorophenyl-phenylether	204	8.152	8.152	(1.092)	168579	50.0000	49.97
94	4-Nitroaniline	138	8.214	8.214	(1.100)	120129	50.0000	54.34
97	4,6-Dinitro-2-methylphenol	198	8.276	8.276	(0.880)	65675	50.0000	48.62
98	N-Nitrosodiphenylamine	169	8.317	8.317	(0.884)	273788	58.6000	47.44
100	Azobenzene	77	8.359	8.359	(0.889)	367990	50.0000	49.15
101	4-Bromophenyl-phenylether	248	8.804	8.804	(0.936)	92973	50.0000	50.02
108	Hexachlorobenzene	284	8.981	8.981	(0.955)	104824	50.0000	50.50
110	Pentachlorophenol	266	9.240	9.240	(0.982)	62906	50.0000	50.72
114	Phenanthrene	178	9.437	9.437	(1.003)	575211	50.0000	47.93
115	Anthracene	178	9.509	9.509	(1.011)	584548	50.0000	48.76
118	Carbazole	167	9.768	9.768	(1.039)	547701	50.0000	50.01
120	Di-n-Butylphthalate	149	10.473	10.473	(1.113)	662234	50.0000	50.26
126	Fluoranthene	202	11.302	11.302	(1.202)	567781	50.0000	52.80
127	Benzidine	184	11.582	11.582	(0.840)	159069	50.0000	18.61
128	Pyrene	202	11.665	11.665	(0.846)	595801	50.0000	45.20
134	3,3'-dimethylbenzidine	212	12.877	12.877	(0.934)	141696	50.0000	18.78
136	Butylbenzylphthalate	149	12.991	12.991	(0.942)	310154	50.0000	46.95
138	Benzo(a)Anthracene	228	13.758	13.758	(0.998)	523382	50.0000	46.59
139	Chrysene	228	13.830	13.830	(1.003)	551943	50.0000	48.03
140	3,3'-Dichlorobenzidine	252	13.799	13.799	(1.001)	198689	50.0000	46.89
141	bis(2-ethylhexyl)Phthalate	149	14.110	14.110	(1.023)	422505	50.0000	46.43
142	Di-n-octylphthalate	149	15.167	15.167	(1.100)	671608	50.0000	46.17
144	Benzo(b)fluoranthene	252	15.582	15.582	(0.963)	474456	50.0000	49.97(Q)
145	Benzo(k)fluoranthene	252	15.623	15.623	(0.966)	540533	50.0000	52.55
147	Benzo(e)pyrene	252	16.007	16.007	(0.990)	515993	50.0000	52.12
148	Benzo(a)pyrene	252	16.079	16.079	(0.994)	500123	50.0000	46.46
151	Indeno(1,2,3-cd)pyrene	276	17.810	17.810	(1.101)	429096	50.0000	49.29
152	Dibenzo(a,h)anthracene	278	17.851	17.851	(1.104)	519505	50.0000	53.41
153	Benzo(g,h,i)perylene	276	18.235	18.235	(1.127)	569749	50.0000	54.10

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
M 162 benzo b,k Fluoranthene Totals	252				1114989	50.0000	

QC Flag Legend

Q - Qualifier signal failed the ratio test.

TestAmerica West Sacramento

INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002H.D
Lab Smp Id: HSL_050 ug/ml ICV
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\SV5\C\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;1_8270STD.SUB;10MSSV0314;0;8270F.M

Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	98364	-19.78
2 Naphthalene-d8	530514	265257	1061028	431655	-18.63
3 Acenaphthene-d10	282538	141269	565076	236662	-16.24
4 Phenanthrene-d10	462722	231361	925444	380734	-17.72
5 Chrysene-d12	435850	217925	871700	421719	-3.24
6 Perylene-d12	422284	211142	844568	419419	-0.68

COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.95	3.45	4.45	3.95	0.00
2 Naphthalene-d8	5.37	4.87	5.87	5.37	0.00
3 Acenaphthene-d10	7.47	6.97	7.97	7.47	0.00
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	0.00
5 Chrysene-d12	13.79	13.29	14.29	13.79	0.00
6 Perylene-d12	16.17	15.67	16.67	16.17	0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\SV5\C\ohem\sv5.i\100210.B\HSL1002H.D

Date : 02-OCT-2010 16:11

Client ID: 8270F.H

Sample Info: HSL_050 ug/ml ICV:2::4:::::4

Column phase:

Instrument: sv5.i

Operator: KT

Column diameter: 2.00

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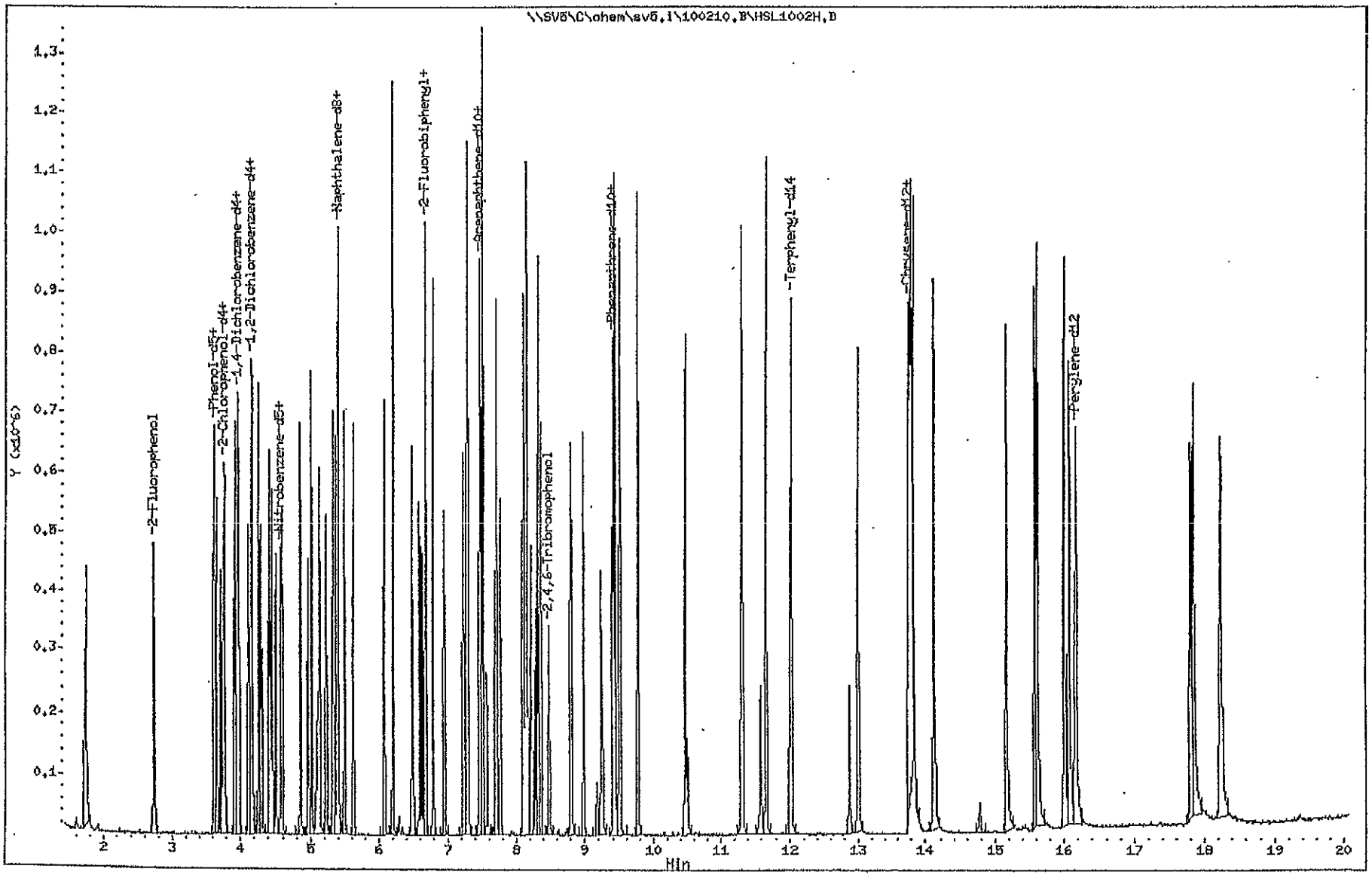

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CONTINUING CALIBRATION COMPOUNDS

Instrument ID: sv5.i Injection Date: 02-OCT-2010 16:36
Lab File ID: HSL1002H1.D Init. Cal. Date(s): 17-AUG-2010 02-OCT-2010
Analysis Type: Init. Cal. Times: 17:32 15:00
Lab Sample ID: Benzidines ICV 50ug Quant Type: ISTD
Method: \\sv5\c\chem\sv5.i\100210.B\8270f.m

COMPOUND	RRF / AMOUNT	RF50	CCAL RRF50	MIN RRF	%D / %DRIFT	MAX %D / %DRIFT	CURVE TYPE
127 Benzidine	0.81067	0.92336	0.92336	0.010	13.89989	50.00000	Averaged
134 3,3'-dimethylbenzidine	0.71564	0.78974	0.78974	0.010	10.35398	50.00000	Averaged
140 3,3'-Dichlorobenzidine	0.40189	0.42433	0.42433	0.010	5.58428	50.00000	Averaged

10-3-10

TestAmerica West Sacramento

Method 8270C

Data file : \\sv5\c\chem\sv5.i\100210.B\HSL1002H1.D
Lab Smp Id: Benzidines ICV 50ug Client Smp ID: 8270F.M
Inj Date : 02-OCT-2010 16:36
Operator : KT Inst ID: sv5.i
Smp Info : Benzidines ICV 50ug/mL;2;;4;;;;4
Misc Info : 3;;0;BenzICV.SUB;10MSSV0342;0;8270F.M
Comment : SOP SAC-MS-0005
Method : \\sv5\c\chem\sv5.i\100210.B\8270f.m
Meth Date : 03-Oct-2010 11:13 truongk Quant Type: ISTD
Cal Date : 17-AUG-2010 21:19 Cal File: AP90817D.D
Als bottle: 9 Continuing Calibration Sample
Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: BenzICV.SUB
Target Version: 4.14
Processing Host: SACP307UM

Compounds	QUANT SIG MASS	RT	EXP RT	REL RT	RESPONSE	AMOUNTS CAL-AMT (NG)	ON-COL (NG)
* 1 1,4-Dichlorobenzene-d4	152	3.954	3.954	(1.000)	115503	40.0000	
* 2 Naphthalene-d8	136	5.364	5.364	(1.000)	480485	40.0000	
* 3 Acenaphthene-d10	164	7.468	7.468	(1.000)	254190	40.0000	
* 4 Phenanthrene-d10	188	9.405	9.405	(1.000)	405333	40.0000	
* 5 Chrysene-d12	240	13.779	13.779	(1.000)	378068	40.0000	
* 6 Perylene-d12	264	16.162	16.162	(1.000)	372382	40.0000	
127 Benzidine	184	11.571	11.571	(0.840)	436364	50.0000	56.95
134 3,3'-dimethylbenzidine	212	12.867	12.867	(0.934)	373217	50.0000	55.18
140 3,3'-Dichlorobenzidine	252	13.799	13.799	(1.002)	200534	50.0000	52.79

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INTERNAL STANDARD COMPOUNDS
AREA AND RT SUMMARY

Instrument ID: sv5.i
Lab File ID: HSL1002H1.D
Lab Smp Id: Benzidines ICV 50ug
Analysis Type: SV
Quant Type: ISTD
Operator: KT
Method File: \\sv5\c\chem\sv5.i\100210.B\8270f.m
Misc Info: 3;;0;BenzICV.SUB;10MSSV0342;0;8270F.M

Calibration Date: 02-OCT-2010
Calibration Time: 13:44
Client Smp ID: 8270F.M
Level:
Sample Type:

Test Mode:
Use Initial Calibration Level 4.

COMPOUND	STANDARD	AREA LIMIT LOWER	AREA LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	122625	61313	245250	115503	-5.81
2 Naphthalene-d8	530514	265257	1061028	480485	-9.43
3 Acenaphthene-d10	282538	141269	565076	254190	-10.03
4 Phenanthrene-d10	462722	231361	925444	405333	-12.40
5 Chrysene-d12	435850	217925	871700	378068	-13.26
6 Perylene-d12	422284	211142	844568	372382	-11.82

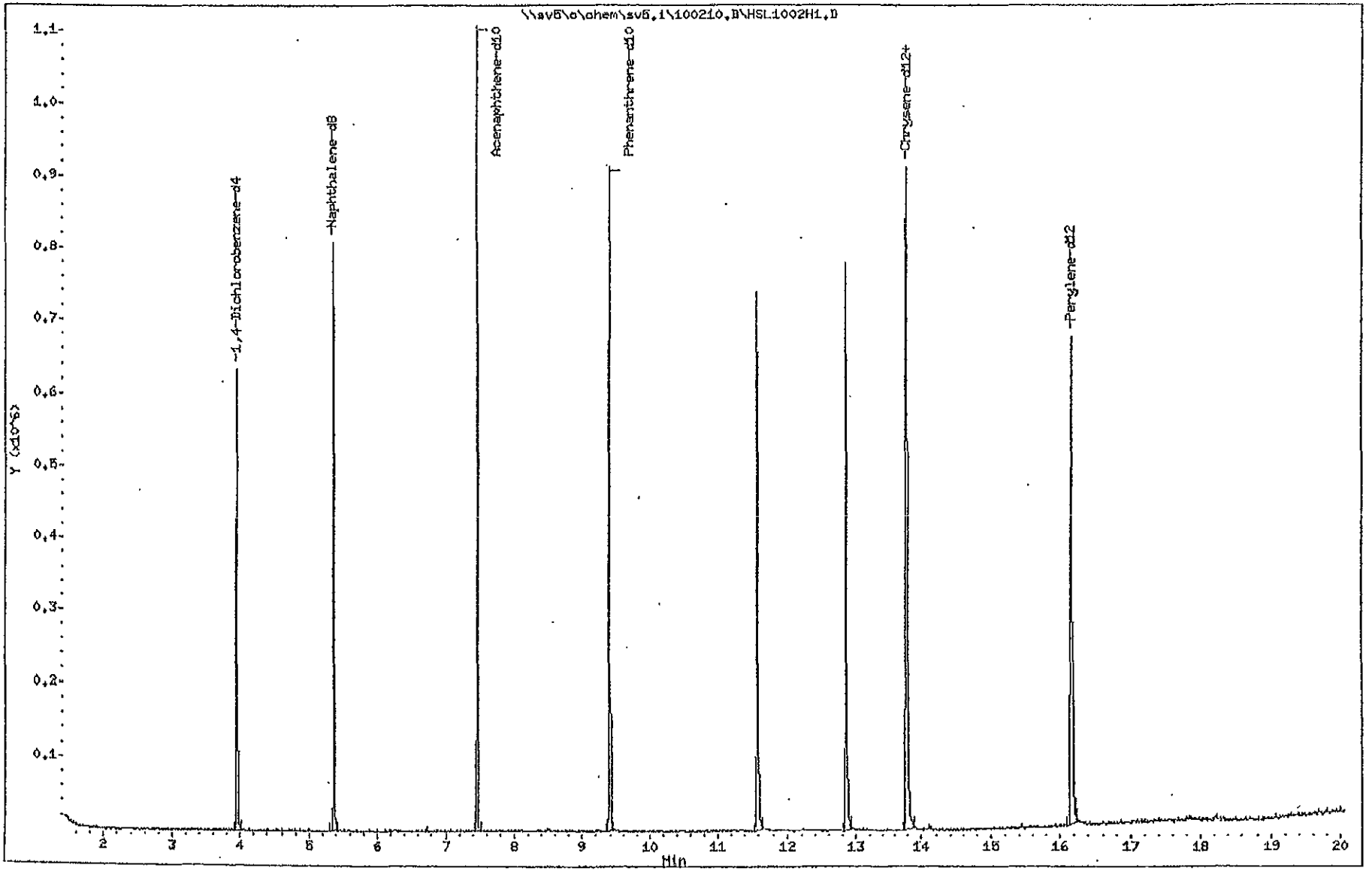
COMPOUND	STANDARD	RT LIMIT LOWER	RT LIMIT UPPER	SAMPLE	%DIFF
1 1,4-Dichlorobenze	3.95	3.45	4.45	3.95	0.00
2 Naphthalene-d8	5.36	4.86	5.86	5.36	0.00
3 Acenaphthene-d10	7.47	6.97	7.97	7.47	0.00
4 Phenanthrene-d10	9.41	8.91	9.91	9.41	0.00
5 Chrysene-d12	13.78	13.28	14.28	13.78	0.00
6 Perylene-d12	16.16	15.66	16.66	16.16	0.00

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = - 50% of internal standard area.
RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: \\sv5\o\ohem\sv5.i\100210.B\HSL1002H1.D
Date : 02-OCT-2010 16:36
Client ID: 8270F.M
Sample Info: Benzidines ICV 60ug/mL;2;;4;;;;;4

Instrument: sv5.i
Operator: KT
Column diameter: 2.00

Column phase:



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INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Origin : Disabled
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:07 sv5.i
Curve Type : Average

Calibration File Names:
Level 1: \\SV5\C\chem\sv5.i\081710.B\AP90817A.D
Level 2: \\SV5\C\chem\sv5.i\081710.B\AP90817B.D
Level 3: \\SV5\C\chem\sv5.i\081710.B\AP90817C.D
Level 4: \\SV5\C\chem\sv5.i\081710.B\AP90817D.D
Level 5: \\SV5\C\chem\sv5.i\081710.B\AP90817E.D
Level 6: \\SV5\C\chem\sv5.i\081710.B\AP90817F.D
Level 7: \\SV5\C\chem\sv5.i\081710.B\AP90817G.D

10/3/10

Original RRF

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	80.000 Level 5	120.000 Level 6	160.000 Level 7	RRF	% RSD
15 N-Nitrosodimethylamine	0.92899	0.88268	0.91048	0.91970	0.93146	0.93816	0.93833	0.92154	2.162
16 Pyridine	1.67117	1.37423	1.59449	1.56610	1.52299	1.53256	1.52623	1.54111	5.856
23 Aniline	2.20796	2.15935	2.19988	2.26058	2.29749	2.33400	2.33783	2.25573	3.098
24 Phenol	2.04111	1.96212	2.02834	2.03430	2.06683	2.06089	2.06740	2.03729	1.802
26 Bis(2-chloroethyl)ether	1.47335	1.38252	1.39491	1.43824	1.42549	1.44300	1.44264	1.42859	2.170
27 2-Chlorophenol	1.52099	1.55595	1.56903	1.58168	1.56789	1.58074	1.57039	1.56381	1.328
28 1,3-Dichlorobenzene	1.68903	1.69173	1.67754	1.73135	1.68641	1.72299	1.72457	1.70337	1.294
29 1,4-Dichlorobenzene	1.77122	1.79861	1.74013	1.76898	1.78200	1.79288	1.81444	1.78118	1.352

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Origin : Disabled
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:07 sv5.i
Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	80.000 Level 5	120.000 Level 6	RRF	% RSD
	160.000 Level 7							
30 Benzyl Alcohol	1.01643	1.03654	0.99182	1.04980	1.07792	1.08952		
	1.09506						1.05101	3.697
31 1,2-Dichlorobenzene	1.62008	1.63185	1.60455	1.68051	1.63410	1.64415		
	1.64691						1.63746	1.459
32 2-Methylphenol	1.40818	1.38930	1.39110	1.42620	1.45565	1.46154		
	1.47889						1.43012	2.506
33 2,2'-oxybis(1-Chloropropane)	2.29602	2.22080	2.28329	2.27928	2.27018	2.27830		
	2.28770						2.27365	1.085
34 4-Methylphenol	1.48606	1.48913	1.46270	1.52239	1.52653	1.55886		
	1.58763						1.51904	2.884
35 Hexachloroethane	0.60925	0.60836	0.60573	0.61394	0.60427	0.59381		
	0.60919						0.60636	1.043
37 N-Nitrosodinpropylamine	0.94498	0.97005	1.01302	1.02370	1.04700	1.03627		
	1.04757						1.01180	3.926
42 Nitrobenzene	0.32855	0.32602	0.32543	0.33083	0.33379	0.33450		
	0.33901						0.33116	1.489
44 Isophorone	0.63431	0.62291	0.61160	0.63344	0.63648	0.65468		
	0.65411						0.63679	2.811
45 2-Nitrophenol	0.18608	0.18833	0.18840	0.20021	0.20022	0.20702		
	0.20508						0.19648	4.423
46 2,4-Dimethyphenol	0.34459	0.34167	0.34307	0.34912	0.34788	0.35962		
	0.35785						0.34911	2.028

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Origin : Disabled
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:07 sv5.i
Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	80.000 Level 5	120.000 Level 6	RRF	% RSD
	160.000 Level 7							
47 Bis(2-chloroethoxy)methane	0.41146	0.37494	0.38565	0.38249	0.38500	0.39859		
	0.38545						0.38908	3.106
49 2,4-Dichlorophenol	0.25434	0.26318	0.27019	0.27037	0.27274	0.28180		
	0.27809						0.27010	3.393
50 Benzoic Acid	0.16747	0.16266	0.17423	0.19357	0.21024	0.22272		
	0.22180						0.19324	13.252
51 1,2,4-Trichlorobenzene	0.29430	0.28827	0.28475	0.29747	0.29189	0.29959		
	0.29091						0.29246	1.760
52 Naphthalene	1.09939	1.12462	1.07435	1.09325	1.09870	1.13821		
	1.10247						1.10443	1.900
54 4-Chloroaniline	0.40751	0.42534	0.43264	0.43910	0.43781	0.44905		
	0.43867						0.43288	3.068
57 Hexachlorobutadiene	0.14295	0.13812	0.14428	0.14415	0.14385	0.14379		
	0.14473						0.14313	1.589
60 4-Chloro-3-Methylphenol	0.29329	0.28866	0.29079	0.30972	0.30295	0.31766		
	0.30839						0.30164	3.644
63 2-Methylnaphthalene	0.68483	0.68064	0.68080	0.70067	0.70560	0.71172		
	0.69217						0.69378	1.797
66 Hexachlorocyclopentadiene	0.26878	0.27757	0.28896	0.29704	0.30236	0.32262		
	0.33186						0.29846	7.645
69 2,4,6-Trichlorophenol	0.31186	0.29820	0.30223	0.31996	0.32305	0.34225		
	0.33638						0.31913	5.157

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Origin : Disabled
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:07 sv5.i
Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	80.000 Level 5	120.000 Level 6	160.000 Level 7	RRF	% RSD
70 2,4,5-Trichlorphenol	0.30823	0.32892	0.33796	0.36298	0.35236	0.35480	0.36135	0.34380	5.807
71 2-Chloronaphthalene	1.13629	1.09411	1.10012	1.14181	1.11220	1.14447	1.15096	1.12571	2.051
73 2-Nitroaniline	0.31576	0.31759	0.33397	0.35205	0.34821	0.35796	0.36278	0.34119	5.573
76 Dimethylphthalate	1.23388	1.25191	1.29803	1.34568	1.31165	1.32891	1.30237	1.29606	3.093
77 Acenaphthylene	1.86531	1.91304	1.91818	2.01646	1.98204	1.99786	2.02968	1.96037	3.150
79 2,6-Dinitrotoluene	0.28347	0.27378	0.29890	0.31220	0.31294	0.32140	0.31106	0.30197	5.786
80 3-Nitroaniline	0.35362	0.34622	0.35978	0.40036	0.38674	0.39559	0.39603	0.37691	6.069
81 Acenaphthene	1.25874	1.22468	1.26733	1.27046	1.21141	1.24781	1.25463	1.24787	1.768
82 2,4-Dinitrophenol	0.10149	0.11058	0.14485	0.16667	0.18378	0.20563	0.20232	0.15933	26.349
83 Dibenzofuran	1.57786	1.62124	1.65200	1.69530	1.65117	1.68450	1.71077	1.65612	2.779
84 4-Nitrophenol	0.12712	0.14148	0.15316	0.16076	0.17130	0.16653	0.17404	0.15634	10.909

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Origin : Disabled
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:07 sv5.i
Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	80.000 Level 5	120.000 Level 6	RRF	% RSD
	160.000 Level 7							
86 2,4-Dinitrotoluene	0.34360	0.35989	0.38479	0.42154	0.41035	0.42305		
	0.43110						0.39633	8.616
91 Fluorene	1.34567	1.33840	1.34292	1.39902	1.38899	1.37835		
	1.40640						1.37139	2.086
92 Diethylphthalate	1.22240	1.29889	1.31549	1.37912	1.31873	1.37345		
	1.38087						1.32699	4.319
93 4-Chlorophenyl-phenylether	0.54964	0.55917	0.56887	0.59265	0.56708	0.57695		
	0.57695						0.57019	2.429
94 4-Nitroaniline	0.33346	0.33747	0.37329	0.38337	0.39216	0.39102		
	0.40452						0.37361	7.424
97 4,6-Dinitro-2-methylphenol	0.09316	0.10533	0.12545	0.13163	0.14105	0.15288		
	0.15229						0.12883	17.707
98 N-Nitrosodiphenylamine	0.57756	0.59736	0.60533	0.60433	0.62172	0.61801		
	0.61968						0.60628	2.577
100 Azobenzene	0.77527	0.76965	0.77321	0.79522	0.80064	0.81892		
	0.77331						0.78660	2.371
101 4-Bromophenyl-phenylether	0.18964	0.18507	0.19281	0.19931	0.19607	0.20581		
	0.19815						0.19527	3.488
108 Hexachlorobenzene	0.22958	0.22054	0.20740	0.21605	0.21731	0.21704		
	0.21854						0.21807	3.009
110 Pentachlorophenol	0.09427	0.09851	0.11582	0.11736	0.13228	0.13923		
	0.13770						0.11931	15.221

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Origin : Disabled
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:07 sv5.i
Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	80.000 Level 5	120.000 Level 6	RRF	% RSD
	160.000 Level 7							
114 Phenanthrene	1.30347	1.26007	1.25408	1.24163	1.24375	1.25610		
	1.26611						1.26074	1.643
115 Anthracene	1.25034	1.21759	1.24206	1.25982	1.27529	1.30214		
	1.26958						1.25955	2.329
118 Carbazole	1.13211	1.12547	1.13694	1.14260	1.17067	1.18192		
	1.16455						1.15061	1.878
120 Di-n-Butylphthalate	1.28492	1.32287	1.36193	1.38164	1.41474	1.43847		
	1.48636						1.38442	4.973
126 Fluoranthene	1.03840	1.07611	1.17216	1.10520	1.15861	1.18294		
	1.17440						1.12969	5.018
127 Benzidine	0.78175	0.76431	0.75250	0.82658	0.82201	0.86375		
	0.86381						0.81067	5.606
128 Pyrene	1.25791	1.23783	1.17078	1.28684	1.25586	1.28463		
	1.25794						1.25025	3.122
134 3,3'-dimethylbenzidine	0.65472	0.64388	0.67361	0.70756	0.73630	0.79414		
	0.79926						0.71564	8.888
136 Butylbenzylphthalate	0.64984	0.60187	0.59142	0.62586	0.61590	0.65233		
	0.64920						0.62663	3.950
138 Benzo(a)Anthracene	1.10169	0.99732	1.03245	1.04489	1.06449	1.10831		
	1.10920						1.06548	4.058
139 Chrysene	1.05284	1.10175	1.06320	1.09705	1.06985	1.12241		
	1.12246						1.08994	2.594

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Origin : Disabled
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:07 sv5.i
Curve Type : Average

Compound	5.000 Level 1	10.000 Level 2	20.000 Level 3	50.000 Level 4	80.000 Level 5	120.000 Level 6	RRF	% RSD
	160.000 Level 7							
140 3,3'-Dichlorobenzidine	0.39148	0.37695	0.39090	0.39906	0.40353	0.42717		
	0.42415						0.40189	4.539
141 bis(2-ethylhexyl)Phthalate	0.91826	0.80897	0.84032	0.85193	0.84371	0.89539		
	0.88354						0.86316	4.348
142 Di-n-octylphthalate	1.34838	1.23185	1.35627	1.34433	1.39356	1.47616		
	1.50770						1.37975	6.651
144 Benzo(b)fluoranthene	0.81012	0.81077	0.82747	0.99930	0.95373	0.91132		
	1.02572						0.90549	10.058
145 Benzo(k)fluoranthene	1.22939	1.16528	1.20022	1.09895	1.14223	1.19597		
	1.10447						1.16236	4.279
147 Benzo(e)pyrene	0.90394	0.92734	0.90757	0.95977	0.96997	0.96929		
	0.97185						0.94425	3.220
148 Benzo(a)pyrene	0.98300	0.97686	0.99402	1.02789	1.07610	1.06275		
	1.06523						1.02655	4.111
151 Indeno(1,2,3-cd)pyrene	0.73783	0.73267	0.73671	0.84698	0.84057	0.93730		
	0.97995						0.83029	12.151
152 Dibenzo(a,h)anthracene	0.88099	0.84384	0.87256	0.92240	0.95990	1.00944		
	1.00392						0.92758	7.071
153 Benzo(g,h,i)perylene	0.96025	0.98457	0.97380	0.99974	1.01731	1.05397		
	1.04026						1.00427	3.452
M 162 benzo b,k Fluoranthene Totals	2.03951	1.97605	2.02770	2.09825	2.09596	2.10729		
	2.13019						2.06785	2.649

TestAmerica West Sacramento

INITIAL CALIBRATION DATA

Start Cal Date : 17-AUG-2010 17:32
End Cal Date : 02-OCT-2010 15:00
Quant Method : ISTD
Origin : Disabled
Target Version : 4.14
Integrator : Falcon
Method file : \\SV5\C\chem\sv5.i\100210.B\8270f.m
Last Edit : 03-Oct-2010 11:07 sv5.i
Curve Type : Average

Compound	5.000 Level 1 / 160.000 Level 7	10.000 Level 2	20.000 Level 3	50.000 Level 4	80.000 Level 5	120.000 Level 6	RRF	% RSD
$ 7 2-Fluorophenol	1.44503 / 1.43635	1.30436	1.38373	1.44170	1.43535	1.42292	1.40952	3.615
$ 8 Phenol-d5	1.72227 / 1.83627	1.67335	1.74151	1.79006	1.80863	1.83864	1.77296	3.520
$ 9 2-Chlorophenol-d4	1.47770 / 1.57804	1.55530	1.53916	1.59414	1.57486	1.57967	1.55698	2.524
$ 10 1,2-Dichlorobenzene-d4	0.95776 / 0.98896	0.98111	0.99827	0.98914	0.99518	0.98547	0.98513	1.356
$ 11 Nitrobenzene-d5	0.33028 / 0.33970	0.34256	0.33065	0.34105	0.33606	0.35127	0.33879	2.162
$ 12 2-Fluorobiphenyl	1.28499 / 1.30010	1.26007	1.27668	1.34206	1.25854	1.29723	1.28852	2.226
$ 13 2,4,6-Tribromophenol	0.15034 / 0.18390	0.16527	0.17466	0.17926	0.17825	0.18501	0.17381	7.052
$ 14 Terphenyl-d14	0.78508 / 0.80107	0.78616	0.73917	0.80441	0.78047	0.81889	0.78789	3.214

Sample Extraction/Preparation Log Copies and Checklists

TestAmerica West Sacramento
Organic Prep Log
8270 Air

ECJ 12/1/10

Box # Air Tox #290 291

Shared QC Batch: N/A

Shares QC With: N/A

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

Internal COC:	
Delivered to Inst.:	12/1/10
Inst Receipt:	

Batch: 0334389

MS Run #:

Prep Date: 11/30/2010 *RUSH*

Method: JZ TO-13

Matrix: S AIR

Extraction: 11 SOXHLET (NONE,Na2SO4)

QC: 3W AMBIENT AIR TESTING

SAC: JZ - S - 11 - 3W

WS-OP-0006

Prep Reagents		
Reagent	Supplier	Lot #
1:1 DCM:Acetone	NA	N/A
DCM	Baker	J34800
Na2SO4	Baker	3640-54B

Soxhlet time on: 16:45 (11/30/10) Soxhlet time off: 8:50 (12/1/10)

Extraction Table

Sample ID	Suff	Work Order	Extraction Hold Time Expires	Sample size	Final Volume 1mL	Final Volume Other	Analysis Hold Time Expires
G0K300000 - 389	B	MANG51AA	12/1/2010	1.0	✓		1/3/2011
G0K300000 - 389	C	MANG51AC	12/1/2010	1.0	✓		1/3/2011
G0K300000 - 389	L	MANG51AD	12/1/2010	1.0	✓		1/3/2011
G0K300434 - 1		MAMLW1AA	12/1/2010	1.0	✓		1/3/2011
G0K300434 - 4		MAML81AA	12/1/2010	1.0	✓		1/3/2011

☑ XAD / PUF / PUF-XAD
☐ Filter
☐ Impinger

Comments/NCMs: QC Media: syp2sv19056/ ECJ 11/30/10
P101910

	ID	Spike Exp Date:	Spiked By:	Witnessed By:	Date:
Surrogate Spike All Samples	500µL / 10AIR0125 / ABN Surr	4/4/11	ECJ		11/30/10 ✓
Spike Mix LCS/LCSD~~AMS/MS~~ ECJ 11/30/10	1.0mL / 10AIR0126 / 8270 LCS Mix	4/9/11	ECJ		11/30/10 ✓
Pre-Spike Standard ~~All Samples~~ MB only ECJ 11/30/10	250µL / 10AIR0128 / 1,2-DCB-d4	4/19/11	ECJ		11/30/10 ✓
Internal Standard All Samples	20mL / 10IS8V0438	11-19-11		VMD	12/3/10
Soxhlet Extraction Analyst/Date	ECJ 11/30/10	Concentration Analyst/Date	ECJ 12/1/10	KD Analyst/Date	ECJ 12/1/10
Liq Liq Extraction Analyst/Date	N/A	KD Temp	85°C	Review Analyst/Date	

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

West Sacramento

Preparation Data Review Checklist

Prep Batch(es) 0334399 **Test:** TO-13

Prep Date: 11/30/10 **Holding Times:** 12/1/10 **NCM:** Y (N)

A. Spike Witness/Batch setup	Spike Witness	Reviewer
1. **Holding times checked?** NCMs filed as appropriate	✓	
2. QAS checked for QC instructions (LCS, LCSD, MS,MSD, etc)	✓	
3. Amount of samples in hood match amount of samples on bench sheet. Sample IDS match.	✓	NA
4. Worksheets have been checked for required spiking compounds	✓	
5. Spiking volumes are correctly documented	✓	
6. Std ID numbers on spike labels match numbers on bench sheet	✓	NA
7. Expiration dates have been checked	✓	
8. Calibration expiration dates on pipettors have been checked	✓	NA
9. Spiker and spike witness have signed and dated bench sheet	✓	
B. Weights and Volumes		
1. Recorded weights are in anticipated range	NA	
2. Balance upload or raw data for weights is included	NA	
3. Weights and volumes have been transcribed correctly to LIMS.	NA	
4. Weights are not targeted to meet exact weights.	NA	
5. Each weight or volume measurement is a unique record (no dittos or line downs)	NA	
C. Standards and Reagents		
1. Lot numbers for all reagents, including clean up stages, are recorded.	NA	
2. Are dates and analysts for cleanups recorded?	NA	
3. Are correct IDs used for standards? Are expiration dates to day/month/year, when listed?	NA	
D. Documentation		
1. Are all nonconformances documented appropriately?	NA	
2. QuantIMs entry correct, including dates and times.	NA	
3. Are all fields completed?	NA	

Spike witness: _____ Date: 11/30/10

2nd Level Reviewer: _____ Date: _____

Comments:

_____

_____

_____

_____

RQC058

TestAmerica Laboratories, Inc.
EXTRACTION BENCH WORKSHEET

Run Date: 12/01/10
Time: 11:47:49

LEV 1	LEV 2		LEV 1	LEV 2	
Y	_	Blank	Y	_	Weights/Volumes
Y	_	Check	Y	_	Spike & Surrogate Worksheet
_	_	MS/MSD	Y	_	Vial contains correct volume
			Y	_	Labels, greenbars, worksheets computer batch: correct & all match
			_	_	Anomalies to Extraction Method

_ Expanded Deliverable
_ COC Completed
_ Bench Sheet Copied
_ Package Submitted to AnalyticalGr
_ Bench Sheet Copied per COC

Extractionist: 403162 erica X. larson

Concentrationist: 403162 erica X. larson

****
* *
* **QC BATCH: 0334389** *
* *
****

PREP DATE: 11/30/10 16:45
COMP DATE: 12/01/10 15:00

Reviewer/Date: LARSONE / 12/01/10

Semivolatile Organics by GCMS in Air (TO-13A)
SOXHLET (NONE,Na2SO4)

EXTR EXPR	ANL DUE	LOT#,MSRUN#/ WORK ORDER	TEST FLGS	EXT	MTH	MATRIX	INIT/FIN WT/VOL	PH"S INIT	PH"S ADJ1	PH"S ADJ2	SOLVENTS EXTRACTION	SOLVENTS VOL	SOLVENTS EXCHANGE	SOLVENTS VOL	SPIKE STANDARD/ SURROGATE ID
12/01/10	12/07/10	G0K300434-001 **MAMLW-1-AA**	R	11	JZ	AIR	**1.0Sample** 1.00mL	NA	NA	NA	DCM	700.0		.0	500UL/10AIR0125/ABN S
COMMENTS:															
12/01/10	12/07/10	G0K300434-004 **MAML8-1-AA**	R	11	JZ	AIR	**1.0Sample** 1.00mL	NA	NA	NA	DCM	700.0		.0	500UL/10AIR0125/ABN S
COMMENTS:															
12/01/10	0/00/00	G0K300000-389 **MANG5-1-AAB**		11	JZ	AIR	**1.0Sample** 1.00mL	NA	NA	NA	DCM	700.0		.0	250UL/10AIR0128/1,2-D 500UL/10AIR0125/ABN S
COMMENTS:															
12/01/10	0/00/00	G0K300000-389 **MANG5-1-ACC**		11	JZ	AIR	**1.0Sample** 1.00mL	NA	NA	NA	DCM	700.0		.0	1.0ML/10AIR0126/8270 500UL/10AIR0125/ABN S
COMMENTS:															
12/01/10	0/00/00	G0K300000-389 **MANG5-1-ADL**	R	11	JZ	AIR	**1.0Sample** 1.00mL	NA	NA	NA	DCM	700.0		.0	1.0ML/10AIR0126/8270 500UL/10AIR0125/ABN S
COMMENTS:															

QC MEDIA:SUP2SV19056/P101910

R = RUSH C = CLP
E = EPA 600 D = EXP.DEL)
M = CLIENT REQ MS/MSD

NUMBER OF WORK ORDERS IN BATCH: 5

TestAmerica West Sacramento

GC/MS Data Review Checklist

Batch: 0334389 **Method ID:** Semivolatile Organics by GCMS in Air (TO-13A)

NCM: (Y) N

A. Calibration/Instrument Run QC	Analyst	Reviewer	N/A
1. ICAL or ICAL Summary and CCV included.	✓	✓	
2. ICAL, CCV Criteria met.	✓	✓	
3. Peaks correctly ID'd by data system.	✓	✓	
4. Copy of logbook for ICAL included	✓	✓	
5. Tune criteria (including tailing factor and breakdown) met and copy included.	✓	✓	
6. Method Number is identified on data.	✓	✓	
B. QA/QC			
1. Method blank, LCS/LCSD and MS/SD frequencies met.	✓	✓	
2. LCS/LCSD and MB data is included.	✓	✓	
3. LCS/LCSD and MB data are within control limits. If not, NCM is present in Clouseau.	✓	✓	
4. MS/MSD data complete.			✓
5. Holding Times were met.	✓	✓	
6. All samples within tune time.	✓	✓	
C. Sample Analysis			
1. Logbook copies for all injections made, including ICV standards and ICAL.	✓	✓	
2. Logbooks/prep sheets properly filled out.	✓	✓	
3. Manual Integrations reviewed and appropriate.	✓	✓	
4. All raw data for samples is included (applies to unused data as well)	✓	✓	
5. All analytes correctly reported.	✓	✓	
6. Correct reporting limits used. (based on client request, prep factors, and dilutions)	✓	✓	
7. Spectra present for all positives.	✓	✓	
D. Documentation			
1. Are all nonconformances documented appropriately?	✓	✓	
2. Quantims entry correct, including dates and times.	✓	✓	
3. Appropriate footnotes used.	✓	✓	

Analyst: [signature] Date: ~~12/6/10~~ 12/7/10

12/7/10 mc

2nd Level Reviewer: [signature] Date: 12/7/10

Comments: _____

AIR, Metals by ICPMS (As and Mn)

Raw Data Package

ICPMS

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

ICP-MS Data Review Checklist
Level I and Level II

Instrument ID (Circle one): **M01** (**M02**)		Method 6020 SOP SAC-MT-0001	
File Number	Batch Numbers	Date	Analyst
101207A2	336286,	12-07-10	SH

Lot Numbers: G0K190601, G0K240587, G0K270427, G0K300434, G0L020446	YES	NO	NA
1. Copy of analysis protocol used included?	✓		
2. ICVs & CCVs within 10% of true value or recal and rerun?	✓		
3. ICB & CCBs < reporting limit or recal and rerun?	✓		
4. 10 samples or less analyzed between calibration checks?	✓		
5. All parameters within linear range?	✓		
6. LCS/LCSD within limits?	✓		
7. Prep blank value < reporting limit or all samples >20x blank?	✓		
8. Internal standard intensities for samples (unless followed by dilution) are > 30% and <120% of the Calibration Blank intensities?	✓		
9. Appropriate dilution factors applied to data?	✓		
10. Matrix spike and spike dup within customer defined limits?			✓
11. Each batch checked for presence of internal standard in samples?	✓		
12. Anomalies entered using Clouseau?			✓

COMMENTS: _____

REVIEWED BY: MTZ	DATA ENTERED BY: SH
DATE: 12/8/10	DATE: 12-7-10

Dataset Report

Perkin Elmer M02
SOP No. SAC-MT-0001
Method: 6020,200.8

User Name: metal
Computer Name: SACP1223
Dataset File Path: e:\elandata\dataset\101207a2\
Report Date/Time: Tuesday, December 07, 2010 14:13:29

The Dataset

Batch ID	Sample ID	Date and Time	Read Type	Description	
	TUNE SHARGRAVE	06:49:13 Tue 07-Dec-10	Sample		
	AUTOLENS SHARGF	06:53:04 Tue 07-Dec-10	Sample		
	DAILY SHARGRAVE	07:34:13 Tue 07-Dec-10	Sample		
	Rinse 2X	09:12:39 Tue 07-Dec-10	Sample		
	Blank	09:15:34 Tue 07-Dec-10	Blank		
	Standard 1	09:18:24 Tue 07-Dec-10	Standard #1		
	ICV	09:20:59 Tue 07-Dec-10	Sample		
	ICV	09:25:16 Tue 07-Dec-10	Sample		
	ICB	10:24:20 Tue 07-Dec-10	Sample		
	LLSTD1	10:26:59 Tue 07-Dec-10	Sample	LLSTD@10X → out Al	
	LLSTD2	10:29:38 Tue 07-Dec-10	Sample	LLSTD@5X	
	ICSA	10:32:16 Tue 07-Dec-10	Sample		
	ICSAB	10:34:53 Tue 07-Dec-10	Sample		
	Rinse	11:11:19 Tue 07-Dec-10	Sample		
336286	MARD8B	11:17:36 Tue 07-Dec-10	Sample	G0L020000-286 BLK	} don't use
336286	MARD8C	11:20:11 Tue 07-Dec-10	Sample	G0L020000-286 LCS	
336286	MARD8L	11:22:45 Tue 07-Dec-10	Sample	G0L020000-286 LCSD	
	CCV 1	11:25:24 Tue 07-Dec-10	Sample		
	CCB 1 > Reread	11:50:24 Tue 07-Dec-10	Sample	> 7 IS	
	CCV 2	12:03:37 Tue 07-Dec-10	Sample		
	CCB 2	12:06:16 Tue 07-Dec-10	Sample		
	LLSTD1	12:08:56 Tue 07-Dec-10	Sample	LLSTD@10X ✓	
335251	MAPEVB	12:23:45 Tue 07-Dec-10	Sample	G0L010000-251 BLK	} report Al
335251/53	MAPEVC	12:26:21 Tue 07-Dec-10	Sample	G0L010000-251 LCS	
335253/51	MAPE7L	12:28:56 Tue 07-Dec-10	Sample	G0L010000-251 LCSD	
335251	MAA80	12:31:29 Tue 07-Dec-10	Sample	G0K190601-3	
335251	MAA80P5	12:34:01 Tue 07-Dec-10	Sample	G0K190601-3 5X	
335251	MAA80Z	12:36:33 Tue 07-Dec-10	Sample	G0K190601-3 PS	
335251	MAA81	12:39:06 Tue 07-Dec-10	Sample	G0K190601-4	
335251	MAKDV	12:41:39 Tue 07-Dec-10	Sample	G0K240587-1	
335251	MAKD2	12:44:12 Tue 07-Dec-10	Sample	G0K240587-2	
	CCV 3	12:46:51 Tue 07-Dec-10	Sample		
	CCB 3	12:49:31 Tue 07-Dec-10	Sample		
	CCV 4	12:55:08 Tue 07-Dec-10	Sample		
	CCB 4	12:57:47 Tue 07-Dec-10	Sample		
336286	MARD8B	13:00:24 Tue 07-Dec-10	Sample	G0L020000-286 BLK	} report As
336286	MARD8C	13:02:59 Tue 07-Dec-10	Sample	G0L020000-286 LCS	
336286	MARD8L	13:05:33 Tue 07-Dec-10	Sample	G0L020000-286 LCSD	
340010	MAWLKB	13:08:09 Tue 07-Dec-10	Sample	G0L060000-10 BLK	} report As, Mn
340010	MAWLKC	13:10:44 Tue 07-Dec-10	Sample	G0L060000-10 LCS	
340010	MAWLKL	13:13:20 Tue 07-Dec-10	Sample	G0L060000-10 LCSD	
340010	MAML1	13:15:54 Tue 07-Dec-10	Sample	G0K300434-2	
340010	MAML1P5	13:18:28 Tue 07-Dec-10	Sample	G0K300434-2 5X	
340010	MAML1Z	13:21:03 Tue 07-Dec-10	Sample	G0K300434-2 PS	
340010	MAML6	13:23:37 Tue 07-Dec-10	Sample	G0K300434-3	
	CCV 5	13:26:17 Tue 07-Dec-10	Sample		
	CCB 5	13:28:56 Tue 07-Dec-10	Sample		

341211	MA0J7B	13:39:35 Tue 07-Dec-10	Sample	G0L070000-211 BLK
341211	MA0J7C	13:42:11 Tue 07-Dec-10	Sample	G0L070000-211 LCS
341211	MA0J7L	13:44:47 Tue 07-Dec-10	Sample	G0L070000-211 LCSD
341211	MAQQ1	13:47:22 Tue 07-Dec-10	Sample	G0L020446-3
341211	MAQQ1P5	13:49:58 Tue 07-Dec-10	Sample	G0L020446-3 5X
341211	MAQQ1Z	13:52:33 Tue 07-Dec-10	Sample	G0L020446-3 PS
341211	MAQQ4	13:55:09 Tue 07-Dec-10	Sample	G0L020446-4
341211	MAQRA	13:57:44 Tue 07-Dec-10	Sample	G0L020446-7
341211	MAQRH	14:00:19 Tue 07-Dec-10	Sample	G0L020446-10
	CCV 6	14:02:57 Tue 07-Dec-10	Sample	
	CCB 6	14:05:37 Tue 07-Dec-10	Sample	

report As, Mn

TAL West Sac

RUN SUMMARY

Method: 6020 (SOP: SAC-MT-001) Instrument: M02 Reported: 12/07/10 14:26:16

File ID: 101207A2 **Analyst:** hargraves

#	Sample ID	Lot No.	Batch		DF	Analyzed Date	Comment	Q
1	Rinse 2X				2.0	12/07/10 09:12		☐
2	Blank				1.0	12/07/10 09:15		☐
3	Standard1				1.0	12/07/10 09:18		☐
4	ICV				1.0	12/07/10 09:20		☐
5	ICV				1.0	12/07/10 09:25		☐
6	ICB				1.0	12/07/10 10:24		☐
7	LLSTD1				1.0	12/07/10 10:26		☐
8	LLSTD2				1.0	12/07/10 10:29		☐
9	ICSA				1.0	12/07/10 10:32		☐
10	ICSAB				1.0	12/07/10 10:34		☐
11	Rinse				1.0	12/07/10 11:11		☐
12	MARD8B	G0L020000	0336286	2A	1.0	12/07/10 11:17		☐
13	MARD8C	G0L020000	0336286	2A	1.0	12/07/10 11:20		☐
14	MARD8L	G0L020000	0336286	2A	1.0	12/07/10 11:22		☐
15	CCV 1				1.0	12/07/10 11:25		☐
16	CCB 1				1.0	12/07/10 11:50		☐
19	CCV 2				1.0	12/07/10 12:03		☐
20	CCB 2				1.0	12/07/10 12:06		☐
21	LLSTD1				1.0	12/07/10 12:08		☐
22	MAPEVB	G0L010000	0335251	2A	1.0	12/07/10 12:23		☐
23	MAPEVC	G0L010000	0335251	2A	1.0	12/07/10 12:26		☐
24	MAPE7L	G0L010000	0335253	2A	1.0	12/07/10 12:28		☐
25	MAA80	G0K190601-3	0335251	2A	1.0	12/07/10 12:31		☐
26	MAA80P5	G0K190601	0335251		5.0	12/07/10 12:34		☐
27	MAA80Z	G0K190601-3	0335251		1.0	12/07/10 12:36		☐
28	MAA81	G0K190601-4	0335251	2A	1.0	12/07/10 12:39		☐
29	MAKDV	G0K240587-1	0335251	2A	1.0	12/07/10 12:41		☐
30	MAKD2	G0K240587-2	0335251	2A	1.0	12/07/10 12:44		☐
31	CCV 3				1.0	12/07/10 12:46		☐
32	CCB 3				1.0	12/07/10 12:49		☐
33	CCV 4				1.0	12/07/10 12:55		☐
34	CCB 4				1.0	12/07/10 12:57		☐
35	MARD8B	G0L020000	0336286	2A	1.0	12/07/10 13:00		☐
36	MARD8C	G0L020000	0336286	2A	1.0	12/07/10 13:02		☐
37	MARD8L	G0L020000	0336286	2A	1.0	12/07/10 13:05		☐
38	MAWLKB	G0L060000	0340010	2A	1.0	12/07/10 13:08		☐
39	MAWLKC	G0L060000	0340010	2A	1.0	12/07/10 13:10		☐
40	MAWLKL	G0L060000	0340010	2A	1.0	12/07/10 13:13		☐
41	MAML1	G0K300434-2	0340010	2A	1.0	12/07/10 13:15		☐
42	MAML1P5	G0K300434	0340010		5.0	12/07/10 13:18		☐
43	MAML1Z	G0K300434-2	0340010		1.0	12/07/10 13:21		☐
44	MAML6	G0K300434-3	0340010	2A	1.0	12/07/10 13:23		☐
45	CCV 5				1.0	12/07/10 13:26		☐
46	CCB 5				1.0	12/07/10 13:28		☐
47	MA0J7B	G0L070000	0341211	2A	1.0	12/07/10 13:39		☐
48	MA0J7C	G0L070000	0341211	2A	1.0	12/07/10 13:42		☐

Method: 6020 (SOP: SAC-MT-001) Instrument: M02 Reported: 12/07/10 14:26:16

File ID: 101207A2 Analyst: hargraves

#	Sample ID	Lot No.	Batch		DF	Analyzed Date	Comment	Q
49	MA0J7L	G0L070000	0341211	2A	1.0	12/07/10 13:44		☐
50	MAQQ1	G0L020446-3	0341211	2A	1.0	12/07/10 13:47		☐
51	MAQQ1P5	G0L020446	0341211		5.0	12/07/10 13:49		☐
52	MAQQ1Z	G0L020446-3	0341211		1.0	12/07/10 13:52		☐
53	MAQQ4	G0L020446-4	0341211	2A	1.0	12/07/10 13:55		☐
54	MAQRA	G0L020446-7	0341211	2A	1.0	12/07/10 13:57		☐
55	MAQRH	G0L020446-10	0341211	2A	1.0	12/07/10 14:00		☐
56	CCV 6				1.0	12/07/10 14:02		☐
57	CCB 6				1.0	12/07/10 14:05		☐

TAL West Sac | INTERNAL STANDARD SUMMARY

Method: 6020 (SOP: SAC-MT-001) | M02 (M02) | Reported: 12/07/10 14:26:16

File ID: 101207A2 | Analyst: hargraves

#	Sample ID	Analyzed Date	Germanium	Q
1	Rinse 2X	12/07/10 09:12	95.6	☐
2	Blank	12/07/10 09:15	100.0	☑
3	Standard1	12/07/10 09:18	99.6	☑
4	ICV	12/07/10 09:20	99.1	☑
5	ICV	12/07/10 09:25	99.9	☑
6	ICB	12/07/10 10:24	103.4	☑
7	LLSTD1	12/07/10 10:26	105.9	☑
8	LLSTD2	12/07/10 10:29	106.7	☑
9	ICSA	12/07/10 10:32	91.2	☑
10	ICSAB	12/07/10 10:34	97.1	☑
11	Rinse	12/07/10 11:11	126.2	☑
12	MARD8B	12/07/10 11:17	122.5	☑
13	MARD8C	12/07/10 11:20	114.4	☑
14	MARD8L	12/07/10 11:22	103.9	☑
15	CCV 1	12/07/10 11:25	119.3	☑
16	CCB 1	12/07/10 11:50	133.6	☐
19	CCV 2	12/07/10 12:03	102.5	☑
20	CCB 2	12/07/10 12:06	100.8	☑
21	LLSTD1	12/07/10 12:08	103.8	☑
22	MAPEVB	12/07/10 12:23	94.3	☑
23	MAPEVC	12/07/10 12:26	87.2	☑
24	MAPE7L	12/07/10 12:28	81.7	☑
25	MAA80	12/07/10 12:31	79.5	☑
26	MAA80P5	12/07/10 12:34	96.1	☐
27	MAA80Z	12/07/10 12:36	82.6	☑
28	MAA81	12/07/10 12:39	80.7	☑
29	MAKDV	12/07/10 12:41	85.3	☑
30	MAKD2	12/07/10 12:44	86.8	☑
31	CCV 3	12/07/10 12:46	94.8	☑
32	CCB 3	12/07/10 12:49	99.0	☑
33	CCV 4	12/07/10 12:55	96.7	☑
34	CCB 4	12/07/10 12:57	103.2	☑
35	MARD8B	12/07/10 13:00	97.4	☑
36	MARD8C	12/07/10 13:02	86.2	☑
37	MARD8L	12/07/10 13:05	81.1	☑
38	MAWLKB	12/07/10 13:08	79.0	☑
39	MAWLKC	12/07/10 13:10	80.8	☑
40	MAWLKL	12/07/10 13:13	79.6	☑
41	MAML1	12/07/10 13:15	80.0	☑
42	MAML1P5	12/07/10 13:18	90.0	☐
43	MAML1Z	12/07/10 13:21	77.6	☑
44	MAML6	12/07/10 13:23	78.6	☑
45	CCV 5	12/07/10 13:26	91.9	☑
46	CCB 5	12/07/10 13:28	95.0	☑
47	MA0J7B	12/07/10 13:39	91.8	☑
48	MA0J7C	12/07/10 13:42	82.1	☑

TAL West Sac

INTERNAL STANDARD SUMMARY

Method: 6020 (SOP: SAC-MT-001) M02 (M02) Reported: 12/07/10 14:26:16

File ID: 101207A2 **Analyst: hargraves**

#	Sample ID	Analyzed Date	Germanium	Q
49	MA0J7L	12/07/10 13:44	80.6	☑
50	MAQQ1	12/07/10 13:47	79.0	☑
51	MAQQ1P5	12/07/10 13:49	86.1	☐
52	MAQQ1Z	12/07/10 13:52	79.3	☑
53	MAQQ4	12/07/10 13:55	77.4	☑
54	MAQRA	12/07/10 13:57	80.6	☑
55	MAQRH	12/07/10 14:00	85.8	☑
56	CCV 6	12/07/10 14:02	93.4	☑
57	CCB 6	12/07/10 14:05	97.2	☑

Quantitative Method Report

File Name: 0006020-SH.mth
File Path: E:\elandata\Method\0006020-SH.mth

Timing Parameters

Sweeps/Reading:	50
Readings/Replicate:	1
Number of Replicates:	3
Tuning File:	default.tun
Optimization File:	default.dac
QC Enabled:	Yes
Settling Time:	Normal

Analyte	Mass	Scan Mode	MCA Channels	Dwell Time	Integration Time
Al	26.982	Peak Hopping	1	14.0 ms	700 ms
Ca	43.956	Peak Hopping	1	14.0 ms	700 ms
Mn	54.938	Peak Hopping	1	14.0 ms	700 ms
As	74.922	Peak Hopping	1	20.0 ms	1000 ms
Ge-1	71.922	Peak Hopping	1	14.0 ms	700 ms

Signal Processing

Detector Mode:	Dual
Measurement Units:	Counts
AutoLens:	On
Spectral Peak Processing:	Average
Signal Profile Processing:	Average
Blank Subtraction:	After Internal Standard
Baseline Readings:	0
Smoothing:	Yes, Factor 5

Equations

Analyte	Mass	Corrections
As	74.922	-3.1278 * Se 77 + 1.0177 * Se 78

Calibration Information

Analyte	Mass	Curve Type	Sample Units	Std Units	Std 1	Std 2	Std 3	Std 4
Al	26.982	Linear Thru Zero	ug/L	ug/L	5.1e+003			
Ca	43.956	Linear Thru Zero	ug/L	ug/L	5.1e+003			
Mn	54.938	Linear Thru Zero	ug/L	ug/L	100			
As	74.922	Linear Thru Zero	ug/L	ug/L	100			
Ge-1	71.922	Linear Thru Zero	ug/L	ug/L				

TAL-W. SACRMENTO – Perkin Elmer Elan 6000 ICPMS, M02 – Methods 6020, 200.8

AIR TOX Standards - 4 % HNO3, 0.5 % HCl

Standards for run:

Tuning standard: 4075-25B

Internal standard: 4075-22C

Blank, CCBs: 3185-42D

Standard 1, CCVs: 4075-21E

ICV: 4075-20D

ICSA: 4075-27B

ICSAB: 4075-27C

File Number: 101207A2

Instrument Tuning Report

File Name: default.tun

Sample Information

Sample Date/Time: Tuesday, December 07, 2010 06:49:13

Sample ID: TUNE SHARGRAVE

Analyte	Exact Mass	Meas. Mass	Mass DAC	Meas. Pk. Width	Res. DAC	Custom Res.
Li	7.016	7.027	1580	0.738	2040	
Be	9.012	9.029	2080	0.700	2035	
Mg	23.985	23.979	5728	0.726	2003	
Co	58.933	58.878	14249	0.724	1958	
In	114.904	114.829	27899	0.733	1937	
Ce	139.905	139.929	33978	0.729	1985	
Tl	204.975	204.979	49692	0.727	2189	
Pb	207.977	207.979	50425	0.723	2210	
U	238.050	238.028	57638	0.740	2360	

Elan 6000 Instrument Optomization Report

Path e:\elandata\Optimize
File Name e:\elandata\Optimize\default.dac

Sample Information

Sample Date/Time: Tuesday, December 07, 2010 07:34:13
Sample ID: DAILY SHARGRAVE

Parameter Settings

Nebulizer Gas Flow	0.92
Lens Voltage	9.00
ICP RF Power	1100.00
Analog Stage Voltage	-2000.00
Pulse Stage Voltage	1350.00
Discriminator Threshold	70.00
AC Rod Offset	-7.00
Service DAC 1	60.00
Quadrupole Rod Offset	0.00

AutoLens Calibration

Date:	06:53:04 Tue 07-Dec-10
Sample Filename:	AUTOLENS SHARGRAVE.003
Dataset Pathname:	e:\elandata\Dataset\101207a2\
Lens Voltage Start:	5.50
Lens Voltage End:	10.00
Lens Voltage Step:	0.25
Slope:	0.02126968
Intercept:	6.53696030

Analyte	Mass	Optimum Voltage	Maximum Intensity	# Points
Be	9.010	6.8	3281.1	19
Co	58.935	7.8	86613.6	19
In	114.903	9.0	328775.0	19

Dual Detector Calibration

Date:	08:01:56 Tue 07-Dec-10
Sample Filename:	DAILY SHARGRAVE.1097
Dataset Pathname:	dual detector calibration\
Points Acquired:	37
Lens Vol Start:	-3.00
Lens Vol End:	15.00
Lens Vol Step:	0.50

Analyte	Mass	Gain	N(max)
Li	6.015	10129.15	1235998078.193
Li	7.016	9449.89	1324842000.392
Be	9.012	8857.80	1413399228.860
B	11.009	9146.06	1368852249.810

Na	22.990	9118.55	1372981536.992
Mg	23.985	8534.23	1466986957.713
Mg	24.986	8369.25	1495905758.665
Al	26.982	7911.93	1582369824.835
Si	27.977	8973.22	1395219571.285
P	30.994	7287.08	1718055165.115
K	38.964	7100.33	1763243847.627
Ca	42.959		
Ca	43.956	6971.61	1795798956.472
Sc	44.956	7080.85	1768093466.146
V	50.944	6876.33	1820682362.555
Cr	51.941	6628.40	1888781365.139
Fe	53.940	6525.67	1918515148.090
Mn	54.938	6554.76	1910002204.770
Fe	56.935	6424.60	1948697279.431
Co	58.933	6269.19	1997004380.979
Ni	59.933	6081.55	2058620450.330
Cu	62.930	5973.72	2095780633.374
Cu	64.928	5888.17	2126231105.985
Zn	67.925	5952.75	2103161748.568
Ge	71.922	6127.47	2043191900.987
As	74.922	6124.98	2044023708.276
Se	77.917	6122.66	2044797442.453
Br	78.918		
Se	81.917	6050.71	2069114669.712
Sr	87.906		
Mo	96.906	6128.33	2042906515.828
Ag	106.905	5534.60	2262061270.648
Ag	108.905	5536.88	2261130679.220
Cd	110.904	5626.11	2225268192.020
Cd	113.904	5627.50	2224719836.750
In	114.904	5658.70	2212451551.812
Sn	117.902	5649.45	2216076556.491
Sb	120.904	5651.19	2215394633.559
Ba	134.906	5526.61	2265330490.437
Ho	164.930		
Tm	168.934	5351.68	2339377359.067
Tl	204.975	5104.08	2452862601.143
Pb	207.977	5100.71	2454483234.194
U	238.050	5062.74	2472892833.787

Daily Performance Report

Sample ID: DAILY SHARGRAVE
Sample Date/Time: Tuesday, December 07, 2010 07:34:13
Sample Description:
Sample File:
Method File: E:\elandata\Method\000daily.mth
Dataset File: e:\elandata\dataset\101207a2\DAILY SHARGRAVE.006
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: e:\elandata\Optimize\default.dac
Number of Replicates: 5
Dual Detector Mode: Dual

Summary

	Analyte	Mass	Net Intens. Mean	Net Intens. RSD
	Mg	24	39582.557	0.225
	Rh	103	227757.567	0.697
	Pb	208	271660.442	0.525
⎡>	Ba	138	316167.621	0.392
⎣	Ba++	69	0.014	2.345
⎡>	Ce	140	407072.164	0.396
⎣	CeO	156	0.033	4.799
	Bkgd	220	2.286	34.233
	Li	7	13654.536	2.149
	Be	9	3772.211	2.677
	Co	59	98556.798	0.485
	In	115	356144.422	0.429
	Tl	205	411132.717	1.234

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: Rinse 2X
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 09:12:39
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\Rinse 2X.007
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 6
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al			378104.740	ug/L	0.000
|	44	Ca			3039.059	ug/L	0.000
|	55	Mn			6805.315	ug/L	0.000
|	75	As			5300.785	ug/L	0.000
⌊>	72	Ge-1			755369.417	ug/L	0.000

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: Blank
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 09:15:34
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\Blank.008
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 5
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al			91619.527	ug/L	
|	44	Ca			3216.147	ug/L	
|	55	Mn			2860.643	ug/L	
|	75	As			5747.158	ug/L	
⌊>	72	Ge-1			790209.635	ug/L	

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: Standard 1
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 09:18:24
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\Standard 1.009
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 4
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	5100.000000	1.545	14503124.094	ug/L	91619.527
|	44	Ca	5100.000000	1.114	655965.802	ug/L	3216.147
|	55	Mn	100.000000	0.927	704528.576	ug/L	2860.643
|	75	As	100.000000	0.939	96528.759	ug/L	5747.158
⌊>	72	Ge-1			787378.408	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: ICV
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 09:25:16
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\ICV .011
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 3
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	743.588271	0.760	2197458.404	ug/L	91619.527
	44	Ca	761.281671	0.856	100857.325	ug/L	3216.147
	55	Mn	77.635827	1.570	548743.825	ug/L	2860.643
	75	As	81.596677	0.174	79997.895	ug/L	5747.158
>	72	Ge-1			789143.453	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	99.865

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: ICB
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 10:24:20
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\ICB.012
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 5
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌜	27	Al	-4.317916	10.910	82056.718	ug/L	91619.527
|	44	Ca	9.068673	2.906	4530.947	ug/L	3216.147
|	55	Mn	-0.011123	73.964	2876.317	ug/L	2860.643
|	75	As	0.410560	66.541	6334.551	ug/L	5747.158
⌊>	72	Ge-1			817166.369	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌜	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	103.411

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: LLSTD1
Sample Description: LLSTD@10X
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 10:26:59
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\LLSTD1.013
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 71
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	17.091767	5.323	148354.452	ug/L	91619.527
|	44	Ca	52.030498	3.923	10481.959	ug/L	3216.147
|	55	Mn	0.715607	3.326	8365.495	ug/L	2860.643
|	75	As	1.106023	1.564	7155.287	ug/L	5747.158
⌊>	72	Ge-1			837042.468	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	105.927

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: LLSTD2
Sample Description: LLSTD@5X
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 10:29:38
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\LLSTD2.014
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 72
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	66.838139	0.888	300021.708	ug/L	91619.527
	44	Ca	99.835983	2.159	17112.980	ug/L	3216.147
	55	Mn	1.705385	2.700	15863.749	ug/L	2860.643
	75	As	2.237674	8.661	8307.267	ug/L	5747.158
>	72	Ge-1			843216.075	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	106.708

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: ICSA
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 10:32:16
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\ICSA .015
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 2
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	97070.591213	4.249	250931943.094	ug/L	91619.527
|	44	Ca	101021.709241	3.318	11829049.317	ug/L	3216.147
|	55	Mn	6.220860	3.749	42529.990	ug/L	2860.643
|	75	As	1.129956	58.484	6176.103	ug/L	5747.158
⌊>	72	Ge-1			720572.478	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	91.188

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: ICSAB
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 10:34:53
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\ICSAB.016
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 1
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	92107.819166	0.457	253685266.589	ug/L	91619.527
	44	Ca	98661.893241	0.553	12306507.364	ug/L	3216.147
	55	Mn	100.843584	1.172	692172.885	ug/L	2860.643
	75	As	105.260552	1.795	98693.864	ug/L	5747.158
>	72	Ge-1			767178.960	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	97.085

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: Rinse
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 11:11:19
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\Rinse.017
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 6
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	71.514890	1.236	371606.938	ug/L	91619.527
|	44	Ca	7.342333	15.114	5250.167	ug/L	3216.147
|	55	Mn	0.454851	3.967	7653.936	ug/L	2860.643
|	75	As	0.511864	22.020	7842.901	ug/L	5747.158
⌊>	72	Ge-1			997449.908	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	126.226

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MARD8B
Sample Description: G0L020000-286 BLK
Batch ID: 336286
Sample Date/Time: Tuesday, December 07, 2010 11:17:36
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MARD8B.018
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 100
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	-24.567215	0.566	26882.704	ug/L	91619.527
	44	Ca	190.055925	2.965	33829.384	ug/L	3216.147
	55	Mn	-0.012615	67.446	3393.571	ug/L	2860.643
	75	As	1.390648	5.953	8590.648	ug/L	5747.158
>	72	Ge-1			967882.665	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	122.484

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MARD8C
Sample Description: G0L020000-286 LCS
Batch ID: 336286
Sample Date/Time: Tuesday, December 07, 2010 11:20:11
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MARD8C.019
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 86
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	803.970332	1.435	2711679.148	ug/L	91619.527
|	44	Ca	1005.777932	1.782	151384.609	ug/L	3216.147
|	55	Mn	178.656276	2.761	1441454.789	ug/L	2860.643
|	75	As	183.788282	1.809	198059.339	ug/L	5747.158
⌊>	72	Ge-1			903751.385	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	114.369

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MARD8L
Sample Description: G0L020000-286 LCSD
Batch ID: 336286
Sample Date/Time: Tuesday, December 07, 2010 11:22:45
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MARD8L.020
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 87
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌜	27	Al	819.309454	2.328	2508958.618	ug/L	91619.527
|	44	Ca	1005.629929	1.870	137539.916	ug/L	3216.147
|	55	Mn	182.871238	1.956	1340841.602	ug/L	2860.643
|	75	As	187.705790	1.723	183676.615	ug/L	5747.158
⌊>	72	Ge-1			821051.065	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌜	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	103.903

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCV 1
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 11:25:24
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCV 1.021
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 4
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	4704.520412	0.952	16030226.356	ug/L	91619.527
|	44	Ca	4894.713792	1.889	754014.074	ug/L	3216.147
|	55	Mn	97.820745	1.895	825341.729	ug/L	2860.643
|	75	As	101.283227	2.860	116971.805	ug/L	5747.158
⌊>	72	Ge-1			943099.344	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	119.348

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCB 1
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 11:50:24
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCB 1.025
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 5
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	-29.747050	0.051	9698.052	ug/L	91619.527
	44	Ca	-0.348102	144.847	4237.078	ug/L	3216.147
	55	Mn	-0.278783	1.724	1199.113	ug/L	2860.643
	75	As	0.713673	50.514	8548.720	ug/L	5747.158
>	72	Ge-1			1055614.610	ug/L	790209.635

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	133.587

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: BLK RECAL
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 11:50:24
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCB 1.025
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 5
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al			9698.052	ug/L	
	44	Ca			4237.078	ug/L	
	55	Mn			1199.113	ug/L	
	75	As			8548.720	ug/L	
>	72	Ge-1			1055614.610	ug/L	

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: STD1 RECAL
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 11:25:24
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCV 1.021
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 4
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	5100.000000	0.946	16030226.356	ug/L	9698.052
|	44	Ca	5100.000000	1.889	754014.074	ug/L	4237.078
|	55	Mn	100.000000	1.890	825341.729	ug/L	1199.113
|	75	As	100.000000	2.880	116971.805	ug/L	8548.720
⌊>	72	Ge-1			943099.344	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCV 2
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 12:03:37
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCV 2.026
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 4
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	5073.274368	1.561	18282350.692	ug/L	9698.052
|	44	Ca	5042.541562	2.834	854598.565	ug/L	4237.078
|	55	Mn	97.044717	2.854	918128.007	ug/L	1199.113
|	75	As	98.848480	2.433	132670.491	ug/L	8548.720
⌊>	72	Ge-1			1081582.543	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	102.460

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCB 2
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 12:06:16
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCB 2.027
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 5
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	4.491008	1.942	25689.767	ug/L	9698.052
	44	Ca	3.731183	19.687	4891.549	ug/L	4237.078
	55	Mn	0.048524	10.375	1660.550	ug/L	1199.113
	75	As	-0.051471	507.244	8549.255	ug/L	8548.720
>	72	Ge-1			1063915.140	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	100.786

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: LLSTD1
Sample Description: LLSTD@10X
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 12:08:56
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\LLSTD1.028
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 71
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	50.195283	1.026	193274.699	ug/L	9698.052
	44	Ca	52.996870	1.288	13456.548	ug/L	4237.078
	55	Mn	1.054725	1.337	11346.776	ug/L	1199.113
	75	As	0.488712	44.280	9493.869	ug/L	8548.720
>	72	Ge-1			1095550.560	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	103.783

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCV 3
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 12:46:51
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCV 3.038
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 4
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	4973.962003	0.360	16599974.270	ug/L	9698.052
	44	Ca	5032.375064	0.442	790139.601	ug/L	4237.078
	55	Mn	98.698521	0.025	865037.932	ug/L	1199.113
	75	As	101.410857	0.718	125857.110	ug/L	8548.720
>	72	Ge-1			1001215.977	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	94.847

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCB 3
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 12:49:31
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCB 3.039
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 5
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	5.332247	1.164	28157.830	ug/L	9698.052
	44	Ca	4.915177	17.786	4994.293	ug/L	4237.078
	55	Mn	0.044176	6.731	1590.199	ug/L	1199.113
	75	As	0.080988	605.092	8556.678	ug/L	8548.720
>	72	Ge-1			1044768.752	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	98.973

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCV 4
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 12:55:08
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCV 4.041
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 4
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌜	27	Al	5079.885051	0.820	17282350.653	ug/L	9698.052
¦	44	Ca	5068.248774	0.360	811193.077	ug/L	4237.078
¦	55	Mn	99.129489	0.557	885693.990	ug/L	1199.113
¦	75	As	101.122438	0.560	127965.915	ug/L	8548.720
⌊>	72	Ge-1			1020660.828	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌜	Al	27	
¦	Ca	44	
¦	Mn	55	
¦	As	75	
⌊>	Ge-1	72	96.689

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCB 4
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 12:57:47
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCB 4.042
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 5
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	5.359257	4.783	29449.997	ug/L	9698.052
	44	Ca	3.894593	15.595	5033.990	ug/L	4237.078
	55	Mn	0.047765	24.296	1690.558	ug/L	1199.113
	75	As	-0.296235	173.161	8439.853	ug/L	8548.720
>	72	Ge-1			1089732.021	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	103.232

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MAWLKB
Sample Description: G0L060000-10 BLK
Batch ID: 340010
Sample Date/Time: Tuesday, December 07, 2010 13:08:09
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MAWLKB.046
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 98
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	-1.323528	1.145	3982.913	ug/L	9698.052
	44	Ca	0.447555	237.884	3403.577	ug/L	4237.078
	55	Mn	0.171180	3.759	2194.378	ug/L	1199.113
	75	As	0.644452	49.988	7372.905	ug/L	8548.720
>	72	Ge-1			833619.116	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	78.970

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MAWLKC
Sample Description: G0L060000-10 LCS
Batch ID: 340010
Sample Date/Time: Tuesday, December 07, 2010 13:10:44
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MAWLKC.047
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 90
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	922.205124	2.577	2626029.260	ug/L	9698.052
	44	Ca	952.917880	3.121	130110.345	ug/L	4237.078
	55	Mn	185.878671	2.578	1385795.968	ug/L	1199.113
	75	As	183.882272	2.676	188642.139	ug/L	8548.720
>	72	Ge-1			852580.884	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	80.766

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MAWLKL
Sample Description: G0L060000-10 LCSD
Batch ID: 340010
Sample Date/Time: Tuesday, December 07, 2010 13:13:20
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MAWLKL.048
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 91
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	934.519374	2.122	2624092.505	ug/L	9698.052
	44	Ca	960.814623	1.430	129361.670	ug/L	4237.078
	55	Mn	189.382739	1.620	1392379.725	ug/L	1199.113
	75	As	186.888334	1.930	188972.657	ug/L	8548.720
>	72	Ge-1			840551.699	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	79.627

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MAML1
Sample Description: G0K300434-2
Batch ID: 340010
Sample Date/Time: Tuesday, December 07, 2010 13:15:54
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MAML1.049
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 17
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	192.398018	2.536	548701.223	ug/L	9698.052
|	44	Ca	1221.869828	2.625	164265.456	ug/L	4237.078
|	55	Mn	1758.087168	3.248	12969815.258	ug/L	1199.113
|	75	As	0.923495	18.738	7739.073	ug/L	8548.720
⌊>	72	Ge-1			844307.470	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	79.983

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MAML1P5
Sample Description: G0K300434-2 5X
Batch ID: 340010
Sample Date/Time: Tuesday, December 07, 2010 13:18:28
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MAML1P5.050
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 18
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	40.216849	4.654	135804.504	ug/L	9698.052
	44	Ca	265.720078	4.630	43134.703	ug/L	4237.078
	55	Mn	355.924240	4.336	2952609.736	ug/L	1199.113
	75	As	-0.250521	102.000	7408.875	ug/L	8548.720
>	72	Ge-1			949687.208	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	89.965

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MAML1Z
Sample Description: G0K300434-2 PS
Batch ID: 340010
Sample Date/Time: Tuesday, December 07, 2010 13:21:03
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MAML1Z.051
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 19
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌈	27	Al	1092.818373	2.266	2988507.331	ug/L	9698.052
|	44	Ca	2138.790498	2.823	276499.150	ug/L	4237.078
|	55	Mn	1914.447317	2.500	13704051.939	ug/L	1199.113
|	75	As	184.843338	3.210	182125.196	ug/L	8548.720
⌊>	72	Ge-1			819207.720	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌈	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	77.605

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: MAML6
Sample Description: G0K300434-3
Batch ID: 340010
Sample Date/Time: Tuesday, December 07, 2010 13:23:37
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\MAML6.052
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 20
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
⌜	27	Al	1151.092504	2.408	3188899.635	ug/L	9698.052
|	44	Ca	5394.609379	2.815	701496.151	ug/L	4237.078
|	55	Mn	110.964845	2.477	805683.507	ug/L	1199.113
|	75	As	1.173209	24.843	7846.206	ug/L	8548.720
⌊>	72	Ge-1			830012.056	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
⌜	Al	27	
|	Ca	44	
|	Mn	55	
|	As	75	
⌊>	Ge-1	72	78.628

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCV 5
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 13:26:17
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCV 5.053
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 4
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	5026.028379	0.533	16256534.086	ug/L	9698.052
	44	Ca	5028.073322	0.481	765124.403	ug/L	4237.078
	55	Mn	98.019486	0.224	832600.681	ug/L	1199.113
	75	As	99.424763	1.224	119746.827	ug/L	8548.720
>	72	Ge-1			970352.462	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	91.923

TAL-W.SACRAMENTO - Perkin Elmer Elan 6000 ICPMS M02 - Method 6020,200.8 - QUANTITATIVE ANALYSIS REPORT
SOP No. SAC-MT-0001
Analyst: SHargrave

Sample ID: CCB 5
Sample Description:
Batch ID:
Sample Date/Time: Tuesday, December 07, 2010 13:28:56
Method File: E:\elandata\Method\0006020-SH.mth
Dataset File: e:\elandata\dataset\101207a2\CCB 5.054
Tuning File: e:\elandata\Tuning\default.tun
Optimization File: E:\elandata\Optimize\default.dac
Autosampler Position: 5
Number of Replicates: 3
Dual Detector Mode: Dual
Initial Sample Quantity (mg):
Sample Prep Volume (mL):

Sample Result Summary

	Mass	Analyte	Conc. Mean	Conc. RSD	Meas. Intens. Mean	Sample Unit	Blank Intensity
	27	Al	5.007182	3.738	25941.101	ug/L	9698.052
	44	Ca	6.003347	9.073	4964.936	ug/L	4237.078
	55	Mn	0.057063	11.714	1640.212	ug/L	1199.113
	75	As	-0.270393	85.379	7805.398	ug/L	8548.720
>	72	Ge-1			1003171.648	ug/L	1055614.610

Internal Standard Recoveries

	Analyte	Mass	Int Std % Recovery
	Al	27	
	Ca	44	
	Mn	55	
	As	75	
>	Ge-1	72	95.032

Method: 6020 (SOP: SAC-MT-001) Instrument: M02 Reported: 12/07/10 14:26:16

File ID: 101207A2 **Analyst: hargraves**

#	Sample ID	Lot No.	Batch		DF	Analyzed Date	Comment	Q
1	Rinse 2X				2.0	12/07/10 09:12		☐
2	Blank				1.0	12/07/10 09:15		☐
3	Standard1				1.0	12/07/10 09:18		☐
4	ICV				1.0	12/07/10 09:20		☐
5	ICV				1.0	12/07/10 09:25		☐
6	ICB				1.0	12/07/10 10:24		☐
7	LLSTD1				1.0	12/07/10 10:26		☐
8	LLSTD2				1.0	12/07/10 10:29		☐
9	ICSA				1.0	12/07/10 10:32		☐
10	ICSAB				1.0	12/07/10 10:34		☐
11	Rinse				1.0	12/07/10 11:11		☐
12	MARD8B	G0L020000	0336286	2A	1.0	12/07/10 11:17		☐
13	MARD8C	G0L020000	0336286	2A	1.0	12/07/10 11:20		☐
14	MARD8L	G0L020000	0336286	2A	1.0	12/07/10 11:22		☐
15	CCV 1				1.0	12/07/10 11:25		☐
16	CCB 1				1.0	12/07/10 11:50		☐
19	CCV 2				1.0	12/07/10 12:03		☐
20	CCB 2				1.0	12/07/10 12:06		☐
21	LLSTD1				1.0	12/07/10 12:08		☐
22	MAPEVB	G0L010000	0335251	2A	1.0	12/07/10 12:23		☐
23	MAPEVC	G0L010000	0335251	2A	1.0	12/07/10 12:26		☐
24	MAPE7L	G0L010000	0335253	2A	1.0	12/07/10 12:28		☐
25	MAA80	G0K190601-3	0335251	2A	1.0	12/07/10 12:31		☐
26	MAA80P5	G0K190601	0335251		5.0	12/07/10 12:34		☐
27	MAA80Z	G0K190601-3	0335251		1.0	12/07/10 12:36		☐
28	MAA81	G0K190601-4	0335251	2A	1.0	12/07/10 12:39		☐
29	MAKDV	G0K240587-1	0335251	2A	1.0	12/07/10 12:41		☐
30	MAKD2	G0K240587-2	0335251	2A	1.0	12/07/10 12:44		☐
31	CCV 3				1.0	12/07/10 12:46		☐
32	CCB 3				1.0	12/07/10 12:49		☐
33	CCV 4				1.0	12/07/10 12:55		☐
34	CCB 4				1.0	12/07/10 12:57		☐
35	MARD8B	G0L020000	0336286	2A	1.0	12/07/10 13:00		☐
36	MARD8C	G0L020000	0336286	2A	1.0	12/07/10 13:02		☐
37	MARD8L	G0L020000	0336286	2A	1.0	12/07/10 13:05		☐
38	MAWLKB	G0L060000	0340010	2A	1.0	12/07/10 13:08		☐
39	MAWLKC	G0L060000	0340010	2A	1.0	12/07/10 13:10		☐
40	MAWLKL	G0L060000	0340010	2A	1.0	12/07/10 13:13		☐
41	MAML1	G0K300434-2	0340010	2A	1.0	12/07/10 13:15		☐
42	MAML1P5	G0K300434	0340010		5.0	12/07/10 13:18		☐
43	MAML1Z	G0K300434-2	0340010		1.0	12/07/10 13:21		☐
44	MAML6	G0K300434-3	0340010	2A	1.0	12/07/10 13:23		☐
45	CCV 5				1.0	12/07/10 13:26		☐
46	CCB 5				1.0	12/07/10 13:28		☐
47	MA0J7B	G0L070000	0341211	2A	1.0	12/07/10 13:39		☐
48	MA0J7C	G0L070000	0341211	2A	1.0	12/07/10 13:42		☐

Method: 6020 (SOP: SAC-MT-001) Instrument: M02 Reported: 12/07/10 14:26:16

File ID: 101207A2 Analyst: hargraves

#	Sample ID	Lot No.	Batch		DF	Analyzed Date	Comment	Q
49	MA0J7L	G0L070000	0341211	2A	1.0	12/07/10 13:44		☐
50	MAQQ1	G0L020446-3	0341211	2A	1.0	12/07/10 13:47		☐
51	MAQQ1P5	G0L020446	0341211		5.0	12/07/10 13:49		☐
52	MAQQ1Z	G0L020446-3	0341211		1.0	12/07/10 13:52		☐
53	MAQQ4	G0L020446-4	0341211	2A	1.0	12/07/10 13:55		☐
54	MAQRA	G0L020446-7	0341211	2A	1.0	12/07/10 13:57		☐
55	MAQRH	G0L020446-10	0341211	2A	1.0	12/07/10 14:00		☐
56	CCV 6				1.0	12/07/10 14:02		☐
57	CCB 6				1.0	12/07/10 14:05		☐

Method: 6020 (SOP: SAC-MT-001) M02 (M02) Reported: 12/07/10 14:26:16

File ID: 101207A2 **Analyst: hargraves**

#	Sample ID	Analyzed Date	Germanium	Q
1	Rinse 2X	12/07/10 09:12	95.6	☐
2	Blank	12/07/10 09:15	100.0	☑
3	Standard1	12/07/10 09:18	99.6	☑
4	ICV	12/07/10 09:20	99.1	☑
5	ICV	12/07/10 09:25	99.9	☑
6	ICB	12/07/10 10:24	103.4	☑
7	LLSTD1	12/07/10 10:26	105.9	☑
8	LLSTD2	12/07/10 10:29	106.7	☑
9	ICSA	12/07/10 10:32	91.2	☑
10	ICSAB	12/07/10 10:34	97.1	☑
11	Rinse	12/07/10 11:11	126.2	☑
12	MARD8B	12/07/10 11:17	122.5	☑
13	MARD8C	12/07/10 11:20	114.4	☑
14	MARD8L	12/07/10 11:22	103.9	☑
15	CCV 1	12/07/10 11:25	119.3	☑
16	CCB 1	12/07/10 11:50	133.6	☐
19	CCV 2	12/07/10 12:03	102.5	☑
20	CCB 2	12/07/10 12:06	100.8	☑
21	LLSTD1	12/07/10 12:08	103.8	☑
22	MAPEVB	12/07/10 12:23	94.3	☑
23	MAPEVC	12/07/10 12:26	87.2	☑
24	MAPE7L	12/07/10 12:28	81.7	☑
25	MAA80	12/07/10 12:31	79.5	☑
26	MAA80P5	12/07/10 12:34	96.1	☐
27	MAA80Z	12/07/10 12:36	82.6	☑
28	MAA81	12/07/10 12:39	80.7	☑
29	MAKDV	12/07/10 12:41	85.3	☑
30	MAKD2	12/07/10 12:44	86.8	☑
31	CCV 3	12/07/10 12:46	94.8	☑
32	CCB 3	12/07/10 12:49	99.0	☑
33	CCV 4	12/07/10 12:55	96.7	☑
34	CCB 4	12/07/10 12:57	103.2	☑
35	MARD8B	12/07/10 13:00	97.4	☑
36	MARD8C	12/07/10 13:02	86.2	☑
37	MARD8L	12/07/10 13:05	81.1	☑
38	MAWLKB	12/07/10 13:08	79.0	☑
39	MAWLKC	12/07/10 13:10	80.8	☑
40	MAWLKL	12/07/10 13:13	79.6	☑
41	MAML1	12/07/10 13:15	80.0	☑
42	MAML1P5	12/07/10 13:18	90.0	☐
43	MAML1Z	12/07/10 13:21	77.6	☑
44	MAML6	12/07/10 13:23	78.6	☑
45	CCV 5	12/07/10 13:26	91.9	☑
46	CCB 5	12/07/10 13:28	95.0	☑
47	MA0J7B	12/07/10 13:39	91.8	☑
48	MA0J7C	12/07/10 13:42	82.1	☑

TAL West Sac

INTERNAL STANDARD SUMMARY

Method: 6020 (SOP: SAC-MT-001) M02 (M02) Reported: 12/07/10 14:26:16

File ID: 101207A2 **Analyst: hargraves**

Germanium

#	Sample ID	Analyzed Date		Q
49	MA0J7L	12/07/10 13:44	80.6	☑
50	MAQQ1	12/07/10 13:47	79.0	☑
51	MAQQ1P5	12/07/10 13:49	86.1	☐
52	MAQQ1Z	12/07/10 13:52	79.3	☑
53	MAQQ4	12/07/10 13:55	77.4	☑
54	MAQRA	12/07/10 13:57	80.6	☑
55	MAQRH	12/07/10 14:00	85.8	☑
56	CCV 6	12/07/10 14:02	93.4	☑
57	CCB 6	12/07/10 14:05	97.2	☑

TAL West Sac

CALIBRATION CHECK SUMMARY

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Method: 6020 **Instrument: M02** **Batch: 101207A2**

Sample ID	Type	File - Sequence	Analyzed Date	Q
ICV	ICV	101207A2, 4	12/07/2010 09:20:59	☐
ICV	ICV	101207A2, 5	12/07/2010 09:25:16	☐
ICB	ICB	101207A2, 6	12/07/2010 10:24:20	☐
ICSA	ICSA	101207A2, 9	12/07/2010 10:32:16	☐
ICSAB	ICSAB	101207A2, 10	12/07/2010 10:34:53	☐
CCV 1	CCV	101207A2, 15	12/07/2010 11:25:24	☐
CCB 1	CCB	101207A2, 16	12/07/2010 11:50:24	☐
CCV 2	CCV	101207A2, 19	12/07/2010 12:03:37	☐
CCB 2	CCB	101207A2, 20	12/07/2010 12:06:16	☐
CCV 3	CCV	101207A2, 31	12/07/2010 12:46:51	☐
CCB 3	CCB	101207A2, 32	12/07/2010 12:49:31	☐
CCV 4	CCV	101207A2, 33	12/07/2010 12:55:08	☐
CCB 4	CCB	101207A2, 34	12/07/2010 12:57:47	☐
CCV 5	CCV	101207A2, 45	12/07/2010 13:26:17	☐
CCB 5	CCB	101207A2, 46	12/07/2010 13:28:56	☐
CCV 6	CCV	101207A2, 56	12/07/2010 14:02:57	☐
CCB 6	CCB	101207A2, 57	12/07/2010 14:05:37	☐

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **ICV (ICV)** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 4 Method 6020_
Acquired: 12/07/2010 09:20:59 M02
Calibrated: 12/07/2010 09:15:34 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	Aluminum	27	2727202	938.26	800.00	117	
7439-96-5	Manganese	55	564887	80.568	80.000	101	
7440-38-2	Arsenic	75	79894	82.186	80.000	103	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	783241		☑

Reviewed by: Date:

TAL West Sac

CALIBRATION REPORT

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **ICV (ICV)** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 5 Method 6020_
Acquired: 12/07/2010 09:25:16 M02
Calibrated: 12/07/2010 09:15:34 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	**Aluminum**	27	2197458	743.59	800.00	92.9	
7439-96-5	**Manganese**	55	548744	77.636	80.000	97.0	
7440-38-2	**Arsenic**	75	79998	81.597	80.000	102	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	789143		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **ICB** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 6 Method 6020_
Acquired: 12/07/2010 10:24:20 M02
Calibrated: 12/07/2010 09:15:34 Units: ug/L

CASN	Analyte Name	M/S	Area	Amount	RL	MDL	%RSD	Q
7429-90-5	Aluminum	27	82057	-4.3179	50.0	2.1	0.0	
7439-96-5	Manganese	55	2876	-0.01112	1.0	0.083	0.0	
7440-38-2	Arsenic	75	6335	0.41056	2.0	0.50	0.0	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	817166		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **ICSA** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 9 Method 6020_
Acquired: 12/07/2010 10:32:16 M02
Calibrated: 12/07/2010 10:26:59 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	Aluminum	27	50931943	97071	100000	97.1	☑
7439-96-5	Manganese	55	42530	6.2209		*	
7440-38-2	Arsenic	75	6176	1.1300		*	☑

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	720572		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **ICSAB** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 10 Method 6020_
Acquired: 12/07/2010 10:34:53 M02
Calibrated: 12/07/2010 10:26:59 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	Aluminum	27	53685267	92108	100100	92.0	☑
7439-96-5	Manganese	55	692173	100.84	100.00	101	☑
7440-38-2	Arsenic	75	98694	105.26	100.00	105	☑

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	767179		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCV 1 (CCV)** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 15 Method 6020_
Acquired: 12/07/2010 11:25:24 M02
Calibrated: 12/07/2010 10:26:59 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	**Aluminum**	27	16030226	4704.5	5100.0	92.2	
7439-96-5	**Manganese**	55	825342	97.821	100.00	97.8	
7440-38-2	**Arsenic**	75	116972	101.28	100.00	101	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	**Germanium**	72	943099		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCB 1** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 16 Method 6020_
Acquired: 12/07/2010 11:50:24 M02
Calibrated: 12/07/2010 11:25:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Amount	RL	MDL	%RSD	Q
7429-90-5	Aluminum	27	9698	-29.747	50.0	2.1	0.0	
7439-96-5	Manganese	55	1199	-0.27878	1.0	0.083	0.0	
7440-38-2	Arsenic	75	8549	0.71367	2.0	0.50	0.0	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	1055615		☐

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCV 2 (CCV)** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 19 Method 6020_
Acquired: 12/07/2010 12:03:37 M02
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	Aluminum	27	18282351	5073.3	5100.0	99.5	
7439-96-5	Manganese	55	918128	97.045	100.00	97.0	
7440-38-2	Arsenic	75	132670	98.848	100.00	98.8	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	1081583		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCB 2** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 20 Method 6020_
Acquired: 12/07/2010 12:06:16 M02
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Amount	RL	MDL	%RSD	Q
7429-90-5	Aluminum	27	25690	4.4910	50.0	2.1	0.0	
7439-96-5	Manganese	55	1661	0.04852	1.0	0.083	0.0	
7440-38-2	Arsenic	75	8549	-0.05147	2.0	0.50	0.0	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	1063915		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCV 3 (CCV)** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 31 Method 6020_
Acquired: 12/07/2010 12:46:51 M02
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	Aluminum	27	16599974	4974.0	5100.0	97.5	
7439-96-5	Manganese	55	865038	98.699	100.00	98.7	
7440-38-2	Arsenic	75	125857	101.41	100.00	101	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	1001216		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCB 3** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 32 Method 6020_
Acquired: 12/07/2010 12:49:31 M02
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Amount	RL	MDL	%RSD	Q
7429-90-5	Aluminum	27	28158	5.3322	50.0	2.1	0.0	
7439-96-5	Manganese	55	1590	0.04418	1.0	0.083	0.0	
7440-38-2	Arsenic	75	8557	0.08099	2.0	0.50	0.0	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	1044769		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCV 4 (CCV)** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 33 Method 6020_
Acquired: 12/07/2010 12:55:08 M02
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	**Aluminum**	27	17282351	5079.9	5100.0	99.6	
7439-96-5	**Manganese**	55	885694	99.129	100.00	99.1	
7440-38-2	**Arsenic**	75	127966	101.12	100.00	101	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	**Germanium**	72	1020661		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCB 4** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 34 Method 6020_
Acquired: 12/07/2010 12:57:47 M02
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Amount	RL	MDL	%RSD	Q
7429-90-5	Aluminum	27	29450	5.3593	50.0	2.1	0.0	
7439-96-5	Manganese	55	1691	0.04777	1.0	0.083	0.0	
7440-38-2	Arsenic	75	8440	-0.29624	2.0	0.50	0.0	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	1089732		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCV 5 (CCV)** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 45 Method 6020_
Acquired: 12/07/2010 13:26:17 M02
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Found	True	%R	Q
7429-90-5	Aluminum	27	16256534	5026.0	5100.0	98.5	
7439-96-5	Manganese	55	832601	98.019	100.00	98.0	
7440-38-2	Arsenic	75	119747	99.425	100.00	99.4	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	970352		☑

Reviewed by: Date:

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:26:55

Department: 120 (Metals) Source: MetEdit

Sample: **CCB 5** Mult: **1.00** Dilf: **1.00** Divs: **1.000**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 #46 Method 6020_
Acquired: 12/07/2010 13:28:56 M02
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Amount	RL	MDL	%RSD	Q
7429-90-5	**Aluminum**	27	25941	5.0072	50.0	2.1	0.0	
7439-96-5	**Manganese**	55	1640	0.05706	1.0	0.083	0.0	
7440-38-2	**Arsenic**	75	7805	-0.27039	2.0	0.50	0.0	

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	**Germanium**	72	1003172		☑

Reviewed by: Date:

TAL West Sac

SERIAL DILUTION

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:27:37

Department: 120 (Metals) Source: MetEdit

Sample: **MAML1P5** Serial Dilution: **5.00** Sample Dilution: **1.00**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 42 Method 6020_
Acquired: 12/07/2010 13:18:28 M02 Matrix: AIR
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Dilution	Sample	%Diff.	MDL	Flag	Q
7429-90-5	Aluminum	27	135805	201.08	192.40	4.51		*	
7439-96-5	Manganese	55	2952610	1779.6	1758.1	1.22	0.14	1.2	☑
7440-38-2	Arsenic	75	7409	-1.2526	0.92349	236	0.41	NC	☑

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	949687		☐

* Analyte not requested for this batch, no MDL

NC : Serial dilution concentration < 100 X MDL

E : Difference greater than Limit (10%)

Reviewed by: Date:

TAL West Sac

SAMPLE SPIKE

Method: 6020 (SOP: SAC-MT-001) M02 Reported: 12/07/10 14:27:46

Department: 120 (Metals) Source: MetEdit

Sample: **MAML1Z** Spike Dilution: **1.00** Sample Dilution: **1.00**

Instrument: **ICPMS M02** Channel 262
File: 101207A2 # 43 Method 6020_
Acquired: 12/07/2010 13:21:03 M02 Matrix: AIR
Calibrated: 12/07/2010 11:50:24 Units: ug/L

CASN	Analyte Name	M/S	Area	Amount	Sample	%Rec.	Spike	Flag	Q
7429-90-5	Aluminum	27	2988507	1092.8	192.40	90.0	1000		☑
7439-96-5	Manganese	55	13704052	1914.4	1758.1	78.2	200	*	☑
7440-38-2	Arsenic	75	182125	184.84	0.92349	92.0	200		☑

CASN	ISTD Name	M/S	Area	Amount	Q
7440-56-4	Germanium	72	819208		☑

Reviewed by: Date:

Sample Preparation Log

TestAmerica - West Sacramento
Metals - Air Toxics - Preparation Log

Date: 6-Dec-10 **Analyst:** FF **Matrix:** AIR

Fraction: Filter **SOP:** WS-IP-0001 **Method:** ICPMS

LOT ID		Workorder		Volume Received	Volume Removed	Initial Prep Volume	Final Prep Volume	Batch	Prep Factor
G0L060000	10	MAWLKB	2A	NA	NA	NA	100 mL	340010	1.2
G0L060000	10	MAWLKC	2A	NA	NA	NA	100 mL	340010	1.2
G0L060000	10	MAWLKL	2A	NA	NA	NA	100 mL	340010	1.2
G0K300434	2	MAML1	2A	9 inches	0.75 inches	0.75 inches	100 mL	340010	1.2
G0K300434	3	MAML6	2A	9 inches	0.75 inches	0.75 inches	100 mL	340010	1.2

For the cassette filter digest the whole filter is used.
For 1" filter: factor = 9 (9/1).
For 0.75" filter factor = 12 (9/0.75).

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

West Sacramento

Metals Spiking Documentation Form

Lot #(s): G0K300434

Batch Number: 0340010
EPA Analytical Method ID: 6020
Spiked Date: 12/06/10

MS Sample(s): N/A
EPA Prep Method ID: 3.0, 3.1
Hot Plate / Microwave ID: MEB T II

Analyst Initial/Date: FF 12/6/10
Witness Initial/Date: 12/06/10 NM
Hot Plate Temp Initial: / Final:

Correct Folder ID Witness: N/A
Digestion Cup Lot #: 101029
Filter Paper Lot #: 390428
Thermometer ID: B709
Fin Vol Cup Lot: 100811

Check If Used	Bottle Name	Elements	Stock Concentration (mg/L)	Tracking Number	LCS/LCSD Volume Spiked	MS/SD Volume Spiked	Expiration Date
	ICP Part 1 5% HNO_3	Ca, Mg	5,000				
		Al, As, Ba, Se, Sn, Tl	200				
		Fe,Mo,Ti	100				
		Sb;Co;Pb;Mn;Ni;V;Zn	50				
		Cu	25				
		Cr	20				
		,Be,Cd	5				
		Ag	5.0				
	ICP Part 2 2% HNO_3	K,Na	5,000				
		P,S	1,000				
		B,Li,Sr	100				
	Si H20/Tr HF	Si	1,000			TP 12/05/10	
✓	TACA-1 5% HNO_3	Al,K,Mg,Ca,Na,Fe,P,B	500	3189-65	200uL	N/A	8/31/11
		As,Be,Cd,Cr,Co,Cu,Pb, Mn,Ni,Se,U,V,Zn,Ba,Li Sr	100				
		Ag,Tl	25				
✓	TACA-2 5% HNO_3	Mo, Sb, Sn, Ti	100	3189-66	200 uL	N/A	8/31/11
	Misc. Elements					TP 12/05/10	

Prep Reagents:

Check If Used	Reagent	Supplier	Lot Number	Check If Used	Reagent	Supplier	Lot Number
	70% HNO_3	Mallinckrodt			30% H_2O_2	Mallinckrodt	
	37% HCl	Mallinckrodt			49% HF	Fisher	
✓	3M HNO_3	In-House	4028-32-1		1:1 HCl	In-House	TP 12/05/10

ICP matrix spike and LCS: For final volumes of 100ml, add 1mL from bottles ICP Part 1, ICP Part 2. Add 1ml of Silica (Si) when requested.
ICPMS matrix spike and LCS: For final volumes of 100ml, add 0.2 mL each of TACA-1 and TACA-2.
Amount to spike is as listed above for final volumes of 100ml. If a different final volume is used, increase or decrease the amount you spike proportionally.

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West Sacramento

Preparation Data Review Checklist

Prep Batch(es) 0340010 Test: 6020

Prep Date: 12/06/10 Holding Times: 5/24/11 NCM: Y (N)

A. Spike Witness/Batch setup	Spike Witness	Reviewer
1. **Holding times checked?** NCMs filed as appropriate	✓	✓
2. QAS checked for QC instructions (LCS, LCSD, MS,MSD, etc)	✓	✓
3. Amount of samples in hood match amount of samples on bench sheet. Sample IDS match.	✓	NA
4. Worksheets have been checked for required spiking compounds	✓	✓
5. Spiking volumes are correctly documented	✓	✓
6. Std ID numbers on spike labels match numbers on bench sheet	✓	NA
7. Expiration dates have been checked	✓	✓
8. Calibration expiration dates on pipettors have been checked	✓	NA
9. Spiker and spike witness have signed and dated bench sheet	✓	✓
B. Weights and Volumes		
1. Recorded weights are in anticipated range	NA	NA
2. Balance upload or raw data for weights is included	NA	✓
3. Weights and volumes have been transcribed correctly to LIMS.	NA	✓
4. Weights are not targeted to meet exact weights.	NA	NA
5. Each weight or volume measurement is a unique record (no dittos or line downs)	NA	✓
C. Standards and Reagents		
1. Lot numbers for all reagents, including clean up stages, are recorded.	NA	✓
2. Are dates and analysts for cleanups recorded?	NA	NA
3. Are correct IDs used for standards? Are expiration dates to day/month/year, when listed?	NA	✓
D. Documentation		
1. Are all nonconformances documented appropriately?	NA	NA
2. QuantIMs entry correct, including dates and times.	NA	✓
3. Are all fields completed?	NA	✓

Spike witness: NM Date: 12/06/10

2nd Level Reviewer: SH Date: 12/6/10

Comments:

AIR, TSP-
Total Suspended
Particulates

Raw Data Package

PARTICULATE ANALYSIS

LEVEL 1 & 2 REVIEW CHECKLIST

LAB NUMBERS: G0K300434 (2-3) Batch #: 0341292

ANALYSIS: (circle) TSP/PM10 or METHOD 5

DATE: 12/07/10 **ANALYST:** JZ

LEVEL 1 ANALYSIS REVIEW	YES	NO	NA
1. Samples are in good condition.	✓		
2. Sample filter number matches the folder or petri ID number.	✓		
3. Desiccator temperature and % humidity criteria in control.	✓		
4. Balance calibration criteria met.	✓		
5. Beginning and ending calibration sample bracket weights are in calibration.	✓		
6. Samples reached stable weight.	✓		
7. Samples exceeded 5 consecutive final weighings.		✓	
LEVEL 1 DATA REVIEW			
1. Benchsheet is complete.	✓		
2. QAS or QAPP consulted and followed for client specifics.	✓		
3. Data entered in properly.	✓		
4. Copy of spreadsheet or logbook raw data entry attached to data package.	✓		
5. Analyst observations, HTV's, Anomalies properly documented and attached to data package.			✓

Completed By & Date: JZ 12/07/10

LEVEL 2 REVIEW:			
1. Level 1 checklist complete and verified.	✓		
2. Deviations, Anomalies, Holding times checked and approved.	✓		
3. Reanalysis documented and chemist notified.			✓
4. Client specific criteria met.	✓		
5. Data entry checked and released in Quantims.	✓		
6. Indication on benchsheet or spreadsheet on review and released (dated & signed).	✓		

Completed By & Date: SV 12/7/10

Comments: Dessicator 3A

RQC050 TestAmerica Laboratories, Inc. Run Date: 12/07/10
WET CHEM BATCHSHEET Time: 14:18:35

TestAmerica West Sacramen

PRODUCTION FIGURES - WET CHEM

TOTAL NUMBER	SAMPLE NUMBER	QC	RE-RUN MATRIX	RE-RUN OTHER	MISC NUMBER	TOTAL HOURS	EXPANDED DELIVERABLE

METHOD: AO Particulates in Air, Suspended "TSP HiVol" (APP B)
QC BATCH #: **0341292** INITIALS: DATA ENTRY:
PREP DATE: 11/30/10 12:30 PREP JZ INITIALS JZ
COMP DATE: 12/02/10 17:35 ANAL JZ DATE 12/07/10
USER: PHOMSOPT

Work Order	Lab Number	Structured Analysis	Exp. Del.	Analysis Date	Sample ID:
MAML1-1-AA	G-0K300434-002	XX S 88 AO 3W	M	12/07/10	UW-11242010B
MAML6-1-AA	G-0K300434-003	XX S 88 AO 3W	M	12/07/10	DW-11242010B

Control Limits

TestAmerica

WEST SACRAMENTO

AIR TOXICS GRAVIMETRIC ANALYSES

Lab ID	Filter ID	Initial Weight (g) date/time initials	Initial Weight (g) date/time initials	Final Weight (g) date/time initials	Final Weight (g) date/time initials	Final Weight (g) date/time initials	Final Weight (g) date/time initials	Final Weight (g) date/time initials	Final Weight (g) date/time initials	Wt of Particulate (g)	Initial Wts Stablity Check
	5g wt	4.9999 110110skv1123	4.9998 110110skv1918	4.9996 120210jz1034	4.9998 120210jz1728						
	5g wt	4.9999 110110skv1123	4.9998 110110skv1918	4.9998 120610jz1038	5.0000 120610jz1802						
MAK04 112910jz1010	tron111010-66	4.5511 110110skv1123	4.5507 110110skv1919	4.5630 113010jz1240	4.5629 120110jz1056					0.0122	
	tron111010-67	4.5565 110110skv1124	4.5564 110110skv1919	4.5996 120610jz1040	4.5999 120610jz1808					0.0435	
	tron111010-68	4.5551 110110skv1124	4.5546 110110skv1919	4.5859 120610jz1042	4.5861 120610jz1817					0.0315	
	tron111010-69	4.5467 110110skv1125	4.5464 110110skv1920	4.5832 120610jz1045	4.5834 120610jz1812					0.0370	
	tron111010-70	4.5572 110110skv1125	4.5571 110110skv1920	4.5841 120610jz1047	4.5840 120610jz1819					0.0269	
MAML6 113010jz1230	tron111010-71	4.5539 110110skv1126	4.5539 110110skv1921	4.6496 120210jz1036	4.6498 120210jz1732					0.0959	✓
MAML1 113010jz1230	tron111010-72	4.5504 110110skv1126	4.5500 110110skv1921	4.5713 120210jz1037	4.5714 120210jz1733					0.0214	✓
MAK09 112910jz1010	tron111010-73	4.5577 110110skv1127	4.5572 110110skv1922	4.6235 113010jz1242	4.6227 120110jz1057	4.6222 120210jz1023				0.0650	
	tron111010-74	4.5735 110110skv1127	4.5730 110110skv1922							NC	
MAK08 112910jz1010	tron111010-75	4.4545 110110skv1128	4.4543 110110skv1922	4.5166 113010jz1243	4.5163 120110jz1058					0.0620	
	5g wt	5.0001 110110skv1128	5.0000 110110skv1923	4.9995 113010jz1245	4.9996 120110jz1100	4.9996 120210jz1031				-0.0004	
	5g wt	5.0001 110110skv1128	5.0000 110110skv1923	4.9996 120210jz1039	4.9998 120210jz1735					-0.0002	

PDE115

TestAmerica Laboratories, Inc.
Inorganics Batch Review
QC Batch **0341292**

Date 12/07/2010
Time 13:19:45

Method Code:**AO Particulates in Air, Suspended "TSP HiVol" (APP B)**
Analyst:**Thep Phomsopha**

Work Order	Result	Units	LDL/Dil	Prep. - Anal.	Total Solids	PSRL Flag	R/R	Rounded Result	Output LDL	Dil.
MAML1-1-AA	0.0214	g	0.0005	11/30-12/07/10	.00	N		0.0214	0.0005	1.00
MAML6-1-AA	0.0959	g	0.0005	11/30-12/07/10	.00	N		0.0959	0.0005	1.00

Notes:

PRODUCTION TOTALS

TEST	TOTAL #	SAMPLE #	QC #	MATRIX #	OTHER #	MISC #	HOURS
	0	0	0	0	0	0	.0

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica West Sacramento
Balance Calibration Check Log

WEIGHT #1				WEIGHT #2				DATE	INIT.	WEIGHT ID	P/F *1
Working WT Denomination (g)	OBSERVED WEIGHT (g)	Acceptance limits*2 Lower (g)	Upper (g)	Working WT Denomination (g)	OBSERVED WEIGHT (g)	Acceptance limits*2 Lower (g)	Upper (g)				
0.2000	0.1996	0.1995	0.2005	10.000	10.0002	~~10.1~~ ECF 10/4/10 9.9000	10.100	10/21/10	ECF	QA-011	P
0.2000	0.2003	0.1995	0.2005	10.000	10.0001	9.9000	10.100	10/22/10	ECF	QA-011	P
0.2000	0.2001	0.1995	0.2005	10.0	10.0005	9.9000	10.0100	10/24/10	ECF	QA-011	P
0.2000	0.2002	0.1995	0.2005	10.000	10.0011	9.9000	10.0100	10/25/10	ECF	QA-011	P
0.2000	0.2000	0.1995	0.2005	10.0	10.0004	9.9000	10.0100	10/26/10	ECF	QA-011	P
0.2000	0.2001	0.1995	0.2005	10.0	9.9997	9.9000	10.0100	10/27/10	ECF	QA-011	P
0.2000	0.1998	0.1995	0.2005	10.0	10.0000	9.9000	10.0100	10/28/10	ECF	QA-011	P
0.2000	0.1999	0.1995	0.2005	10.0	9.9999	9.9000	10.0100	10/29/10	ECF	QA-011	P
0.2000	0.1996	0.1995	0.2005	10.0	9.9993	9.9000	10.0100	11/1/10	ECF	QA-011	P
0.2000	0.1996	0.1995	0.2005	10.0	9.9999	9.9000	10.0100	11/2/10	ECF	QA-011	P
0.2000	0.2001	0.1995	0.2005	10.0	10.0004	9.9000	10.0100	11/3/10	ECF	QA-011	P
0.2000	0.2000	0.1995	0.2005	10.0	10.0004	9.9000	10.0100	11/4/10	ECF	QA-011	P
0.2000	0.1999	0.1995	0.2005	10.0	10.0004	9.9000	10.0100	11/5/10	ECF	QA-011	P
0.2000	0.1999	0.1995	0.2005	10.0	10.0000	9.9000	10.0100	11/8/10	ECF	QA-011	P

1 P= Pass, F= Fail. The observed weight must be with in the listed tolerances in order to pass. If calibration check values fall outside acceptance limits, the balance is considered to be out of calibration.
a) Do not move or use the balance
b) Attach a sign instructing others not to use the balance (see front of logbook).
c) Notify the QA department.

*2 Balance Tolerances (grams):

Denomination	Range	Denomination	Range
0.2000	0.1995 – 0.2005	10	9.9000 – 10.100
0.5000	0.4995 – 0.5005	20	19.8000 – 20.200
1	0.9900 – 1.0100	50	49.5000 – 50.500
2	1.9800 – 2.0200	100	99.0000 – 101.000
5	4.9500 – 5.0500		

Calibration range is (+/-) 1% for top loading balances. The above tolerances have been rounded to meet balance read out capability.

*3 When performing Method 1664A, the following Class 1 weights and tolerances must be used (in grams).

Denomination	Range
0.0020	0.0018 – 0.0022
1	0.9950 – 1.0050
Calibration range is (±) 10% for 2 mg weight and (±) 0.5% for 1 g weight. The above tolerances have been modified to meet balance read out capability.	

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica West Sacramento
Balance Calibration Check Log

WEIGHT #1				WEIGHT #2				DATE	INIT.	WEIGHT ID	P/F *1
Working WT Denomination (g)	OBSERVED WEIGHT (g)	Acceptance limits*2 Lower (g)	Upper (g)	Working WT Denomination (g)	OBSERVED WEIGHT (g)	Acceptance limits*2 Lower (g)	Upper (g)				
0.2000	0.2001	0.1995	0.2005	10.0000	9.9994	9.9000	10.1000	11/30/10	JZ	QA 011	P
0.2000	0.2002	0.1995	0.2005	10.0000	9.9995	9.9000	10.1000	12/1/10	JZ	QA 011	P
0.2000	0.2000	0.1995	0.2005	10.0000	9.9994	9.9000	10.1000	12/2/10	BN	QA-11	P
0.2000	0.2000	0.1995	0.2005	10.0000	9.9997	9.9000	10.1000	12/3/10	SV	QA-11	P
0.2000	0.2002	0.1995	0.2005	10.0000	10.0003	9.9000	10.1000	12/6/10	JZ	QA 011	P

1 P= Pass, F= Fail. The observed weight must be with in the listed tolerances in order to pass. If calibration check values fall outside acceptance limits, the balance is considered to be out of calibration
a) Do not move or use the balance
b) Attach a sign instructing others not to use the balance (see front of logbook).
c) Notify the QA department.

*2 Balance Tolerances (grams):

Denomination	Range	Denomination	Range
0.2000	0.1995 – 0.2005	10	9.9000 – 10.100
0.5000	0.4995 – 0.5005	20	19.8000 – 20.200
1	0.9900 – 1.0100	50	49.5000 – 50.500
2	1.9800 – 2.0200	100	99.0000 – 101.000
5	4.9500 – 5.0500		

Calibration range is (+/-) 1% for top loading balances. The above tolerances have been rounded to meet balance read out capability.

*3 When performing Method 1664A, the following Class 1 weights and tolerances must be used (in grams).

Denomination	Range
0.0020	0.0018 – 0.0022
1	0.9950 – 1.0050
Calibration range is (±) 10% for 2 mg weight and (±) 0.5% for 1 g weight. The above tolerances have been modified to meet balance read out capability.	

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica West Sacramento
Air Toxics
Desiccator Humidity/Temperature Logbook

Desiccator #		1			2			3			4			5			6			7			Amb	
Date	Init	T	RH	FN	T	RH	FN	T	RH	FN	T	RH	FN	T	RH	FN	T	RH	FN	T	RH	FN	T	RH
10/27/10	ECf	70	33	—	71	28	—	72	35	—	71	37	—	71	38	—	73	35	—	72	31	—	73	45
10/28/10	ECf	70	33	—	71	28	—	72	36	—	72	41	(1)	71	42	(1)	73	35	—	72	32		73	56
10/29/10	ECf	69	33	—	69	28	—	70	37	—	69	28	—	70	28	—	72	35	—	72	31	—	72	43
10/26/10	ECf	68	33	—	68	28	—	70	36	—	68	28	—	69	28	—	70	35	—	70	32	—	70	34
10/27/10	ECf	67	33	—	68	28	—	69	35	—	68	28	—	68	28	—	70	34	—	70	31	—	70	30
10/28/10	ECf	65	34	—	66	29	—	66	36	—	65	29	—	66	29	—	68	34	—	66	31	—	68	40
10/29/10	ECf	65	33	—	66	29	—	67	37	—	65	29	—	66	29	—	68	34	—	66	31	—	68	48
10/1/10	ECf	65	34	—	65	30	—	66	40	—	65	29	—	66	29	—	68	35	—	66	32	—	68	49
11/2/10	ECf	65	34	—	66	32	—	67	41		65	30	—	66	29	—	68	35	—	68	32	—	68	50
11/3/10	ECf	65	34	—	66	34	—	66	42		65	29	—	66	29	—	68	36	—	66	33	—	68	51
11/4/10	ECf	68	33	—	69	36	(1) @ 17:30 ECf 11/4/10	70	42	(1)	68	29	—	69	29	—	70	37	—	70	32	—	72	49
11/5/10	ECf	69	33	—	69	28	—	70	43	(1)	69	30	—	70	32	—	70	37	—	70	33	—	72	49
11/5/10 (PM)	ECf	72	33	—	72	27	(2)	74	28	—	72	29	—	73	33	—	73	37	—	73	32	—	75	43
11/7/10	SV	69	33	—	70	28	—	71	27	(2)	70	31	—	70	36	—	72	37	—	72	32	—	72	54
11/8/10	ECf	66	34	—	66	29	—	67	28	—	66	32	—	67	35	—	68	36	—	68	33	—	68	~~55~~ 37 ECf 11/8/10

Abbreviations: T = Temperature (°F) RH = Relative Humidity (%) FN = Foot Note
Limits: RH 33± 5% Temperature 22± 5 °C or 71.6± 9°F
Foot Notes: 1 = Desiccant Changed 2 = Desiccator < 28% Humidity

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica West Sacramento
Air Toxics
Desiccator Humidity/Temperature Logbook

Desiccator #		1			2			3			4			5			6			7			Amb	
Date	Init	T	RH	FN	T	RH	FN	T	RH	FN	T	RH	FN	T	RH	FN	T	RH	FN	T	RH	FN	T	RH
11-30-10	AV	65	30	-	65	33	-	67	28	-	66	33	-	66	29	-	68	33	-	68	32	-	68	29
12/1/10	ECf	64	31	—	65	33	—	66	28	—	65	33	—	66	31	-	68	33	—	68	32	—	68	30
12/2/10	ECf	64	31	—	65	33	—	66	28	—	65	33	—	65	32	—	68	33	—	68	32	—	68	32
12-3-10	AV	67	31	—	68	33	-	69	27	(2)	67	34	-	68	32	-	70	33	—	70	32	-	70	36
12/6/10	ECf	67	32	—	68	37	—	69	30	—	68	36	—	68	35	—	70	34	—	70	32	—	70	47
12/7/10	ECf	66	34	—	67	36	—	68	30	—	67	36	—	67	34	—	70	34	—	68	32	—	70	38

Abbreviations: T = Temperature (°F) RH = Relative Humidity (%) FN = Foot Note
Limits: RH 33± 5% Temperature 22± 5 °C or 71.6± 9°F
Foot Notes: 1 = Desiccant Changed 2 = Desiccator < 28% Humidity

Requested By: VALMORES

Batch	Lot/Sample ID	Analysis Code	W/O#	Group	Message
0341292					Release Requested
0341292					Successfully Released