TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

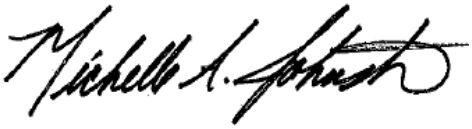
ANALYTICAL REPORT

Job Number: 280-67711-2

Job Description: Fort Wingate, New Mexico

For:
Sundance Consulting, Inc
6700 Jefferson Blvd NE
Albuquerque, NM 87109

Attention: John Nance



Approved for release.
Michelle A Johnston
Project Manager II
5/1/2015 3:10 PM

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05/01/2015

cc: Elizabeth Farias
Jim Lockhart
Ben Moayyad
Mr. Doug Scott

The test results in this report relate only to the samples in this report and meet all requirements of NELAC, with any exceptions noted. Pursuant to NELAP, this report shall not be reproduced except in full, without the written approval of the laboratory. All questions regarding this report should be directed to the TestAmerica Denver Project Manager.

The Lab Certification ID# is 4025.

Reporting limits are adjusted for sample size used, dilutions and moisture content if applicable.

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CASE NARRATIVE
Client: Sundance Consulting, Inc.
Project: Fort Wingate, New Mexico
Report Number: 280-67711-2

With the exceptions noted as flags or footnotes, standard analytical protocols were followed in the analysis of the samples and no problems were encountered or anomalies observed. In addition all laboratory quality control samples were within established control limits, with any exceptions noted below. Each sample was analyzed to achieve the lowest possible reporting limit within the constraints of the method. In some cases, due to interference or analytes present at high concentrations, samples were diluted. For diluted samples, the reporting limits are adjusted relative to the dilution required.

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

All holding times were met and proper preservation noted for the methods performed on these samples, unless otherwise detailed in the individual sections below.

Sample Receipt

The following report contains the analytical results for three water samples received April 11, 2015, according to documented sample acceptance procedures. The samples were received at temperatures of 1.0°C, 1.4°C, 5.6°C, 0.9°C and 1.2°C.

The 8081A page of the chain-of-custody lists the sample ID as DMW43042015, but all other pages of the COC lists the ID as DTW43042015 (280-67711-2). The sample ID was logged as DTW43042015. The client was notified on April 13, 2015.

Additional samples/analyses requested on the chain-of-custody are reported under separate cover (280-67711-1).

No other anomalies were encountered during sample receipt.

GC/MS Semivolatiles - 8270D

Samples TMW43042015 (280-67711-1), DTW43042015 (280-67711-2) and TMW47042015 (280-67711-4) were analyzed for semivolatile organic compounds (GC-MS) in accordance with EPA SW-846 Method 8270D. The samples were prepared on 04/15/2015 and analyzed on 04/18/2015.

Please note the Caprolactam data are reported under separate cover, as the laboratory does not hold DOD ELAP certification for this compound. The laboratory does not maintain quarterly QC requirements for precision, accuracy and detections.

Reporting limits and method detection limits have been adjusted accordingly for the initial volumes extracted.

A deviation from the Standard Operating Procedure (SOP) occurred. The details are as follows:
The samples were extracted per SW-846 3520C for 18 hours at a pH of 1-2 and 18 hours at a pH of 11-12; however, the laboratory's SOP requires the samples be extracted for 24 hours at each pH range. The laboratory's SOP is in the process of being revised for these changes.

Due to an analyst oversight, the MS/MSD was spiked with the LCS Main and LCS Supp after the acid had been added. Per the SOP, the acid is supposed to be added after the samples are spiked.

No other analytical or quality issues were noted, other than those described in the Definitions/Glossary page.

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Instrument ID: SMS_G6 Analysis Batch Number: 272059

Lab Sample ID: ICIS 280-272059/3 Client Sample ID:

Date Analyzed: 02/25/15 11:53 Lab File ID: G6_16637.D GC Column: Vf-5MS (30.25) ID: 0.25(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Caprolactam	6.43	Split Peak	kiekeld	04/10/15 08:28
Indeno[1,2,3-cd]pyrene	21.13	Shouldering	kiekeld	04/10/15 08:28

Lab Sample ID: STD020 280-272059/6 IC Client Sample ID:

Date Analyzed: 02/25/15 13:12 Lab File ID: G6_16640.D GC Column: Vf-5MS (30.25) ID: 0.25(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Caprolactam	6.41	Split Peak	hoeflera	04/10/15 09:04

Lab Sample ID: STD050 280-272059/7 IC Client Sample ID:

Date Analyzed: 02/25/15 13:39 Lab File ID: G6_16641.D GC Column: Vf-5MS (30.25) ID: 0.25(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Caprolactam	6.42	Split Peak	hoeflera	04/10/15 09:09
Indeno[1,2,3-cd]pyrene	21.12	Shouldering	hoeflera	04/10/15 09:09

Lab Sample ID: STD120 280-272059/8 IC Client Sample ID:

Date Analyzed: 02/25/15 14:06 Lab File ID: G6_16642.D GC Column: Vf-5MS (30.25) ID: 0.25(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Caprolactam	6.43	Split Peak	hoeflera	04/10/15 09:12
Indeno[1,2,3-cd]pyrene	21.13	Shouldering	hoeflera	04/10/15 09:12

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Instrument ID: SMS_G6 Analysis Batch Number: 272059

Lab Sample ID: STD160 280-272059/9 IC Client Sample ID:

Date Analyzed: 02/25/15 14:32 Lab File ID: G6_16643.D GC Column: Vf-5MS (30.25) ID: 0.25(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Caprolactam	6.45	Split Peak	hoeflera	04/10/15 09:14
Indeno[1,2,3-cd]pyrene	21.15	Shouldering	hoeflera	04/10/15 09:14

Lab Sample ID: STD200 280-272059/10 IC Client Sample ID:

Date Analyzed: 02/25/15 14:59 Lab File ID: G6_16644.D GC Column: Vf-5MS (30.25) ID: 0.25(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Caprolactam	6.45	Split Peak	hoeflera	04/10/15 09:22
Indeno[1,2,3-cd]pyrene	21.16	Shouldering	hoeflera	04/10/15 09:22

Lab Sample ID: ICV 280-272059/11 Client Sample ID:

Date Analyzed: 02/25/15 15:26 Lab File ID: G6_16645.D GC Column: Vf-5MS (30.25) ID: 0.25(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Indeno[1,2,3-cd]pyrene	21.13	Shouldering	hoeflera	04/10/15 09:24

GC/MS SEMI VOA MANUAL INTEGRATION SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Instrument ID: SMS_G6 Analysis Batch Number: 273380

Lab Sample ID: CCV 280-273380/3 Client Sample ID:

Date Analyzed: 04/18/15 15:20 Lab File ID: G6_17414.D GC Column: Vf-5MS (30.25) ID: 0.25(mm)

COMPOUND NAME	RETENTION TIME	MANUAL INTEGRATION		
		REASON	ANALYST	DATE
Indeno[1,2,3-cd]pyrene	20.86	Shouldering	hoeflera	04/18/15 16:20

SAMPLE SUMMARY

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received
280-67711-1	TMW43042015	Water	04/10/2015 0825	04/11/2015 0930
280-67711-1MS	TMW43042015	Water	04/10/2015 0825	04/11/2015 0930
280-67711-1MSD	TMW43042015	Water	04/10/2015 0825	04/11/2015 0930
280-67711-2	DTW43042015	Water	04/10/2015 0825	04/11/2015 0930
280-67711-4	TMW47042015	Water	04/10/2015 0930	04/11/2015 0930

EXECUTIVE SUMMARY - Detections

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Lab Sample ID Analyte	Client Sample ID	Result	Qualifier	Reporting Limit	Units	Method

No Detections

METHOD SUMMARY

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Semivolatile Organic Compounds (GC/MS)	TAL DEN	SW846 8270D	
Liquid-Liquid Extraction (Continuous)			SW846 3520C

Lab References:

TAL DEN = TestAmerica Denver

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

METHOD / ANALYST SUMMARY

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Method	Analyst	Analyst ID
SW846 8270D	Hoefler, Alexandra F	AFH

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Client Sample ID: **TMW43042015**

Lab Sample ID: 280-67711-1
Client Matrix: Water

Date Sampled: 04/10/2015 0825
Date Received: 04/11/2015 0930

8270D Semivolatile Organic Compounds (GC/MS)

Analysis Method:	8270D	Analysis Batch:	280-273380	Instrument ID:	SMS_G6
Prep Method:	3520C	Prep Batch:	280-272870	Lab File ID:	G6_17428.D
Dilution:	1.0			Initial Weight/Volume:	925 mL
Analysis Date:	04/18/2015 2113			Final Weight/Volume:	1 mL
Prep Date:	04/15/2015 1850			Injection Volume:	0.5 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Caprolactam	2.7	U	2.7	5.4

Surrogate	%Rec	Qualifier	Acceptance Limits
2,4,6-Tribromophenol (Surr)	67		48 - 135
2-Fluorobiphenyl	64		48 - 135
2-Fluorophenol (Surr)	65		41 - 135
Nitrobenzene-d5 (Surr)	66		42 - 135
Phenol-d5 (Surr)	66		46 - 135
Terphenyl-d14 (Surr)	65		20 - 135

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Client Sample ID: **DTW43042015**

Lab Sample ID: 280-67711-2
Client Matrix: Water

Date Sampled: 04/10/2015 0825
Date Received: 04/11/2015 0930

8270D Semivolatile Organic Compounds (GC/MS)

Analysis Method:	8270D	Analysis Batch:	280-273380	Instrument ID:	SMS_G6
Prep Method:	3520C	Prep Batch:	280-272870	Lab File ID:	G6_17431.D
Dilution:	1.0			Initial Weight/Volume:	1029 mL
Analysis Date:	04/18/2015 2232			Final Weight/Volume:	1 mL
Prep Date:	04/15/2015 1850			Injection Volume:	0.5 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Caprolactam	2.4	U	2.4	4.9

Surrogate	%Rec	Qualifier	Acceptance Limits
2,4,6-Tribromophenol (Surr)	67		48 - 135
2-Fluorobiphenyl	66		48 - 135
2-Fluorophenol (Surr)	65		41 - 135
Nitrobenzene-d5 (Surr)	66		42 - 135
Phenol-d5 (Surr)	66		46 - 135
Terphenyl-d14 (Surr)	64		20 - 135

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Client Sample ID: **TMW47042015**

Lab Sample ID: 280-67711-4
Client Matrix: Water

Date Sampled: 04/10/2015 0930
Date Received: 04/11/2015 0930

8270D Semivolatile Organic Compounds (GC/MS)

Analysis Method:	8270D	Analysis Batch:	280-273380	Instrument ID:	SMS_G6
Prep Method:	3520C	Prep Batch:	280-272870	Lab File ID:	G6_17432.D
Dilution:	1.0			Initial Weight/Volume:	963.2 mL
Analysis Date:	04/18/2015 2259			Final Weight/Volume:	1 mL
Prep Date:	04/15/2015 1850			Injection Volume:	0.5 uL

Analyte	Result (ug/L)	Qualifier	MDL	RL
Caprolactam	2.6	U	2.6	5.2

Surrogate	%Rec	Qualifier	Acceptance Limits
2,4,6-Tribromophenol (Surr)	60		48 - 135
2-Fluorobiphenyl	53		48 - 135
2-Fluorophenol (Surr)	51		41 - 135
Nitrobenzene-d5 (Surr)	53		42 - 135
Phenol-d5 (Surr)	53		46 - 135
Terphenyl-d14 (Surr)	63		20 - 135

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Surrogate Recovery Report

8270D Semivolatile Organic Compounds (GC/MS)

Client Matrix: Water

Lab Sample ID	Client Sample ID	2FP %Rec	PHL %Rec	NBZ %Rec	FBP %Rec	TBP %Rec	TPH %Rec
280-67711-1	TMW43042015	65	66	66	64	67	65
280-67711-2	DTW43042015	65	66	66	66	67	64
280-67711-4	TMW47042015	51	53	53	53	60	63
MB 280-272870/1-A		64	64	65	62	65	64
LCS 280-272870/2-A		60	60	61	61	68	62
280-67711-1 MS	TMW43042015 MS	58	60	58	61	72	65
280-67711-1 MSD	TMW43042015 MSD	62	62	63	61	66	62

Surrogate	Acceptance Limits
2FP = 2-Fluorophenol (Surr)	41-135
PHL = Phenol-d5 (Surr)	46-135
NBZ = Nitrobenzene-d5 (Surr)	42-135
FBP = 2-Fluorobiphenyl	48-135
TBP = 2,4,6-Tribromophenol (Surr)	48-135
TPH = Terphenyl-d14 (Surr)	20-135

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Method Blank - Batch: 280-272870

Method: 8270D
Preparation: 3520C

Lab Sample ID:	MB 280-272870/1-A	Analysis Batch:	280-273380	Instrument ID:	SMS_G6
Client Matrix:	Water	Prep Batch:	280-272870	Lab File ID:	G6_17420.D
Dilution:	1.0	Leach Batch:	N/A	Initial Weight/Volume:	1000 mL
Analysis Date:	04/18/2015 1749	Units:	ug/L	Final Weight/Volume:	1 mL
Prep Date:	04/15/2015 1850			Injection Volume:	0.5 uL
Leach Date:	N/A				

Analyte	Result	Qual	MDL	RL
Caprolactam	2.5	U	2.5	5.0

Surrogate	% Rec	Acceptance Limits
2,4,6-Tribromophenol (Surr)	65	48 - 135
2-Fluorobiphenyl	62	48 - 135
2-Fluorophenol (Surr)	64	41 - 135
Nitrobenzene-d5 (Surr)	65	42 - 135
Phenol-d5 (Surr)	64	46 - 135
Terphenyl-d14 (Surr)	64	20 - 135

Lab Control Sample - Batch: 280-272870

Method: 8270D
Preparation: 3520C

Lab Sample ID:	LCS 280-272870/2-A	Analysis Batch:	280-273380	Instrument ID:	SMS_G6
Client Matrix:	Water	Prep Batch:	280-272870	Lab File ID:	G6_17421.D
Dilution:	1.0	Leach Batch:	N/A	Initial Weight/Volume:	1000 mL
Analysis Date:	04/18/2015 1816	Units:	ug/L	Final Weight/Volume:	1 mL
Prep Date:	04/15/2015 1850			Injection Volume:	0.5 uL
Leach Date:	N/A				

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Caprolactam	80.0	52.3	65	64 - 120	

Surrogate	% Rec	Acceptance Limits
2,4,6-Tribromophenol (Surr)	68	48 - 135
2-Fluorobiphenyl	61	48 - 135
2-Fluorophenol (Surr)	60	41 - 135
Nitrobenzene-d5 (Surr)	61	42 - 135
Phenol-d5 (Surr)	60	46 - 135
Terphenyl-d14 (Surr)	62	20 - 135

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Matrix Spike/
Matrix Spike Duplicate Recovery Report - Batch: 280-272870

Method: 8270D
Preparation: 3520C

MS Lab Sample ID:	280-67711-1	Analysis Batch:	280-273380	Instrument ID:	SMS_G6
Client Matrix:	Water	Prep Batch:	280-272870	Lab File ID:	G6_17429.D
Dilution:	1.0	Leach Batch:	N/A	Initial Weight/Volume:	885 mL
Analysis Date:	04/18/2015 2139			Final Weight/Volume:	1 mL
Prep Date:	04/15/2015 1850			Injection Volume:	0.5 uL
Leach Date:	N/A				

MSD Lab Sample ID:	280-67711-1	Analysis Batch:	280-273380	Instrument ID:	SMS_G6
Client Matrix:	Water	Prep Batch:	280-272870	Lab File ID:	G6_17430.D
Dilution:	1.0	Leach Batch:	N/A	Initial Weight/Volume:	903.3 mL
Analysis Date:	04/18/2015 2206			Final Weight/Volume:	1 mL
Prep Date:	04/15/2015 1850			Injection Volume:	0.5 uL
Leach Date:	N/A				

	% Rec.						
Analyte	MS	MSD	Limit	RPD	RPD Limit	MS Qual	MSD Qual
Caprolactam	69	66	64 - 120	6	30		

Surrogate	MS % Rec	MSD % Rec	Acceptance Limits
2,4,6-Tribromophenol (Surr)	72	66	48 - 135
2-Fluorobiphenyl	61	61	48 - 135
2-Fluorophenol (Surr)	58	62	41 - 135
Nitrobenzene-d5 (Surr)	58	63	42 - 135
Phenol-d5 (Surr)	60	62	46 - 135
Terphenyl-d14 (Surr)	65	62	20 - 135

Matrix Spike/
Matrix Spike Duplicate Recovery Report - Batch: 280-272870

Method: 8270D
Preparation: 3520C

MS Lab Sample ID:	280-67711-1	Units:	ug/L	MSD Lab Sample ID:	280-67711-1
Client Matrix:	Water			Client Matrix:	Water
Dilution:	1.0			Dilution:	1.0
Analysis Date:	04/18/2015 2139			Analysis Date:	04/18/2015 2206
Prep Date:	04/15/2015 1850			Prep Date:	04/15/2015 1850
Leach Date:	N/A			Leach Date:	N/A

Analyte	Sample Result/Qual		MS Spike Amount	MSD Spike Amount	MS Result/Qual	MSD Result/Qual
Caprolactam	2.7	U	90.4	88.6	62.1	58.4

DATA REPORTING QUALIFIERS

Client: Sundance Consulting, Inc Job Number: 280-67711-2

Lab Section	Qualifier	Description
GC/MS Semi VOA		
	U	Indicates the analyte was analyzed for but not detected.

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC/MS Semi VOA					
Prep Batch: 280-272870					
LCS 280-272870/2-A	Lab Control Sample	T	Water	3520C	
MB 280-272870/1-A	Method Blank	T	Water	3520C	
280-67711-1	TMW43042015	T	Water	3520C	
280-67711-1MS	Matrix Spike	T	Water	3520C	
280-67711-1MSD	Matrix Spike Duplicate	T	Water	3520C	
280-67711-2	DTW43042015	T	Water	3520C	
280-67711-4	TMW47042015	T	Water	3520C	
Analysis Batch:280-273380					
LCS 280-272870/2-A	Lab Control Sample	T	Water	8270D	280-272870
MB 280-272870/1-A	Method Blank	T	Water	8270D	280-272870
280-67711-1	TMW43042015	T	Water	8270D	280-272870
280-67711-1MS	Matrix Spike	T	Water	8270D	280-272870
280-67711-1MSD	Matrix Spike Duplicate	T	Water	8270D	280-272870
280-67711-2	DTW43042015	T	Water	8270D	280-272870
280-67711-4	TMW47042015	T	Water	8270D	280-272870

Report Basis

T = Total

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Laboratory Chronicle

Lab ID: 280-67711-1 **Client ID:** TMW43042015

Sample Date/Time: 04/10/2015 08:25 Received Date/Time: 04/11/2015 09:30

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
P:3520C	280-67711-C-1-A		280-273380	280-272870	04/15/2015 18:50	1	TAL DEN	JRK
A:8270D	280-67711-C-1-A		280-273380	280-272870	04/18/2015 21:13	1	TAL DEN	AFH

Lab ID: 280-67711-1 **Client ID:** TMW43042015

Sample Date/Time: 04/10/2015 08:25 Received Date/Time: 04/11/2015 09:30

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
P:3520C	280-67711-B-1-B MS		280-273380	280-272870	04/15/2015 18:50	1	TAL DEN	JRK
A:8270D	280-67711-B-1-B MS		280-273380	280-272870	04/18/2015 21:39	1	TAL DEN	AFH

Lab ID: 280-67711-1 **Client ID:** TMW43042015

Sample Date/Time: 04/10/2015 08:25 Received Date/Time: 04/11/2015 09:30

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
P:3520C	280-67711-L-1-B MSD		280-273380	280-272870	04/15/2015 18:50	1	TAL DEN	JRK
A:8270D	280-67711-L-1-B MSD		280-273380	280-272870	04/18/2015 22:06	1	TAL DEN	AFH

Lab ID: 280-67711-2 **Client ID:** DTW43042015

Sample Date/Time: 04/10/2015 08:25 Received Date/Time: 04/11/2015 09:30

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
P:3520C	280-67711-B-2-A		280-273380	280-272870	04/15/2015 18:50	1	TAL DEN	JRK
A:8270D	280-67711-B-2-A		280-273380	280-272870	04/18/2015 22:32	1	TAL DEN	AFH

Lab ID: 280-67711-4 **Client ID:** TMW47042015

Sample Date/Time: 04/10/2015 09:30 Received Date/Time: 04/11/2015 09:30

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
P:3520C	280-67711-C-4-A		280-273380	280-272870	04/15/2015 18:50	1	TAL DEN	JRK
A:8270D	280-67711-C-4-A		280-273380	280-272870	04/18/2015 22:59	1	TAL DEN	AFH

Lab ID: MB **Client ID:** N/A

Sample Date/Time: N/A Received Date/Time: N/A

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
P:3520C	MB 280-272870/1-A		280-273380	280-272870	04/15/2015 18:50	1	TAL DEN	JRK
A:8270D	MB 280-272870/1-A		280-273380	280-272870	04/18/2015 17:49	1	TAL DEN	AFH

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Laboratory Chronicle

Lab ID: **LCS** **Client ID:** **N/A**

Sample Date/Time: N/A Received Date/Time: N/A

Method	Bottle ID	Run	Analysis Batch	Prep Batch	Date Prepared / Analyzed	Dil	Lab	Analyst
P:3520C	LCS 280-272870/2-A		280-273380	280-272870	04/15/2015 18:50	1	TAL DEN	JRK
A:8270D	LCS 280-272870/2-A		280-273380	280-272870	04/18/2015 18:16	1	TAL DEN	AFH

Lab References:

TAL DEN = TestAmerica Denver

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
8270_LCS_Main_00020	03/27/16	03/27/15	P&T Methanol, Lot MethanolP&T_00111	250 mL	MS-569729_00013	5 mL	1,1'-Biphenyl	80 ug/mL
							1,2,4,5-Tetrachlorobenzene	80 ug/mL
							1,2,4-Trichlorobenzene	80 ug/mL
							1,2-Dichlorobenzene	80 ug/mL
							1,2-Diphenylhydrazine	80.878 ug/mL
							1,3-Dichlorobenzene	80 ug/mL
							1,3-Dinitrobenzene	80 ug/mL
							1,4-Dichlorobenzene	80 ug/mL
							1,4-Dioxane	80 ug/mL
							1-Methylnaphthalene	80 ug/mL
							2,2'-oxybis[1-chloropropane]	80 ug/mL
							2,3,4,6-Tetrachlorophenol	80 ug/mL
							2,4,5-Trichlorophenol	80 ug/mL
							2,4,6-Trichlorophenol	80 ug/mL
							2,4-Dichlorophenol	80 ug/mL
							2,4-Dimethylphenol	80 ug/mL
							2,4-Dinitrophenol	160 ug/mL
							2,4-Dinitrotoluene	80 ug/mL
							2,6-Dichlorophenol	80 ug/mL
							2,6-Dinitrotoluene	80 ug/mL
							2-Chloronaphthalene	80 ug/mL
							2-Chlorophenol	80 ug/mL
							2-Methylnaphthalene	80 ug/mL
							2-Methylphenol	80 ug/mL
							2-Nitroaniline	80 ug/mL
							2-Nitrophenol	80 ug/mL
							3 & 4 Methylphenol	80 ug/mL
							3-Methylphenol	80 ug/mL
							3-Nitroaniline	80 ug/mL
							4,6-Dinitro-2-methylphenol	160 ug/mL
							4-Bromophenyl phenyl ether	80 ug/mL
							4-Chloro-3-methylphenol	80 ug/mL
							4-Chloroaniline	80 ug/mL
							4-Chlorophenyl phenyl ether	80 ug/mL
							4-Methylphenol	80 ug/mL
							4-Nitroaniline	80 ug/mL
							4-Nitrophenol	160 ug/mL
							Acenaphthene	80 ug/mL
							Acenaphthylene	80 ug/mL
							Acetophenone	80 ug/mL
							Aniline	80 ug/mL
							Anthracene	80 ug/mL
							Azobenzene	80 ug/mL
							Benzo[a]anthracene	80 ug/mL
							Benzo[a]pyrene	80 ug/mL
							Benzo[b]fluoranthene	80 ug/mL
							Benzo[g,h,i]perylene	80 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Benzo[k]fluoranthene	80 ug/mL
							Benzyl alcohol	80 ug/mL
							Bis(2-chloroethoxy)methane	80 ug/mL
							Bis(2-chloroethyl)ether	80 ug/mL
							Bis(2-ethylhexyl) phthalate	80 ug/mL
							Butyl benzyl phthalate	80 ug/mL
							Carbazole	80 ug/mL
							Chrysene	80 ug/mL
							Di-n-butyl phthalate	80 ug/mL
							Di-n-octyl phthalate	80 ug/mL
							Dibenz(a,h)anthracene	80 ug/mL
							Dibenzofuran	80 ug/mL
							Diethyl phthalate	80 ug/mL
							Dimethyl phthalate	80 ug/mL
							Diphenylamine	137.072 ug/mL
							Fluoranthene	80 ug/mL
							Fluorene	80 ug/mL
							Hexachlorobenzene	80 ug/mL
							Hexachlorobutadiene	80 ug/mL
							Hexachlorocyclopentadiene	80 ug/mL
							Hexachloroethane	80 ug/mL
							Hexadecane	80 ug/mL
							Indeno[1,2,3-cd]pyrene	80 ug/mL
							Isophorone	80 ug/mL
							n-Decane	80 ug/mL
							N-Nitrosodi-n-propylamine	80 ug/mL
							N-Nitrosodimethylamine	80 ug/mL
							N-Nitrosodiphenylamine	160 ug/mL
							n-Octadecane	80 ug/mL
							Naphthalene	80 ug/mL
							Nitrobenzene	80 ug/mL
							Pentachlorophenol	160 ug/mL
							Phenanthrene	80 ug/mL
							Phenol	80 ug/mL
							Pyrene	80 ug/mL
							Pyridine	80 ug/mL
					MS-569729_00014	5 mL	1,1'-Biphenyl	80 ug/mL
							1,2,4,5-Tetrachlorobenzene	80 ug/mL
							1,2,4-Trichlorobenzene	80 ug/mL
							1,2-Dichlorobenzene	80 ug/mL
							1,2-Diphenylhydrazine	80.878 ug/mL
							1,3-Dichlorobenzene	80 ug/mL
							1,3-Dinitrobenzene	80 ug/mL
							1,4-Dichlorobenzene	80 ug/mL
							1,4-Dioxane	80 ug/mL
							1-Methylnaphthalene	80 ug/mL
							2,2'-oxybis[1-chloropropane]	80 ug/mL
							2,3,4,6-Tetrachlorophenol	80 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2,4,5-Trichlorophenol	80 ug/mL
							2,4,6-Trichlorophenol	80 ug/mL
							2,4-Dichlorophenol	80 ug/mL
							2,4-Dimethylphenol	80 ug/mL
							2,4-Dinitrophenol	160 ug/mL
							2,4-Dinitrotoluene	80 ug/mL
							2,6-Dichlorophenol	80 ug/mL
							2,6-Dinitrotoluene	80 ug/mL
							2-Chloronaphthalene	80 ug/mL
							2-Chlorophenol	80 ug/mL
							2-Methylnaphthalene	80 ug/mL
							2-Methylphenol	80 ug/mL
							2-Nitroaniline	80 ug/mL
							2-Nitrophenol	80 ug/mL
							3 & 4 Methylphenol	80 ug/mL
							3-Methylphenol	80 ug/mL
							3-Nitroaniline	80 ug/mL
							4,6-Dinitro-2-methylphenol	160 ug/mL
							4-Bromophenyl phenyl ether	80 ug/mL
							4-Chloro-3-methylphenol	80 ug/mL
							4-Chloroaniline	80 ug/mL
							4-Chlorophenyl phenyl ether	80 ug/mL
							4-Methylphenol	80 ug/mL
							4-Nitroaniline	80 ug/mL
							4-Nitrophenol	160 ug/mL
							Acenaphthene	80 ug/mL
							Acenaphthylene	80 ug/mL
							Acetophenone	80 ug/mL
							Aniline	80 ug/mL
							Anthracene	80 ug/mL
							Azobenzene	80 ug/mL
							Benzo[a]anthracene	80 ug/mL
							Benzo[a]pyrene	80 ug/mL
							Benzo[b]fluoranthene	80 ug/mL
							Benzo[g,h,i]perylene	80 ug/mL
							Benzo[k]fluoranthene	80 ug/mL
							Benzyl alcohol	80 ug/mL
							Bis(2-chloroethoxy)methane	80 ug/mL
							Bis(2-chloroethyl)ether	80 ug/mL
							Bis(2-ethylhexyl) phthalate	80 ug/mL
							Butyl benzyl phthalate	80 ug/mL
							Carbazole	80 ug/mL
							Chrysene	80 ug/mL
							Di-n-butyl phthalate	80 ug/mL
							Di-n-octyl phthalate	80 ug/mL
							Dibenz(a,h)anthracene	80 ug/mL
							Dibenzofuran	80 ug/mL
							Diethyl phthalate	80 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Dimethyl phthalate	80 ug/mL
							Diphenylamine	137.072 ug/mL
							Fluoranthene	80 ug/mL
							Fluorene	80 ug/mL
							Hexachlorobenzene	80 ug/mL
							Hexachlorobutadiene	80 ug/mL
							Hexachlorocyclopentadiene	80 ug/mL
							Hexachloroethane	80 ug/mL
							Hexadecane	80 ug/mL
							Indeno[1,2,3-cd]pyrene	80 ug/mL
							Isophorone	80 ug/mL
							n-Decane	80 ug/mL
							N-Nitrosodi-n-propylamine	80 ug/mL
							N-Nitrosodimethylamine	80 ug/mL
							N-Nitrosodiphenylamine	160 ug/mL
							n-Octadecane	80 ug/mL
							Naphthalene	80 ug/mL
							Nitrobenzene	80 ug/mL
							Pentachlorophenol	160 ug/mL
							Phenanthrene	80 ug/mL
							Phenol	80 ug/mL
							Pyrene	80 ug/mL
							Pyridine	80 ug/mL
					MS-569729_00015	5 mL	1,1'-Biphenyl	80 ug/mL
							1,2,4,5-Tetrachlorobenzene	80 ug/mL
							1,2,4-Trichlorobenzene	80 ug/mL
							1,2-Dichlorobenzene	80 ug/mL
							1,2-Diphenylhydrazine	80.878 ug/mL
							1,3-Dichlorobenzene	80 ug/mL
							1,3-Dinitrobenzene	80 ug/mL
							1,4-Dichlorobenzene	80 ug/mL
							1,4-Dioxane	80 ug/mL
							1-Methylnaphthalene	80 ug/mL
							2,2'-oxybis[1-chloropropane]	80 ug/mL
							2,3,4,6-Tetrachlorophenol	80 ug/mL
							2,4,5-Trichlorophenol	80 ug/mL
							2,4,6-Trichlorophenol	80 ug/mL
							2,4-Dichlorophenol	80 ug/mL
							2,4-Dimethylphenol	80 ug/mL
							2,4-Dinitrophenol	160 ug/mL
							2,4-Dinitrotoluene	80 ug/mL
							2,6-Dichlorophenol	80 ug/mL
							2,6-Dinitrotoluene	80 ug/mL
							2-Chloronaphthalene	80 ug/mL
							2-Chlorophenol	80 ug/mL
							2-Methylnaphthalene	80 ug/mL
							2-Methylphenol	80 ug/mL
							2-Nitroaniline	80 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Nitrophenol	80 ug/mL
							3 & 4 Methylphenol	80 ug/mL
							3-Methylphenol	80 ug/mL
							3-Nitroaniline	80 ug/mL
							4,6-Dinitro-2-methylphenol	160 ug/mL
							4-Bromophenyl phenyl ether	80 ug/mL
							4-Chloro-3-methylphenol	80 ug/mL
							4-Chloroaniline	80 ug/mL
							4-Chlorophenyl phenyl ether	80 ug/mL
							4-Methylphenol	80 ug/mL
							4-Nitroaniline	80 ug/mL
							4-Nitrophenol	160 ug/mL
							Acenaphthene	80 ug/mL
							Acenaphthylene	80 ug/mL
							Acetophenone	80 ug/mL
							Aniline	80 ug/mL
							Anthracene	80 ug/mL
							Azobenzene	80 ug/mL
							Benzo[a]anthracene	80 ug/mL
							Benzo[a]pyrene	80 ug/mL
							Benzo[b]fluoranthene	80 ug/mL
							Benzo[g,h,i]perylene	80 ug/mL
							Benzo[k]fluoranthene	80 ug/mL
							Benzyl alcohol	80 ug/mL
							Bis(2-chloroethoxy)methane	80 ug/mL
							Bis(2-chloroethyl)ether	80 ug/mL
							Bis(2-ethylhexyl) phthalate	80 ug/mL
							Butyl benzyl phthalate	80 ug/mL
							Carbazole	80 ug/mL
							Chrysene	80 ug/mL
							Di-n-butyl phthalate	80 ug/mL
							Di-n-octyl phthalate	80 ug/mL
							Dibenz(a,h)anthracene	80 ug/mL
							Dibenzofuran	80 ug/mL
							Diethyl phthalate	80 ug/mL
							Dimethyl phthalate	80 ug/mL
							Diphenylamine	137.072 ug/mL
							Fluoranthene	80 ug/mL
							Fluorene	80 ug/mL
							Hexachlorobenzene	80 ug/mL
							Hexachlorobutadiene	80 ug/mL
							Hexachlorocyclopentadiene	80 ug/mL
							Hexachloroethane	80 ug/mL
							Hexadecane	80 ug/mL
							Indeno[1,2,3-cd]pyrene	80 ug/mL
							Isophorone	80 ug/mL
							n-Decane	80 ug/mL
							N-Nitrosodi-n-propylamine	80 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							N-Nitrosodimethylamine	80 ug/mL
							N-Nitrosodiphenylamine	160 ug/mL
							n-Octadecane	80 ug/mL
							Naphthalene	80 ug/mL
							Nitrobenzene	80 ug/mL
							Pentachlorophenol	160 ug/mL
							Phenanthrene	80 ug/mL
							Phenol	80 ug/mL
							Pyrene	80 ug/mL
							Pyridine	80 ug/mL
					MS-569729_00016	5 mL	1,1'-Biphenyl	80 ug/mL
							1,2,4,5-Tetrachlorobenzene	80 ug/mL
							1,2,4-Trichlorobenzene	80 ug/mL
							1,2-Dichlorobenzene	80 ug/mL
							1,2-Diphenylhydrazine	80.878 ug/mL
							1,3-Dichlorobenzene	80 ug/mL
							1,3-Dinitrobenzene	80 ug/mL
							1,4-Dichlorobenzene	80 ug/mL
							1,4-Dioxane	80 ug/mL
							1-Methylnaphthalene	80 ug/mL
							2,2'-oxybis[1-chloropropane]	80 ug/mL
							2,3,4,6-Tetrachlorophenol	80 ug/mL
							2,4,5-Trichlorophenol	80 ug/mL
							2,4,6-Trichlorophenol	80 ug/mL
							2,4-Dichlorophenol	80 ug/mL
							2,4-Dimethylphenol	80 ug/mL
							2,4-Dinitrophenol	160 ug/mL
							2,4-Dinitrotoluene	80 ug/mL
							2,6-Dichlorophenol	80 ug/mL
							2,6-Dinitrotoluene	80 ug/mL
							2-Chloronaphthalene	80 ug/mL
							2-Chlorophenol	80 ug/mL
							2-Methylnaphthalene	80 ug/mL
							2-Methylphenol	80 ug/mL
							2-Nitroaniline	80 ug/mL
							2-Nitrophenol	80 ug/mL
							3 & 4 Methylphenol	80 ug/mL
							3-Methylphenol	80 ug/mL
							3-Nitroaniline	80 ug/mL
							4,6-Dinitro-2-methylphenol	160 ug/mL
							4-Bromophenyl phenyl ether	80 ug/mL
							4-Chloro-3-methylphenol	80 ug/mL
							4-Chloroaniline	80 ug/mL
							4-Chlorophenyl phenyl ether	80 ug/mL
							4-Methylphenol	80 ug/mL
							4-Nitroaniline	80 ug/mL
							4-Nitrophenol	160 ug/mL
							Acenaphthene	80 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Acenaphthylene	80 ug/mL
							Acetophenone	80 ug/mL
							Aniline	80 ug/mL
							Anthracene	80 ug/mL
							Azobenzene	80 ug/mL
							Benzo[a]anthracene	80 ug/mL
							Benzo[a]pyrene	80 ug/mL
							Benzo[b]fluoranthene	80 ug/mL
							Benzo[g,h,i]perylene	80 ug/mL
							Benzo[k]fluoranthene	80 ug/mL
							Benzyl alcohol	80 ug/mL
							Bis(2-chloroethoxy)methane	80 ug/mL
							Bis(2-chloroethyl)ether	80 ug/mL
							Bis(2-ethylhexyl) phthalate	80 ug/mL
							Butyl benzyl phthalate	80 ug/mL
							Carbazole	80 ug/mL
							Chrysene	80 ug/mL
							Di-n-butyl phthalate	80 ug/mL
							Di-n-octyl phthalate	80 ug/mL
							Dibenz(a,h)anthracene	80 ug/mL
							Dibenzofuran	80 ug/mL
							Diethyl phthalate	80 ug/mL
							Dimethyl phthalate	80 ug/mL
							Diphenylamine	137.072 ug/mL
							Fluoranthene	80 ug/mL
							Fluorene	80 ug/mL
							Hexachlorobenzene	80 ug/mL
							Hexachlorobutadiene	80 ug/mL
							Hexachlorocyclopentadiene	80 ug/mL
							Hexachloroethane	80 ug/mL
							Hexadecane	80 ug/mL
							Indeno[1,2,3-cd]pyrene	80 ug/mL
							Isophorone	80 ug/mL
							n-Decane	80 ug/mL
							N-Nitrosodi-n-propylamine	80 ug/mL
							N-Nitrosodimethylamine	80 ug/mL
							N-Nitrosodiphenylamine	160 ug/mL
							n-Octadecane	80 ug/mL
							Naphthalene	80 ug/mL
							Nitrobenzene	80 ug/mL
							Pentachlorophenol	160 ug/mL
							Phenanthrene	80 ug/mL
							Phenol	80 ug/mL
							Pyrene	80 ug/mL
							Pyridine	80 ug/mL
					MS-569731_00005	5 mL	Benzoic acid	80 ug/mL
							Indene	80 ug/mL
					MS-569731_00006	5 mL	Benzoic acid	80 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Indene	80 ug/mL
.MS-569729_00013	05/31/16		Restek, Lot A0107399		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dichlorophenol	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Diphenylamine	1713.4 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Hexadecane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							n-Decane	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							N-Nitrosodiphenylamine	2000 ug/mL
							n-Octadecane	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
.MS-569729_00014	05/31/16		Restek, Lot A0107399		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dichlorophenol	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Dimethyl phthalate	1000 ug/mL
							Diphenylamine	1713.4 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Hexadecane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							n-Decane	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							N-Nitrosodiphenylamine	2000 ug/mL
							n-Octadecane	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
.MS-569729_00015	05/31/16		Restek, Lot A0107399		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dichlorophenol	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Diphenylamine	1713.4 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Hexadecane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							n-Decane	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							N-Nitrosodimethylamine	1000 ug/mL
							N-Nitrosodiphenylamine	2000 ug/mL
							n-Octadecane	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
.MS-569729_00016	05/31/16		Restek, Lot A0107399		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dichlorophenol	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Diphenylamine	1713.4 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Hexadecane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							n-Decane	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							N-Nitrosodiphenylamine	2000 ug/mL
							n-Octadecane	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
.MS-569731_00005	06/30/16		Restek, Lot A0107943		(Purchased Reagent)		Benzoic acid	2000 ug/mL
							Indene	2000 ug/mL
.MS-569731_00006	06/30/16		Restek, Lot A0107943		(Purchased Reagent)		Benzoic acid	2000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Indene	2000 ug/mL
8270_LCS_Supp_00103	04/20/15	04/13/15	P&T Methanol, Lot MethanolP&T_00112	50 mL	MS-569730_00016	2 mL	3,3'-Dichlorobenzidine	80 ug/mL
							Benzidine	80 ug/mL
					MS-569732_00011	2 mL	Atrazine	80 ug/mL
							Benzaldehyde	80 ug/mL
							Caprolactam	80 ug/mL
.MS-569730_00016	02/11/16		Restek, Lot A0108709		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Benzidine	2000 ug/mL
.MS-569732_00011	06/30/16		Restek, Lot A0108035		(Purchased Reagent)		Atrazine	2000 ug/mL
							Benzaldehyde	2000 ug/mL
							Caprolactam	2000 ug/mL
8270Surrogate_00078	04/07/16	04/07/15	ACETONE, Lot Acetone_000115	1000 mL	8270SurStkHL_00078	5 mL	2,4,6 - Tribromophenol	100 ug/mL
							2,4,6-Tribromophenol (Surr)	100 ug/mL
							2-Fluorobiphenyl	100 ug/mL
							2-Fluorophenol (Surr)	100 ug/mL
							Nitrobenzene-d5 (Surr)	100 ug/mL
							Phenol-d5 (Surr)	100 ug/mL
							Phenol-d6	100 ug/mL
							Terphenyl-d14 (Surr)	100 ug/mL
					8270SurStkHL_00103	5 mL	2,4,6 - Tribromophenol	100 ug/mL
							2,4,6-Tribromophenol (Surr)	100 ug/mL
							2-Fluorobiphenyl	100 ug/mL
							2-Fluorophenol (Surr)	100 ug/mL
							Nitrobenzene-d5 (Surr)	100 ug/mL
							Phenol-d5 (Surr)	100 ug/mL
							Phenol-d6	100 ug/mL
							Terphenyl-d14 (Surr)	100 ug/mL
					8270SurStkHL_00104	5 mL	2,4,6 - Tribromophenol	100 ug/mL
							2,4,6-Tribromophenol (Surr)	100 ug/mL
							2-Fluorobiphenyl	100 ug/mL
							2-Fluorophenol (Surr)	100 ug/mL
							Nitrobenzene-d5 (Surr)	100 ug/mL
							Phenol-d5 (Surr)	100 ug/mL
							Phenol-d6	100 ug/mL
							Terphenyl-d14 (Surr)	100 ug/mL
					8270SurStkHL_00105	5 mL	2,4,6 - Tribromophenol	100 ug/mL
							2,4,6-Tribromophenol (Surr)	100 ug/mL
							2-Fluorobiphenyl	100 ug/mL
							2-Fluorophenol (Surr)	100 ug/mL
							Nitrobenzene-d5 (Surr)	100 ug/mL
							Phenol-d5 (Surr)	100 ug/mL
							Phenol-d6	100 ug/mL
							Terphenyl-d14 (Surr)	100 ug/mL
.8270SurStkHL_00078	02/23/18		Restek, Lot A093638		(Purchased Reagent)		2,4,6 - Tribromophenol	5000 ug/mL
							2,4,6-Tribromophenol (Surr)	5000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Phenol-d6	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
.8270SurStkHL_00103	05/31/19		Restek, Lot A0103615		(Purchased Reagent)		2,4,6 - Tribromophenol	5000 ug/mL
							2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Phenol-d6	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
.8270SurStkHL_00104	05/31/19		Restek, Lot A0103615		(Purchased Reagent)		2,4,6 - Tribromophenol	5000 ug/mL
							2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Phenol-d6	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
.8270SurStkHL_00105	05/31/19		Restek, Lot A0103615		(Purchased Reagent)		2,4,6 - Tribromophenol	5000 ug/mL
							2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Phenol-d6	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
MS-HSLA004_00015	04/30/15	02/13/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	10 uL	2,4,6-Tribromophenol (Surr)	4 ug/mL
							2-Fluorobiphenyl	4 ug/mL
							2-Fluorophenol (Surr)	4 ug/mL
							Nitrobenzene-d5 (Surr)	4 ug/mL
							Phenol-d5 (Surr)	4 ug/mL
							Terphenyl-d14 (Surr)	4 ug/mL
							1,1'-Biphenyl	4 ug/mL
							1,2,4,5-Tetrachlorobenzene	4 ug/mL
							1,2,4-Trichlorobenzene	4 ug/mL
							1,2-Dichlorobenzene	4 ug/mL
							1,2-Diphenylhydrazine	4.0439 ug/mL
							1,3-Dichlorobenzene	4 ug/mL
							1,3-Dinitrobenzene	4 ug/mL
							1,4-Dichlorobenzene	4 ug/mL
							1,4-Dioxane	4 ug/mL
							1-Methylnaphthalene	4 ug/mL
							2,2'-oxybis[1-chloropropane]	4 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2,3,4,6-Tetrachlorophenol	4 ug/mL
							2,4,5-Trichlorophenol	4 ug/mL
							2,4,6-Trichlorophenol	4 ug/mL
							2,4-Dichlorophenol	4 ug/mL
							2,4-Dimethylphenol	4 ug/mL
							2,4-Dinitrophenol	8 ug/mL
							2,4-Dinitrotoluene	4 ug/mL
							2,6-Dinitrotoluene	4 ug/mL
							2-Chloronaphthalene	4 ug/mL
							2-Chlorophenol	4 ug/mL
							2-Methylnaphthalene	4 ug/mL
							2-Methylphenol	4 ug/mL
							2-Nitroaniline	4 ug/mL
							2-Nitrophenol	4 ug/mL
							3 & 4 Methylphenol	4 ug/mL
							3-Methylphenol	4 ug/mL
							3-Nitroaniline	4 ug/mL
							4,6-Dinitro-2-methylphenol	8 ug/mL
							4-Bromophenyl phenyl ether	4 ug/mL
							4-Chloro-3-methylphenol	4 ug/mL
							4-Chloroaniline	4 ug/mL
							4-Chlorophenyl phenyl ether	4 ug/mL
							4-Methylphenol	4 ug/mL
							4-Nitroaniline	4 ug/mL
							4-Nitrophenol	8 ug/mL
							Acenaphthene	4 ug/mL
							Acenaphthylene	4 ug/mL
							Acetophenone	4 ug/mL
							Aniline	4 ug/mL
							Anthracene	4 ug/mL
							Azobenzene	4 ug/mL
							Benzo[a]anthracene	4 ug/mL
							Benzo[a]pyrene	4 ug/mL
							Benzo[b]fluoranthene	4 ug/mL
							Benzo[g,h,i]perylene	4 ug/mL
							Benzo[k]fluoranthene	4 ug/mL
							Benzyl alcohol	4 ug/mL
							Bis(2-chloroethoxy)methane	4 ug/mL
							Bis(2-chloroethyl)ether	4 ug/mL
							Bis(2-ethylhexyl) phthalate	4 ug/mL
							Butyl benzyl phthalate	4 ug/mL
							Carbazole	4 ug/mL
							Chrysene	4 ug/mL
							Di-n-butyl phthalate	4 ug/mL
							Di-n-octyl phthalate	4 ug/mL
							Dibenz(a,h)anthracene	4 ug/mL
							Dibenzofuran	4 ug/mL
							Diethyl phthalate	4 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Dimethyl phthalate	4 ug/mL
							Fluoranthene	4 ug/mL
							Fluorene	4 ug/mL
							Hexachlorobenzene	4 ug/mL
							Hexachlorobutadiene	4 ug/mL
							Hexachlorocyclopentadiene	4 ug/mL
							Hexachloroethane	4 ug/mL
							Indeno[1,2,3-cd]pyrene	4 ug/mL
							Isophorone	4 ug/mL
							N-Nitrosodi-n-propylamine	4 ug/mL
							N-Nitrosodimethylamine	4 ug/mL
							Naphthalene	4 ug/mL
							Nitrobenzene	4 ug/mL
							Pentachlorophenol	8 ug/mL
							Phenanthrene	4 ug/mL
							Phenol	4 ug/mL
							Pyrene	4 ug/mL
							Pyridine	4 ug/mL
							3,3'-Dichlorobenzidine	4 ug/mL
							Caprolactam	4 ug/mL
							Benzoic acid	8 ug/mL
							Famphur	4 ug/mL
							N-Nitrosodiphenylamine	4 ug/mL
					MS-IS_00007	50 uL	1,4-Dichlorobenzene-d4	40 ug/mL
							Acenaphthene-d10	40 ug/mL
							Chrysene-d12	40 ug/mL
							Naphthalene-d8	40 ug/mL
							Perylene-d12	40 ug/mL
							Phenanthrene-d10	40 ug/mL
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL
							Benzo[a]pyrene	200 ug/mL
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
					MS-567673_00096	1 mL	3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
					MS-567674_00038	2 mL	Benzoic acid	400 ug/mL
					MS-568023_00008	1 mL	Famphur	200 ug/mL
					MS-568038_00008	1 mL	N-Nitrosodiphenylamine	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Caprolactam	2000 ug/mL
..MS-567674_00038	02/29/16		Restek, Lot A093441		(Purchased Reagent)		Benzoic acid	2000 ug/mL
..MS-568023_00008	02/29/16		Restek, Lot A0101191		(Purchased Reagent)		Famphur	2000 ug/mL
..MS-568038_00008	02/28/17		Restek, Lot A0101189		(Purchased Reagent)		N-Nitrosodiphenylamine	2000 ug/mL
.MS-IS_00007	12/08/15	12/08/14	Methylene Chloride, Lot 71006	250 mL	MS-567684_00016	35 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
					MS-567684_00017	15 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
..MS-567684_00016	02/28/18		Restek, Lot A093676		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
..MS-567684_00017	12/31/17		Restek, Lot A092546		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
MS-HSLA010_00015	04/30/15	02/13/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	25 uL	2,4,6-Tribromophenol (Surr)	10 ug/mL
							2-Fluorobiphenyl	10 ug/mL
							2-Fluorophenol (Surr)	10 ug/mL
							Nitrobenzene-d5 (Surr)	10 ug/mL
							Phenol-d5 (Surr)	10 ug/mL
							Terphenyl-d14 (Surr)	10 ug/mL
							1,1'-Biphenyl	10 ug/mL
							1,2,4,5-Tetrachlorobenzene	10 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,2,4-Trichlorobenzene	10 ug/mL
							1,2-Dichlorobenzene	10 ug/mL
							1,2-Diphenylhydrazine	10.1097 ug/mL
							1,3-Dichlorobenzene	10 ug/mL
							1,3-Dinitrobenzene	10 ug/mL
							1,4-Dichlorobenzene	10 ug/mL
							1,4-Dioxane	10 ug/mL
							1-Methylnaphthalene	10 ug/mL
							2,2'-oxybis[1-chloropropane]	10 ug/mL
							2,3,4,6-Tetrachlorophenol	10 ug/mL
							2,4,5-Trichlorophenol	10 ug/mL
							2,4,6-Trichlorophenol	10 ug/mL
							2,4-Dichlorophenol	10 ug/mL
							2,4-Dimethylphenol	10 ug/mL
							2,4-Dinitrophenol	20 ug/mL
							2,4-Dinitrotoluene	10 ug/mL
							2,6-Dinitrotoluene	10 ug/mL
							2-Chloronaphthalene	10 ug/mL
							2-Chlorophenol	10 ug/mL
							2-Methylnaphthalene	10 ug/mL
							2-Methylphenol	10 ug/mL
							2-Nitroaniline	10 ug/mL
							2-Nitrophenol	10 ug/mL
							3 & 4 Methylphenol	10 ug/mL
							3-Methylphenol	10 ug/mL
							3-Nitroaniline	10 ug/mL
							4,6-Dinitro-2-methylphenol	20 ug/mL
							4-Bromophenyl phenyl ether	10 ug/mL
							4-Chloro-3-methylphenol	10 ug/mL
							4-Chloroaniline	10 ug/mL
							4-Chlorophenyl phenyl ether	10 ug/mL
							4-Methylphenol	10 ug/mL
							4-Nitroaniline	10 ug/mL
							4-Nitrophenol	20 ug/mL
							Acenaphthene	10 ug/mL
							Acenaphthylene	10 ug/mL
							Acetophenone	10 ug/mL
							Aniline	10 ug/mL
							Anthracene	10 ug/mL
							Azobenzene	10 ug/mL
							Benzo[a]anthracene	10 ug/mL
							Benzo[a]pyrene	10 ug/mL
							Benzo[b]fluoranthene	10 ug/mL
							Benzo[g,h,i]perylene	10 ug/mL
							Benzo[k]fluoranthene	10 ug/mL
							Benzyl alcohol	10 ug/mL
							Bis(2-chloroethoxy)methane	10 ug/mL
							Bis(2-chloroethyl)ether	10 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Bis(2-ethylhexyl) phthalate	10 ug/mL
							Butyl benzyl phthalate	10 ug/mL
							Carbazole	10 ug/mL
							Chrysene	10 ug/mL
							Di-n-butyl phthalate	10 ug/mL
							Di-n-octyl phthalate	10 ug/mL
							Dibenz(a,h)anthracene	10 ug/mL
							Dibenzofuran	10 ug/mL
							Diethyl phthalate	10 ug/mL
							Dimethyl phthalate	10 ug/mL
							Fluoranthene	10 ug/mL
							Fluorene	10 ug/mL
							Hexachlorobenzene	10 ug/mL
							Hexachlorobutadiene	10 ug/mL
							Hexachlorocyclopentadiene	10 ug/mL
							Hexachloroethane	10 ug/mL
							Indeno[1,2,3-cd]pyrene	10 ug/mL
							Isophorone	10 ug/mL
							N-Nitrosodi-n-propylamine	10 ug/mL
							N-Nitrosodimethylamine	10 ug/mL
							Naphthalene	10 ug/mL
							Nitrobenzene	10 ug/mL
							Pentachlorophenol	20 ug/mL
							Phenanthrene	10 ug/mL
							Phenol	10 ug/mL
							Pyrene	10 ug/mL
							Pyridine	10 ug/mL
							3,3'-Dichlorobenzidine	10 ug/mL
							Caprolactam	10 ug/mL
							Benzoic acid	20 ug/mL
							Famphur	10 ug/mL
							N-Nitrosodiphenylamine	10 ug/mL
					MS-IS_00007	50 uL	1,4-Dichlorobenzene-d4	40 ug/mL
							Acenaphthene-d10	40 ug/mL
							Chrysene-d12	40 ug/mL
							Naphthalene-d8	40 ug/mL
							Perylene-d12	40 ug/mL
							Phenanthrene-d10	40 ug/mL
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL
							Benzo[a]pyrene	200 ug/mL
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
					MS-567673_00096	1 mL	3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
					MS-567674_00038	2 mL	Benzoic acid	400 ug/mL
					MS-568023_00008	1 mL	Famphur	200 ug/mL
					MS-568038_00008	1 mL	N-Nitrosodiphenylamine	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added	Dimethyl phthalate	1000 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Caprolactam	2000 ug/mL
..MS-567674_00038	02/29/16		Restek, Lot A093441		(Purchased Reagent)		Benzoic acid	2000 ug/mL
..MS-568023_00008	02/29/16		Restek, Lot A0101191		(Purchased Reagent)		Famphur	2000 ug/mL
..MS-568038_00008	02/28/17		Restek, Lot A0101189		(Purchased Reagent)		N-Nitrosodiphenylamine	2000 ug/mL
.MS-IS_00007	12/08/15	12/08/14	Methylene Chloride, Lot 71006	250 mL	MS-567684_00016	35 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
					MS-567684_00017	15 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
..MS-567684_00016	02/28/18		Restek, Lot A093676		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
..MS-567684_00017	12/31/17		Restek, Lot A092546		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
MS-HSLA020_00015	04/30/15	02/13/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	50 uL	2,4,6-Tribromophenol (Surr)	20 ug/mL
							2-Fluorobiphenyl	20 ug/mL
							2-Fluorophenol (Surr)	20 ug/mL
							Nitrobenzene-d5 (Surr)	20 ug/mL
							Phenol-d5 (Surr)	20 ug/mL
							Terphenyl-d14 (Surr)	20 ug/mL
							1,1'-Biphenyl	20 ug/mL
							1,2,4,5-Tetrachlorobenzene	20 ug/mL
							1,2,4-Trichlorobenzene	20 ug/mL
							1,2-Dichlorobenzene	20 ug/mL
							1,2-Diphenylhydrazine	20.2195 ug/mL
							1,3-Dichlorobenzene	20 ug/mL
							1,3-Dinitrobenzene	20 ug/mL
							1,4-Dichlorobenzene	20 ug/mL
							1,4-Dioxane	20 ug/mL
							1-Methylnaphthalene	20 ug/mL
							2,2'-oxybis[1-chloropropane]	20 ug/mL
							2,3,4,6-Tetrachlorophenol	20 ug/mL
							2,4,5-Trichlorophenol	20 ug/mL
							2,4,6-Trichlorophenol	20 ug/mL
							2,4-Dichlorophenol	20 ug/mL
							2,4-Dimethylphenol	20 ug/mL
							2,4-Dinitrophenol	40 ug/mL
							2,4-Dinitrotoluene	20 ug/mL
							2,6-Dinitrotoluene	20 ug/mL
							2-Chloronaphthalene	20 ug/mL
							2-Chlorophenol	20 ug/mL
							2-Methylnaphthalene	20 ug/mL
							2-Methylphenol	20 ug/mL
							2-Nitroaniline	20 ug/mL
							2-Nitrophenol	20 ug/mL
							3 & 4 Methylphenol	20 ug/mL
							3-Methylphenol	20 ug/mL
							3-Nitroaniline	20 ug/mL
							4,6-Dinitro-2-methylphenol	40 ug/mL
							4-Bromophenyl phenyl ether	20 ug/mL
							4-Chloro-3-methylphenol	20 ug/mL
							4-Chloroaniline	20 ug/mL
							4-Chlorophenyl phenyl ether	20 ug/mL
							4-Methylphenol	20 ug/mL
							4-Nitroaniline	20 ug/mL
							4-Nitrophenol	40 ug/mL
							Acenaphthene	20 ug/mL
							Acenaphthylene	20 ug/mL
							Acetophenone	20 ug/mL
							Aniline	20 ug/mL
							Anthracene	20 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Azobenzene	20 ug/mL
							Benzo[a]anthracene	20 ug/mL
							Benzo[a]pyrene	20 ug/mL
							Benzo[b]fluoranthene	20 ug/mL
							Benzo[g,h,i]perylene	20 ug/mL
							Benzo[k]fluoranthene	20 ug/mL
							Benzyl alcohol	20 ug/mL
							Bis(2-chloroethoxy)methane	20 ug/mL
							Bis(2-chloroethyl)ether	20 ug/mL
							Bis(2-ethylhexyl) phthalate	20 ug/mL
							Butyl benzyl phthalate	20 ug/mL
							Carbazole	20 ug/mL
							Chrysene	20 ug/mL
							Di-n-butyl phthalate	20 ug/mL
							Di-n-octyl phthalate	20 ug/mL
							Dibenz(a,h)anthracene	20 ug/mL
							Dibenzofuran	20 ug/mL
							Diethyl phthalate	20 ug/mL
							Dimethyl phthalate	20 ug/mL
							Fluoranthene	20 ug/mL
							Fluorene	20 ug/mL
							Hexachlorobenzene	20 ug/mL
							Hexachlorobutadiene	20 ug/mL
							Hexachlorocyclopentadiene	20 ug/mL
							Hexachloroethane	20 ug/mL
							Indeno[1,2,3-cd]pyrene	20 ug/mL
							Isophorone	20 ug/mL
							N-Nitrosodi-n-propylamine	20 ug/mL
							N-Nitrosodimethylamine	20 ug/mL
							Naphthalene	20 ug/mL
							Nitrobenzene	20 ug/mL
							Pentachlorophenol	40 ug/mL
							Phenanthrene	20 ug/mL
							Phenol	20 ug/mL
							Pyrene	20 ug/mL
							Pyridine	20 ug/mL
							3,3'-Dichlorobenzidine	20 ug/mL
							Caprolactam	20 ug/mL
							Benzoic acid	40 ug/mL
							Famphur	20 ug/mL
							N-Nitrosodiphenylamine	20 ug/mL
					MS-IS_00007	50 uL	1,4-Dichlorobenzene-d4	40 ug/mL
							Acenaphthene-d10	40 ug/mL
							Chrysene-d12	40 ug/mL
							Naphthalene-d8	40 ug/mL
							Perylene-d12	40 ug/mL
							Phenanthrene-d10	40 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL
							Benzo[a]pyrene	200 ug/mL
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
					MS-567673_00096	1 mL	3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
					MS-567674_00038	2 mL	Benzoic acid	400 ug/mL
					MS-568023_00008	1 mL	Famphur	200 ug/mL
					MS-568038_00008	1 mL	N-Nitrosodiphenylamine	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added	Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Caprolactam	2000 ug/mL
..MS-567674_00038	02/29/16		Restek, Lot A093441		(Purchased Reagent)		Benzoic acid	2000 ug/mL
..MS-568023_00008	02/29/16		Restek, Lot A0101191		(Purchased Reagent)		Famphur	2000 ug/mL
..MS-568038_00008	02/28/17		Restek, Lot A0101189		(Purchased Reagent)		N-Nitrosodiphenylamine	2000 ug/mL
.MS-IS_00007	12/08/15	12/08/14	Methylene Chloride, Lot 71006	250 mL	MS-567684_00016	35 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
					MS-567684_00017	15 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
..MS-567684_00016	02/28/18		Restek, Lot A093676		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
..MS-567684_00017	12/31/17		Restek, Lot A092546		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
MS-HSLA050_00015	04/30/15	02/13/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	125 uL	2,4,6-Tribromophenol (Surr)	50 ug/mL
							2-Fluorobiphenyl	50 ug/mL
							2-Fluorophenol (Surr)	50 ug/mL
							Nitrobenzene-d5 (Surr)	50 ug/mL
							Phenol-d5 (Surr)	50 ug/mL
							Terphenyl-d14 (Surr)	50 ug/mL
							1,1'-Biphenyl	50 ug/mL
							1,2,4,5-Tetrachlorobenzene	50 ug/mL
							1,2,4-Trichlorobenzene	50 ug/mL
							1,2-Dichlorobenzene	50 ug/mL
							1,2-Diphenylhydrazine	50.5487 ug/mL
							1,3-Dichlorobenzene	50 ug/mL
							1,3-Dinitrobenzene	50 ug/mL
							1,4-Dichlorobenzene	50 ug/mL
							1,4-Dioxane	50 ug/mL
							1-Methylnaphthalene	50 ug/mL
							2,2'-oxybis[1-chloropropane]	50 ug/mL
							2,3,4,6-Tetrachlorophenol	50 ug/mL
							2,4,5-Trichlorophenol	50 ug/mL
							2,4,6-Trichlorophenol	50 ug/mL
							2,4-Dichlorophenol	50 ug/mL
							2,4-Dimethylphenol	50 ug/mL
							2,4-Dinitrophenol	100 ug/mL
							2,4-Dinitrotoluene	50 ug/mL
							2,6-Dinitrotoluene	50 ug/mL
							2-Chloronaphthalene	50 ug/mL
							2-Chlorophenol	50 ug/mL
							2-Methylnaphthalene	50 ug/mL
							2-Methylphenol	50 ug/mL
							2-Nitroaniline	50 ug/mL
							2-Nitrophenol	50 ug/mL
							3 & 4 Methylphenol	50 ug/mL
							3-Methylphenol	50 ug/mL
							3-Nitroaniline	50 ug/mL
							4,6-Dinitro-2-methylphenol	100 ug/mL
							4-Bromophenyl phenyl ether	50 ug/mL
							4-Chloro-3-methylphenol	50 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							4-Chloroaniline	50 ug/mL
							4-Chlorophenyl phenyl ether	50 ug/mL
							4-Methylphenol	50 ug/mL
							4-Nitroaniline	50 ug/mL
							4-Nitrophenol	100 ug/mL
							Acenaphthene	50 ug/mL
							Acenaphthylene	50 ug/mL
							Acetophenone	50 ug/mL
							Aniline	50 ug/mL
							Anthracene	50 ug/mL
							Azobenzene	50 ug/mL
							Benzo[a]anthracene	50 ug/mL
							Benzo[a]pyrene	50 ug/mL
							Benzo[b]fluoranthene	50 ug/mL
							Benzo[g,h,i]perylene	50 ug/mL
							Benzo[k]fluoranthene	50 ug/mL
							Benzyl alcohol	50 ug/mL
							Bis(2-chloroethoxy)methane	50 ug/mL
							Bis(2-chloroethyl)ether	50 ug/mL
							Bis(2-ethylhexyl) phthalate	50 ug/mL
							Butyl benzyl phthalate	50 ug/mL
							Carbazole	50 ug/mL
							Chrysene	50 ug/mL
							Di-n-butyl phthalate	50 ug/mL
							Di-n-octyl phthalate	50 ug/mL
							Dibenz(a,h)anthracene	50 ug/mL
							Dibenzofuran	50 ug/mL
							Diethyl phthalate	50 ug/mL
							Dimethyl phthalate	50 ug/mL
							Fluoranthene	50 ug/mL
							Fluorene	50 ug/mL
							Hexachlorobenzene	50 ug/mL
							Hexachlorobutadiene	50 ug/mL
							Hexachlorocyclopentadiene	50 ug/mL
							Hexachloroethane	50 ug/mL
							Indeno[1,2,3-cd]pyrene	50 ug/mL
							Isophorone	50 ug/mL
							N-Nitrosodi-n-propylamine	50 ug/mL
							N-Nitrosodimethylamine	50 ug/mL
							Naphthalene	50 ug/mL
							Nitrobenzene	50 ug/mL
							Pentachlorophenol	100 ug/mL
							Phenanthrene	50 ug/mL
							Phenol	50 ug/mL
							Pyrene	50 ug/mL
							Pyridine	50 ug/mL
							3,3'-Dichlorobenzidine	50 ug/mL
							Caprolactam	50 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Benzoic acid	100 ug/mL
							Famphur	50 ug/mL
							N-Nitrosodiphenylamine	50 ug/mL
					MS-IS_00007	50 uL	1,4-Dichlorobenzene-d4	40 ug/mL
							Acenaphthene-d10	40 ug/mL
							Chrysene-d12	40 ug/mL
							Naphthalene-d8	40 ug/mL
							Perylene-d12	40 ug/mL
							Phenanthrene-d10	40 ug/mL
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL
							Benzo[a]pyrene	200 ug/mL
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
					MS-567673_00096	1 mL	3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
					MS-567674_00038	2 mL	Benzoic acid	400 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
					MS-568023_00008	1 mL	Famphur	200 ug/mL
					MS-568038_00008	1 mL	N-Nitrosodiphenylamine	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Caprolactam	2000 ug/mL
..MS-567674_00038	02/29/16		Restek, Lot A093441		(Purchased Reagent)		Benzoic acid	2000 ug/mL
..MS-568023_00008	02/29/16		Restek, Lot A0101191		(Purchased Reagent)		Famphur	2000 ug/mL
..MS-568038_00008	02/28/17		Restek, Lot A0101189		(Purchased Reagent)		N-Nitrosodiphenylamine	2000 ug/mL
.MS-IS_00007	12/08/15	12/08/14	Methylene Chloride, Lot 71006	250 mL	MS-567684_00016	35 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Phenanthrene-d10	400 ug/mL
					MS-567684_00017	15 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
..MS-567684_00016	02/28/18		Restek, Lot A093676		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
..MS-567684_00017	12/31/17		Restek, Lot A092546		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
MS-HSLA080_00015	04/30/15	02/13/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	200 uL	2,4,6-Tribromophenol (Surr)	80 ug/mL
							2-Fluorobiphenyl	80 ug/mL
							2-Fluorophenol (Surr)	80 ug/mL
							Nitrobenzene-d5 (Surr)	80 ug/mL
							Phenol-d5 (Surr)	80 ug/mL
							Terphenyl-d14 (Surr)	80 ug/mL
							1,1'-Biphenyl	80 ug/mL
							1,2,4,5-Tetrachlorobenzene	80 ug/mL
							1,2,4-Trichlorobenzene	80 ug/mL
							1,2-Dichlorobenzene	80 ug/mL
							1,2-Diphenylhydrazine	80.878 ug/mL
							1,3-Dichlorobenzene	80 ug/mL
							1,3-Dinitrobenzene	80 ug/mL
							1,4-Dichlorobenzene	80 ug/mL
							1,4-Dioxane	80 ug/mL
							1-Methylnaphthalene	80 ug/mL
							2,2'-oxybis[1-chloropropane]	80 ug/mL
							2,3,4,6-Tetrachlorophenol	80 ug/mL
							2,4,5-Trichlorophenol	80 ug/mL
							2,4,6-Trichlorophenol	80 ug/mL
							2,4-Dichlorophenol	80 ug/mL
							2,4-Dimethylphenol	80 ug/mL
							2,4-Dinitrophenol	160 ug/mL
							2,4-Dinitrotoluene	80 ug/mL
							2,6-Dinitrotoluene	80 ug/mL
							2-Chloronaphthalene	80 ug/mL
							2-Chlorophenol	80 ug/mL
							2-Methylnaphthalene	80 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Methylphenol	80 ug/mL
							2-Nitroaniline	80 ug/mL
							2-Nitrophenol	80 ug/mL
							3 & 4 Methylphenol	80 ug/mL
							3-Methylphenol	80 ug/mL
							3-Nitroaniline	80 ug/mL
							4,6-Dinitro-2-methylphenol	160 ug/mL
							4-Bromophenyl phenyl ether	80 ug/mL
							4-Chloro-3-methylphenol	80 ug/mL
							4-Chloroaniline	80 ug/mL
							4-Chlorophenyl phenyl ether	80 ug/mL
							4-Methylphenol	80 ug/mL
							4-Nitroaniline	80 ug/mL
							4-Nitrophenol	160 ug/mL
							Acenaphthene	80 ug/mL
							Acenaphthylene	80 ug/mL
							Acetophenone	80 ug/mL
							Aniline	80 ug/mL
							Anthracene	80 ug/mL
							Azobenzene	80 ug/mL
							Benzo[a]anthracene	80 ug/mL
							Benzo[a]pyrene	80 ug/mL
							Benzo[b]fluoranthene	80 ug/mL
							Benzo[g,h,i]perylene	80 ug/mL
							Benzo[k]fluoranthene	80 ug/mL
							Benzyl alcohol	80 ug/mL
							Bis(2-chloroethoxy)methane	80 ug/mL
							Bis(2-chloroethyl)ether	80 ug/mL
							Bis(2-ethylhexyl) phthalate	80 ug/mL
							Butyl benzyl phthalate	80 ug/mL
							Carbazole	80 ug/mL
							Chrysene	80 ug/mL
							Di-n-butyl phthalate	80 ug/mL
							Di-n-octyl phthalate	80 ug/mL
							Dibenz(a,h)anthracene	80 ug/mL
							Dibenzofuran	80 ug/mL
							Diethyl phthalate	80 ug/mL
							Dimethyl phthalate	80 ug/mL
							Fluoranthene	80 ug/mL
							Fluorene	80 ug/mL
							Hexachlorobenzene	80 ug/mL
							Hexachlorobutadiene	80 ug/mL
							Hexachlorocyclopentadiene	80 ug/mL
							Hexachloroethane	80 ug/mL
							Indeno[1,2,3-cd]pyrene	80 ug/mL
							Isophorone	80 ug/mL
							N-Nitrosodi-n-propylamine	80 ug/mL
							N-Nitrosodimethylamine	80 ug/mL

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SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Naphthalene	80 ug/mL
							Nitrobenzene	80 ug/mL
							Pentachlorophenol	160 ug/mL
							Phenanthrene	80 ug/mL
							Phenol	80 ug/mL
							Pyrene	80 ug/mL
							Pyridine	80 ug/mL
							3,3'-Dichlorobenzidine	80 ug/mL
							Caprolactam	80 ug/mL
							Benzoic acid	160 ug/mL
							Famphur	80 ug/mL
							N-Nitrosodiphenylamine	80 ug/mL
					MS-IS_00007	50 uL	1,4-Dichlorobenzene-d4	40 ug/mL
							Acenaphthene-d10	40 ug/mL
							Chrysene-d12	40 ug/mL
							Naphthalene-d8	40 ug/mL
							Perylene-d12	40 ug/mL
							Phenanthrene-d10	40 ug/mL
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL
							Benzo[a]pyrene	200 ug/mL
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
					MS-567673_00096	1 mL	3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
					MS-567674_00038	2 mL	Benzoic acid	400 ug/mL
					MS-568023_00008	1 mL	Famphur	200 ug/mL
					MS-568038_00008	1 mL	N-Nitrosodiphenylamine	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Caprolactam	2000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
..MS-567674_00038	02/29/16		Restek, Lot A093441		(Purchased Reagent)		Benzoic acid	2000 ug/mL
..MS-568023_00008	02/29/16		Restek, Lot A0101191		(Purchased Reagent)		Famphur	2000 ug/mL
..MS-568038_00008	02/28/17		Restek, Lot A0101189		(Purchased Reagent)		N-Nitrosodiphenylamine	2000 ug/mL
.MS-IS_00007	12/08/15	12/08/14	Methylene Chloride, Lot 71006	250 mL	MS-567684_00016	35 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
					MS-567684_00017	15 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
..MS-567684_00016	02/28/18		Restek, Lot A093676		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
..MS-567684_00017	12/31/17		Restek, Lot A092546		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
MS-HSLA120_00015	04/30/15	02/13/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	300 uL	2,4,6-Tribromophenol (Surr)	120 ug/mL
							2-Fluorobiphenyl	120 ug/mL
							2-Fluorophenol (Surr)	120 ug/mL
							Nitrobenzene-d5 (Surr)	120 ug/mL
							Phenol-d5 (Surr)	120 ug/mL
							Terphenyl-d14 (Surr)	120 ug/mL
							1,1'-Biphenyl	120 ug/mL
							1,2,4,5-Tetrachlorobenzene	120 ug/mL
							1,2,4-Trichlorobenzene	120 ug/mL
							1,2-Dichlorobenzene	120 ug/mL
							1,2-Diphenylhydrazine	121.317 ug/mL
							1,3-Dichlorobenzene	120 ug/mL
							1,3-Dinitrobenzene	120 ug/mL
							1,4-Dichlorobenzene	120 ug/mL
							1,4-Dioxane	120 ug/mL
							1-Methylnaphthalene	120 ug/mL
							2,2'-oxybis[1-chloropropane]	120 ug/mL
							2,3,4,6-Tetrachlorophenol	120 ug/mL
							2,4,5-Trichlorophenol	120 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2,4,6-Trichlorophenol	120 ug/mL
							2,4-Dichlorophenol	120 ug/mL
							2,4-Dimethylphenol	120 ug/mL
							2,4-Dinitrophenol	240 ug/mL
							2,4-Dinitrotoluene	120 ug/mL
							2,6-Dinitrotoluene	120 ug/mL
							2-Chloronaphthalene	120 ug/mL
							2-Chlorophenol	120 ug/mL
							2-Methylnaphthalene	120 ug/mL
							2-Methylphenol	120 ug/mL
							2-Nitroaniline	120 ug/mL
							2-Nitrophenol	120 ug/mL
							3 & 4 Methylphenol	120 ug/mL
							3-Methylphenol	120 ug/mL
							3-Nitroaniline	120 ug/mL
							4,6-Dinitro-2-methylphenol	240 ug/mL
							4-Bromophenyl phenyl ether	120 ug/mL
							4-Chloro-3-methylphenol	120 ug/mL
							4-Chloroaniline	120 ug/mL
							4-Chlorophenyl phenyl ether	120 ug/mL
							4-Methylphenol	120 ug/mL
							4-Nitroaniline	120 ug/mL
							4-Nitrophenol	240 ug/mL
							Acenaphthene	120 ug/mL
							Acenaphthylene	120 ug/mL
							Acetophenone	120 ug/mL
							Aniline	120 ug/mL
							Anthracene	120 ug/mL
							Azobenzene	120 ug/mL
							Benzo[a]anthracene	120 ug/mL
							Benzo[a]pyrene	120 ug/mL
							Benzo[b]fluoranthene	120 ug/mL
							Benzo[g,h,i]perylene	120 ug/mL
							Benzo[k]fluoranthene	120 ug/mL
							Benzyl alcohol	120 ug/mL
							Bis(2-chloroethoxy)methane	120 ug/mL
							Bis(2-chloroethyl)ether	120 ug/mL
							Bis(2-ethylhexyl) phthalate	120 ug/mL
							Butyl benzyl phthalate	120 ug/mL
							Carbazole	120 ug/mL
							Chrysene	120 ug/mL
							Di-n-butyl phthalate	120 ug/mL
							Di-n-octyl phthalate	120 ug/mL
							Dibenz(a,h)anthracene	120 ug/mL
							Dibenzofuran	120 ug/mL
							Diethyl phthalate	120 ug/mL
							Dimethyl phthalate	120 ug/mL
							Fluoranthene	120 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Fluorene	120 ug/mL
							Hexachlorobenzene	120 ug/mL
							Hexachlorobutadiene	120 ug/mL
							Hexachlorocyclopentadiene	120 ug/mL
							Hexachloroethane	120 ug/mL
							Indeno[1,2,3-cd]pyrene	120 ug/mL
							Isophorone	120 ug/mL
							N-Nitrosodi-n-propylamine	120 ug/mL
							N-Nitrosodimethylamine	120 ug/mL
							Naphthalene	120 ug/mL
							Nitrobenzene	120 ug/mL
							Pentachlorophenol	240 ug/mL
							Phenanthrene	120 ug/mL
							Phenol	120 ug/mL
							Pyrene	120 ug/mL
							Pyridine	120 ug/mL
							3,3'-Dichlorobenzidine	120 ug/mL
							Caprolactam	120 ug/mL
							Benzoic acid	240 ug/mL
							Famphur	120 ug/mL
							N-Nitrosodiphenylamine	120 ug/mL
					MS-IS_00007	50 uL	1,4-Dichlorobenzene-d4	40 ug/mL
							Acenaphthene-d10	40 ug/mL
							Chrysene-d12	40 ug/mL
							Naphthalene-d8	40 ug/mL
							Perylene-d12	40 ug/mL
							Phenanthrene-d10	40 ug/mL
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL
							Benzo[a]pyrene	200 ug/mL
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
					MS-567673_00096	1 mL	3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
					MS-567674_00038	2 mL	Benzoic acid	400 ug/mL
					MS-568023_00008	1 mL	Famphur	200 ug/mL
					MS-568038_00008	1 mL	N-Nitrosodiphenylamine	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Caprolactam	2000 ug/mL
..MS-567674_00038	02/29/16		Restek, Lot A093441		(Purchased Reagent)		Benzoic acid	2000 ug/mL
..MS-568023_00008	02/29/16		Restek, Lot A0101191		(Purchased Reagent)		Famphur	2000 ug/mL
..MS-568038_00008	02/28/17		Restek, Lot A0101189		(Purchased Reagent)		N-Nitrosodiphenylamine	2000 ug/mL
.MS-IS_00007	12/08/15	12/08/14	Methylene Chloride, Lot 71006	250 mL	MS-567684_00016	35 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
					MS-567684_00017	15 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
..MS-567684_00016	02/28/18		Restek, Lot A093676		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
..MS-567684_00017	12/31/17		Restek, Lot A092546		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
MS-HSLA160_00015	04/30/15	02/13/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	400 uL	2,4,6-Tribromophenol (Surr)	160 ug/mL
							2-Fluorobiphenyl	160 ug/mL
							2-Fluorophenol (Surr)	160 ug/mL
							Nitrobenzene-d5 (Surr)	160 ug/mL
							Phenol-d5 (Surr)	160 ug/mL
							Terphenyl-d14 (Surr)	160 ug/mL
							1,1'-Biphenyl	160 ug/mL
							1,2,4,5-Tetrachlorobenzene	160 ug/mL
							1,2,4-Trichlorobenzene	160 ug/mL
							1,2-Dichlorobenzene	160 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,2-Diphenylhydrazine	161.756 ug/mL
							1,3-Dichlorobenzene	160 ug/mL
							1,3-Dinitrobenzene	160 ug/mL
							1,4-Dichlorobenzene	160 ug/mL
							1,4-Dioxane	160 ug/mL
							1-Methylnaphthalene	160 ug/mL
							2,2'-oxybis[1-chloropropane]	160 ug/mL
							2,3,4,6-Tetrachlorophenol	160 ug/mL
							2,4,5-Trichlorophenol	160 ug/mL
							2,4,6-Trichlorophenol	160 ug/mL
							2,4-Dichlorophenol	160 ug/mL
							2,4-Dimethylphenol	160 ug/mL
							2,4-Dinitrophenol	320 ug/mL
							2,4-Dinitrotoluene	160 ug/mL
							2,6-Dinitrotoluene	160 ug/mL
							2-Chloronaphthalene	160 ug/mL
							2-Chlorophenol	160 ug/mL
							2-Methylnaphthalene	160 ug/mL
							2-Methylphenol	160 ug/mL
							2-Nitroaniline	160 ug/mL
							2-Nitrophenol	160 ug/mL
							3 & 4 Methylphenol	160 ug/mL
							3-Methylphenol	160 ug/mL
							3-Nitroaniline	160 ug/mL
							4,6-Dinitro-2-methylphenol	320 ug/mL
							4-Bromophenyl phenyl ether	160 ug/mL
							4-Chloro-3-methylphenol	160 ug/mL
							4-Chloroaniline	160 ug/mL
							4-Chlorophenyl phenyl ether	160 ug/mL
							4-Methylphenol	160 ug/mL
							4-Nitroaniline	160 ug/mL
							4-Nitrophenol	320 ug/mL
							Acenaphthene	160 ug/mL
							Acenaphthylene	160 ug/mL
							Acetophenone	160 ug/mL
							Aniline	160 ug/mL
							Anthracene	160 ug/mL
							Azobenzene	160 ug/mL
							Benzo[a]anthracene	160 ug/mL
							Benzo[a]pyrene	160 ug/mL
							Benzo[b]fluoranthene	160 ug/mL
							Benzo[g,h,i]perylene	160 ug/mL
							Benzo[k]fluoranthene	160 ug/mL
							Benzyl alcohol	160 ug/mL
							Bis(2-chloroethoxy)methane	160 ug/mL
							Bis(2-chloroethyl)ether	160 ug/mL
							Bis(2-ethylhexyl) phthalate	160 ug/mL
							Butyl benzyl phthalate	160 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Carbazole	160 ug/mL
							Chrysene	160 ug/mL
							Di-n-butyl phthalate	160 ug/mL
							Di-n-octyl phthalate	160 ug/mL
							Dibenz(a,h)anthracene	160 ug/mL
							Dibenzofuran	160 ug/mL
							Diethyl phthalate	160 ug/mL
							Dimethyl phthalate	160 ug/mL
							Fluoranthene	160 ug/mL
							Fluorene	160 ug/mL
							Hexachlorobenzene	160 ug/mL
							Hexachlorobutadiene	160 ug/mL
							Hexachlorocyclopentadiene	160 ug/mL
							Hexachloroethane	160 ug/mL
							Indeno[1,2,3-cd]pyrene	160 ug/mL
							Isophorone	160 ug/mL
							N-Nitrosodi-n-propylamine	160 ug/mL
							N-Nitrosodimethylamine	160 ug/mL
							Naphthalene	160 ug/mL
							Nitrobenzene	160 ug/mL
							Pentachlorophenol	320 ug/mL
							Phenanthrene	160 ug/mL
							Phenol	160 ug/mL
							Pyrene	160 ug/mL
							Pyridine	160 ug/mL
							3,3'-Dichlorobenzidine	160 ug/mL
							Caprolactam	160 ug/mL
							Benzoic acid	320 ug/mL
							Famphur	160 ug/mL
							N-Nitrosodiphenylamine	160 ug/mL
					MS-IS_00007	50 uL	1,4-Dichlorobenzene-d4	40 ug/mL
							Acenaphthene-d10	40 ug/mL
							Chrysene-d12	40 ug/mL
							Naphthalene-d8	40 ug/mL
							Perylene-d12	40 ug/mL
							Phenanthrene-d10	40 ug/mL
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL
							Benzo[a]pyrene	200 ug/mL
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
					MS-567673_00096	1 mL	3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
					MS-567674_00038	2 mL	Benzoic acid	400 ug/mL
					MS-568023_00008	1 mL	Famphur	200 ug/mL
					MS-568038_00008	1 mL	N-Nitrosodiphenylamine	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Fluoranthene	1000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Caprolactam	2000 ug/mL
..MS-567674_00038	02/29/16		Restek, Lot A093441		(Purchased Reagent)		Benzoic acid	2000 ug/mL
..MS-568023_00008	02/29/16		Restek, Lot A0101191		(Purchased Reagent)		Famphur	2000 ug/mL
..MS-568038_00008	02/28/17		Restek, Lot A0101189		(Purchased Reagent)		N-Nitrosodiphenylamine	2000 ug/mL
.MS-IS_00007	12/08/15	12/08/14	Methylene Chloride, Lot 71006	250 mL	MS-567684_00016	35 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
					MS-567684_00017	15 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
..MS-567684_00016	02/28/18		Restek, Lot A093676		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
..MS-567684_00017	12/31/17		Restek, Lot A092546		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
MS-HSLA200_00015	04/30/15	02/13/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	500 uL	2,4,6-Tribromophenol (Surr)	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
							1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Benzo[a]pyrene	200 ug/mL
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
							3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
							Benzoic acid	400 ug/mL
							Famphur	200 ug/mL
							N-Nitrosodiphenylamine	200 ug/mL
					MS-IS_00007	50 uL	1,4-Dichlorobenzene-d4	40 ug/mL
							Acenaphthene-d10	40 ug/mL
							Chrysene-d12	40 ug/mL
							Naphthalene-d8	40 ug/mL
							Perylene-d12	40 ug/mL
							Phenanthrene-d10	40 ug/mL
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,1'-Biphenyl	200 ug/mL
							1,2,4,5-Tetrachlorobenzene	200 ug/mL
							1,2,4-Trichlorobenzene	200 ug/mL
							1,2-Dichlorobenzene	200 ug/mL
							1,2-Diphenylhydrazine	202.195 ug/mL
							1,3-Dichlorobenzene	200 ug/mL
							1,3-Dinitrobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							1,4-Dioxane	200 ug/mL
							1-Methylnaphthalene	200 ug/mL
							2,2'-oxybis[1-chloropropane]	200 ug/mL
							2,3,4,6-Tetrachlorophenol	200 ug/mL
							2,4,5-Trichlorophenol	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dichlorophenol	200 ug/mL
							2,4-Dimethylphenol	200 ug/mL
							2,4-Dinitrophenol	400 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2,6-Dinitrotoluene	200 ug/mL
							2-Chloronaphthalene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							2-Nitroaniline	200 ug/mL
							2-Nitrophenol	200 ug/mL
							3 & 4 Methylphenol	200 ug/mL
							3-Methylphenol	200 ug/mL
							3-Nitroaniline	200 ug/mL
							4,6-Dinitro-2-methylphenol	400 ug/mL
							4-Bromophenyl phenyl ether	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Chloroaniline	200 ug/mL
							4-Chlorophenyl phenyl ether	200 ug/mL
							4-Methylphenol	200 ug/mL
							4-Nitroaniline	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Acenaphthylene	200 ug/mL
							Acetophenone	200 ug/mL
							Aniline	200 ug/mL
							Anthracene	200 ug/mL
							Azobenzene	200 ug/mL
							Benzo[a]anthracene	200 ug/mL
							Benzo[a]pyrene	200 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Benzo[b]fluoranthene	200 ug/mL
							Benzo[g,h,i]perylene	200 ug/mL
							Benzo[k]fluoranthene	200 ug/mL
							Benzyl alcohol	200 ug/mL
							Bis(2-chloroethoxy)methane	200 ug/mL
							Bis(2-chloroethyl)ether	200 ug/mL
							Bis(2-ethylhexyl) phthalate	200 ug/mL
							Butyl benzyl phthalate	200 ug/mL
							Carbazole	200 ug/mL
							Chrysene	200 ug/mL
							Di-n-butyl phthalate	200 ug/mL
							Di-n-octyl phthalate	200 ug/mL
							Dibenz(a,h)anthracene	200 ug/mL
							Dibenzofuran	200 ug/mL
							Diethyl phthalate	200 ug/mL
							Dimethyl phthalate	200 ug/mL
							Fluoranthene	200 ug/mL
							Fluorene	200 ug/mL
							Hexachlorobenzene	200 ug/mL
							Hexachlorobutadiene	200 ug/mL
							Hexachlorocyclopentadiene	200 ug/mL
							Hexachloroethane	200 ug/mL
							Indeno[1,2,3-cd]pyrene	200 ug/mL
							Isophorone	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							N-Nitrosodimethylamine	200 ug/mL
							Naphthalene	200 ug/mL
							Nitrobenzene	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Phenanthrene	200 ug/mL
							Phenol	200 ug/mL
							Pyrene	200 ug/mL
							Pyridine	200 ug/mL
					MS-567673_00096	1 mL	3,3'-Dichlorobenzidine	200 ug/mL
							Caprolactam	200 ug/mL
					MS-567674_00038	2 mL	Benzoic acid	400 ug/mL
					MS-568023_00008	1 mL	Famphur	200 ug/mL
					MS-568038_00008	1 mL	N-Nitrosodiphenylamine	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,1'-Biphenyl	1000 ug/mL
							1,2,4,5-Tetrachlorobenzene	1000 ug/mL
							1,2,4-Trichlorobenzene	1000 ug/mL
							1,2-Dichlorobenzene	1000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							1,2-Diphenylhydrazine	1010.97 ug/mL
							1,3-Dichlorobenzene	1000 ug/mL
							1,3-Dinitrobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							1,4-Dioxane	1000 ug/mL
							1-Methylnaphthalene	1000 ug/mL
							2,2'-oxybis[1-chloropropane]	1000 ug/mL
							2,3,4,6-Tetrachlorophenol	1000 ug/mL
							2,4,5-Trichlorophenol	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dichlorophenol	1000 ug/mL
							2,4-Dimethylphenol	1000 ug/mL
							2,4-Dinitrophenol	2000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2,6-Dinitrotoluene	1000 ug/mL
							2-Chloronaphthalene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							2-Nitroaniline	1000 ug/mL
							2-Nitrophenol	1000 ug/mL
							3 & 4 Methylphenol	1000 ug/mL
							3-Methylphenol	1000 ug/mL
							3-Nitroaniline	1000 ug/mL
							4,6-Dinitro-2-methylphenol	2000 ug/mL
							4-Bromophenyl phenyl ether	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Chloroaniline	1000 ug/mL
							4-Chlorophenyl phenyl ether	1000 ug/mL
							4-Methylphenol	1000 ug/mL
							4-Nitroaniline	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Acenaphthylene	1000 ug/mL
							Acetophenone	1000 ug/mL
							Aniline	1000 ug/mL
							Anthracene	1000 ug/mL
							Azobenzene	1000 ug/mL
							Benzo[a]anthracene	1000 ug/mL
							Benzo[a]pyrene	1000 ug/mL
							Benzo[b]fluoranthene	1000 ug/mL
							Benzo[g,h,i]perylene	1000 ug/mL
							Benzo[k]fluoranthene	1000 ug/mL
							Benzyl alcohol	1000 ug/mL
							Bis(2-chloroethoxy)methane	1000 ug/mL
							Bis(2-chloroethyl)ether	1000 ug/mL
							Bis(2-ethylhexyl) phthalate	1000 ug/mL
							Butyl benzyl phthalate	1000 ug/mL

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Carbazole	1000 ug/mL
							Chrysene	1000 ug/mL
							Di-n-butyl phthalate	1000 ug/mL
							Di-n-octyl phthalate	1000 ug/mL
							Dibenz(a,h)anthracene	1000 ug/mL
							Dibenzofuran	1000 ug/mL
							Diethyl phthalate	1000 ug/mL
							Dimethyl phthalate	1000 ug/mL
							Fluoranthene	1000 ug/mL
							Fluorene	1000 ug/mL
							Hexachlorobenzene	1000 ug/mL
							Hexachlorobutadiene	1000 ug/mL
							Hexachlorocyclopentadiene	1000 ug/mL
							Hexachloroethane	1000 ug/mL
							Indeno[1,2,3-cd]pyrene	1000 ug/mL
							Isophorone	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							N-Nitrosodimethylamine	1000 ug/mL
							Naphthalene	1000 ug/mL
							Nitrobenzene	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Phenanthrene	1000 ug/mL
							Phenol	1000 ug/mL
							Pyrene	1000 ug/mL
							Pyridine	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		3,3'-Dichlorobenzidine	2000 ug/mL
							Caprolactam	2000 ug/mL
..MS-567674_00038	02/29/16		Restek, Lot A093441		(Purchased Reagent)		Benzoic acid	2000 ug/mL
..MS-568023_00008	02/29/16		Restek, Lot A0101191		(Purchased Reagent)		Famphur	2000 ug/mL
..MS-568038_00008	02/28/17		Restek, Lot A0101189		(Purchased Reagent)		N-Nitrosodiphenylamine	2000 ug/mL
.MS-IS_00007	12/08/15	12/08/14	Methylene Chloride, Lot 71006	250 mL	MS-567684_00016	35 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
					MS-567684_00017	15 mL	1,4-Dichlorobenzene-d4	400 ug/mL
							Acenaphthene-d10	400 ug/mL
							Chrysene-d12	400 ug/mL
							Naphthalene-d8	400 ug/mL
							Perylene-d12	400 ug/mL
							Phenanthrene-d10	400 ug/mL
..MS-567684_00016	02/28/18		Restek, Lot A093676		(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Phenanthrene-d10	2000 ug/mL
..MS-567684_00017	12/31/17	Restek, Lot A092546			(Purchased Reagent)		1,4-Dichlorobenzene-d4	2000 ug/mL
							Acenaphthene-d10	2000 ug/mL
							Chrysene-d12	2000 ug/mL
							Naphthalene-d8	2000 ug/mL
							Perylene-d12	2000 ug/mL
							Phenanthrene-d10	2000 ug/mL
MS-HSLACCV080_00038	04/30/15	02/10/15	Methylene Chloride, Lot 87836	0.5 mL	MS-HSLA_STK_00012	200 uL	2,4,6-Tribromophenol (Surr)	80 ug/mL
							2-Fluorobiphenyl	80 ug/mL
							2-Fluorophenol (Surr)	80 ug/mL
							Nitrobenzene-d5 (Surr)	80 ug/mL
							Phenol-d5 (Surr)	80 ug/mL
							Terphenyl-d14 (Surr)	80 ug/mL
							1,2,4-Trichlorobenzene	80 ug/mL
							1,4-Dichlorobenzene	80 ug/mL
							2,4,6-Trichlorophenol	80 ug/mL
							2,4-Dinitrotoluene	80 ug/mL
							2-Chlorophenol	80 ug/mL
							2-Methylnaphthalene	80 ug/mL
							2-Methylphenol	80 ug/mL
							4-Chloro-3-methylphenol	80 ug/mL
							4-Nitrophenol	160 ug/mL
							Acenaphthene	80 ug/mL
							Anthracene	80 ug/mL
							N-Nitrosodi-n-propylamine	80 ug/mL
							Pentachlorophenol	160 ug/mL
							Pyrene	80 ug/mL
							Caprolactam	80 ug/mL
.MS-HSLA_STK_00012	04/30/15	11/20/14	Methylene Chloride, Lot 71006	10 mL	8270SurStkHL_00002	0.4 mL	2,4,6-Tribromophenol (Surr)	200 ug/mL
							2-Fluorobiphenyl	200 ug/mL
							2-Fluorophenol (Surr)	200 ug/mL
							Nitrobenzene-d5 (Surr)	200 ug/mL
							Phenol-d5 (Surr)	200 ug/mL
							Terphenyl-d14 (Surr)	200 ug/mL
					MS-567672_00091	2 mL	1,2,4-Trichlorobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Anthracene	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							Pentachlorophenol	400 ug/mL
							Pyrene	200 ug/mL
					MS-567673_00096	1 mL	Caprolactam	200 ug/mL
..8270SurStkHL_00002	01/31/18		Restek, Lot A092712		(Purchased Reagent)		2,4,6-Tribromophenol (Surr)	5000 ug/mL
							2-Fluorobiphenyl	5000 ug/mL
							2-Fluorophenol (Surr)	5000 ug/mL
							Nitrobenzene-d5 (Surr)	5000 ug/mL
							Phenol-d5 (Surr)	5000 ug/mL
							Terphenyl-d14 (Surr)	5000 ug/mL
..MS-567672_00091	05/31/15		Restek, Lot A093709		(Purchased Reagent)		1,2,4-Trichlorobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Anthracene	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Pyrene	1000 ug/mL
..MS-567673_00096	05/31/15		Restek, Lot A093755		(Purchased Reagent)		Caprolactam	2000 ug/mL
MS-HSLB1B3SSV_00022	05/31/15	03/11/15	Methylene Chloride, Lot 87975	0.5 mL	MS-HSLB1_STK_00005	250 uL	1,2,4-Trichlorobenzene	100 ug/mL
							1,4-Dichlorobenzene	100 ug/mL
							2,4,6-Trichlorophenol	100 ug/mL
							2,4-Dinitrotoluene	100 ug/mL
							2-Chlorophenol	100 ug/mL
							2-Methylnaphthalene	100 ug/mL
							2-Methylphenol	100 ug/mL
							4-Chloro-3-methylphenol	100 ug/mL
							4-Nitrophenol	200 ug/mL
							Acenaphthene	100 ug/mL
							Anthracene	100 ug/mL
							N-Nitrosodi-n-propylamine	100 ug/mL
							Pentachlorophenol	200 ug/mL
							Pyrene	100 ug/mL
.MS-HSLB1_STK_00005	05/31/15	03/11/15	Methylene Chloride, Lot 87975	10 mL	MS-567672SEC_00012	2 mL	1,2,4-Trichlorobenzene	200 ug/mL
							1,4-Dichlorobenzene	200 ug/mL
							2,4,6-Trichlorophenol	200 ug/mL
							2,4-Dinitrotoluene	200 ug/mL
							2-Chlorophenol	200 ug/mL
							2-Methylnaphthalene	200 ug/mL
							2-Methylphenol	200 ug/mL
							4-Chloro-3-methylphenol	200 ug/mL

REAGENT TRACEABILITY SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Reagent ID	Exp Date	Prep Date	Dilutant Used	Reagent Final Volume	Parent Reagent		Analyte	Concentration
					Reagent ID	Volume Added		
							4-Nitrophenol	400 ug/mL
							Acenaphthene	200 ug/mL
							Anthracene	200 ug/mL
							N-Nitrosodi-n-propylamine	200 ug/mL
							Pentachlorophenol	400 ug/mL
							Pyrene	200 ug/mL
..MS-567672SEC_00012	05/31/15	Restek, Lot A094002			(Purchased Reagent)		1,2,4-Trichlorobenzene	1000 ug/mL
							1,4-Dichlorobenzene	1000 ug/mL
							2,4,6-Trichlorophenol	1000 ug/mL
							2,4-Dinitrotoluene	1000 ug/mL
							2-Chlorophenol	1000 ug/mL
							2-Methylnaphthalene	1000 ug/mL
							2-Methylphenol	1000 ug/mL
							4-Chloro-3-methylphenol	1000 ug/mL
							4-Nitrophenol	2000 ug/mL
							Acenaphthene	1000 ug/mL
							Anthracene	1000 ug/mL
							N-Nitrosodi-n-propylamine	1000 ug/mL
							Pentachlorophenol	2000 ug/mL
							Pyrene	1000 ug/mL
MS-HSLB2SSV_00020	05/16/15	03/11/15	Methylene Chloride, Lot 87975	0.5 mL	MS-HSLB2_STK_00004	250 uL	Caprolactam	100 ug/mL
.MS-HSLB2_STK_00004	05/16/15	05/16/14	Methylene Chloride, Lot 51144	10 mL	MS-567673SEC_00009	1 mL	Caprolactam	200 ug/mL
..MS-567673SEC_00009	07/31/15	Restek, Lot A0101015			(Purchased Reagent)		Caprolactam	2000 ug/mL

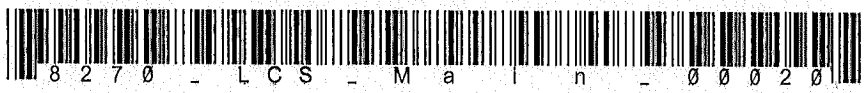
TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		
Comment:	Place all stocks in a sonication bath for 5 minutes before use. Dilute to 250mL in P&T MeOH 20mL of MS-569729 "8270 List 1 Std 1 MegaMix" 10mL of MS-569731 "Benzoic Acid" 1 year expiration Exchange /MeCl2 verification All Standards in fridge.		

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
1,1'-Biphenyl	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2,4,5-Tetrachlorobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2,4-Trichlorobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2-Dichlorobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,3-Dichlorobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,3-Dinitrobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,4-Dichlorobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,4-Dioxane	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1-Methylnaphthalene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,2'-oxybis[1-chloropropane]	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,3,4,6-Tetrachlorophenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4,5-Trichlorophenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4,6-Trichlorophenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dichlorophenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dimethylphenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dinitrophenol	MS-569729_00013	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL



Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
2,4-Dinitrotoluene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,6-Dichlorophenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,6-Dinitrotoluene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Chloronaphthalene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Chlorophenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Methylnaphthalene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Methylphenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Nitroaniline	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Nitrophenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3 & 4 Methylphenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3-Methylphenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3-Nitroaniline	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4,6-Dinitro-2-methylphenol	MS-569729_00013	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
4-Bromophenyl phenyl ether	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chloro-3-methylphenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chloroaniline	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

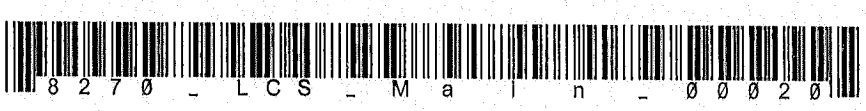
Reagent ID: 8270_LCS_Main_00020

		Expiration Date:	03/27/2016
Description:	USE IN CONJUCTION WITH 8270_LCS_S	Laboratory:	TestAmerica Denver
No. of Bottles:	2	Prepared By:	Isberg, Kyle
Storage Location:	North Prep	Solvent:	P&T Methanol
Reagent Volume:	250.000 mL	Solvent Lot:	MethanolP&T_00111
Creation Date:	03/27/2015		
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
4-Chlorophenyl phenyl ether	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Methylphenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Nitroaniline	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Nitrophenol	MS-569729_00013	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
Acenaphthene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Acenaphthylene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Acetophenone	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Aniline	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Anthracene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Azobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[a]anthracene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[a]pyrene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[b]fluoranthene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[g,h,i]perylene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[k]fluoranthene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzyl alcohol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL



Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Bis(2-chloroethoxy)methane	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-chloroethyl)ether	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-ethylhexyl) phthalate	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Butyl benzyl phthalate	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Carbazole	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Chrysene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dibenz(a,h)anthracene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dibenzofuran	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Diethyl phthalate	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dimethyl phthalate	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Di-n-butyl phthalate	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Di-n-octyl phthalate	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Diphenylamine	MS-569729_00013	05/31/2016	1713.40000	ug/mL	137.07200	ug/mL
Fluoranthene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Fluorene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		
Comment:	Place all stocks in a sonication bath for 5 minutes before use. Dilute to 250mL in P&T MeOH 20mL of MS-569729 "8270 List 1 Std 1 MegaMix" 10mL of MS-569731 "Benzoic Acid" 1 year expiration Exchange /MeCl2 verification All Standards in fridge.		

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Hexachlorobutadiene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorocyclopentadiene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachloroethane	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexadecane	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Indeno[1,2,3-cd]pyrene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Isophorone	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Naphthalene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
n-Decane	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Nitrobenzene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodimethylamine	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodi-n-propylamine	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodiphenylamine	MS-569729_00013	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
n-Octadecane	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pentachlorophenol	MS-569729_00013	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
Phenanthrene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Phenol	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

Reagent ID: 8270_LCS_Main_00020

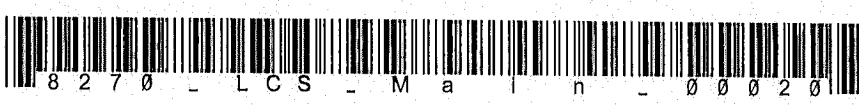
		Expiration Date:	03/27/2016
Description:	USE IN CONJUCTION WITH 8270_LCS_S	Laboratory:	TestAmerica Denver
No. of Bottles:	2	Prepared By:	Isberg, Kyle
Storage Location:	North Prep	Solvent:	P&T Methanol
Reagent Volume:	250.000 mL	Solvent Lot:	MethanolP&T_00111
Creation Date:	03/27/2015		
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Pyrene	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pyridine	MS-569729_00013	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,1'-Biphenyl	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2,4,5-Tetrachlorobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2,4-Trichlorobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2-Dichlorobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,3-Dichlorobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,3-Dinitrobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,4-Dichlorobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,4-Dioxane	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1-Methylnaphthalene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,2'-oxybis[1-chloropropane]	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,3,4,6-Tetrachlorophenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4,5-Trichlorophenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4,6-Trichlorophenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dichlorophenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING



Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		
Comment:	Place all stocks in a sonication bath for 5 minutes before use. Dilute to 250mL in P&T MeOH 20mL of MS-569729 "8270 List 1 Std 1 MegaMix" 10mL of MS-569731 "Benzoic Acid" 1 year expiration Exchange /MeCl2 verification All Standards in fridge.		

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
2,4-Dimethylphenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dinitrophenol	MS-569729_00014	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
2,4-Dinitrotoluene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,6-Dichlorophenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,6-Dinitrotoluene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Chloronaphthalene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Chlorophenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Methylnaphthalene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Methylphenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Nitroaniline	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Nitrophenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3 & 4 Methylphenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3-Methylphenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3-Nitroaniline	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4,6-Dinitro-2-methylphenol	MS-569729_00014	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
4-Bromophenyl phenyl ether	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

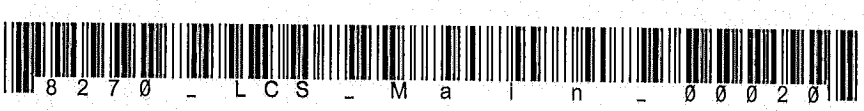
Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
4-Chloro-3-methylphenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chloroaniline	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chlorophenyl phenyl ether	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Methylphenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Nitroaniline	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Nitrophenol	MS-569729_00014	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
Acenaphthene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Acenaphthylene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Acetophenone	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Aniline	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Anthracene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Azobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[a]anthracene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[a]pyrene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[b]fluoranthene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[g,h,i]perylene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL



Reagent ID: 8270_LCS_Main_00020

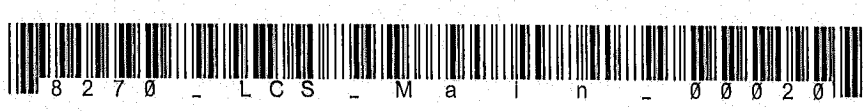
Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Benzo[k]fluoranthene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzyl alcohol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-chloroethoxy)methane	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-chloroethyl)ether	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-ethylhexyl) phthalate	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Butyl benzyl phthalate	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Carbazole	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Chrysene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dibenz(a,h)anthracene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dibenzofuran	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Diethyl phthalate	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dimethyl phthalate	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Di-n-butyl phthalate	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Di-n-octyl phthalate	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Diphenylamine	MS-569729_00014	05/31/2016	1713.40000	ug/mL	137.07200	ug/mL
Fluoranthene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING



Reagent ID: 8270_LCS_Main_00020

		Expiration Date:	03/27/2016
Description:	USE IN CONJUCTION WITH 8270_LCS_S	Laboratory:	TestAmerica Denver
No. of Bottles:	2	Prepared By:	Isberg, Kyle
Storage Location:	North Prep	Solvent:	P&T Methanol
Reagent Volume:	250.000 mL	Solvent Lot:	MethanolP&T_00111
Creation Date:	03/27/2015		
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Fluorene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorobutadiene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorocyclopentadiene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachloroethane	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexadecane	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Indeno[1,2,3-cd]pyrene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Isophorone	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Naphthalene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
n-Decane	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Nitrobenzene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodimethylamine	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodi-n-propylamine	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodiphenylamine	MS-569729_00014	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
n-Octadecane	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pentachlorophenol	MS-569729_00014	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL

Reagent ID: 8270_LCS_Main_00020

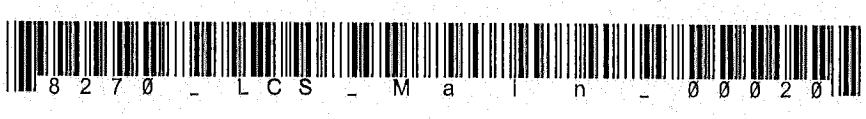
Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Phenanthrene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Phenol	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pyrene	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pyridine	MS-569729_00014	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,1'-Biphenyl	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2,4,5-Tetrachlorobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2,4-Trichlorobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2-Dichlorobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,3-Dichlorobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,3-Dinitrobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,4-Dichlorobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,4-Dioxane	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1-Methylnaphthalene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,2'-oxybis[1-chloropropane]	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,3,4,6-Tetrachlorophenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4,5-Trichlorophenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING



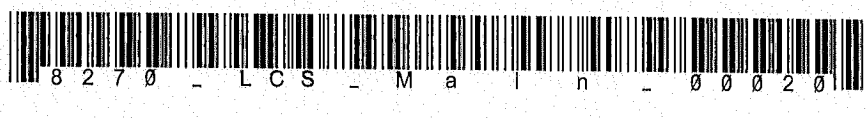
Reagent ID: 8270_LCS_Main_00020

		Expiration Date:	03/27/2016
Description:	USE IN CONJUCTION WITH 8270_LCS_S	Laboratory:	TestAmerica Denver
No. of Bottles:	2	Prepared By:	Isberg, Kyle
Storage Location:	North Prep	Solvent:	P&T Methanol
Reagent Volume:	250.000 mL	Solvent Lot:	MethanolP&T_00111
Creation Date:	03/27/2015		
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
2,4,6-Trichlorophenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dichlorophenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dimethylphenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dinitrophenol	MS-569729_00015	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
2,4-Dinitrotoluene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,6-Dichlorophenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,6-Dinitrotoluene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Chloronaphthalene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Chlorophenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Methylnaphthalene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Methylphenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Nitroaniline	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Nitrophenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3 & 4 Methylphenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3-Methylphenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3-Nitroaniline	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL



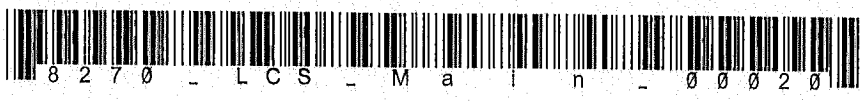
Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
4,6-Dinitro-2-methylphenol	MS-569729_00015	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
4-Bromophenyl phenyl ether	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chloro-3-methylphenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chloroaniline	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chlorophenyl phenyl ether	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Methylphenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Nitroaniline	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Nitrophenol	MS-569729_00015	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
Acenaphthene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Acenaphthylene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Acetophenone	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Aniline	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Anthracene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Azobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[a]anthracene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[a]pyrene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL



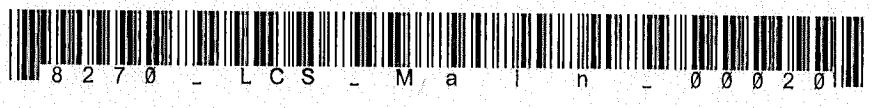
Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Benzo[b]fluoranthene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[g,h,i]perylene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[k]fluoranthene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzyl alcohol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-chloroethoxy)methane	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-chloroethyl)ether	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-ethylhexyl) phthalate	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Butyl benzyl phthalate	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Carbazole	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Chrysene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dibenz(a,h)anthracene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dibenzofuran	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Diethyl phthalate	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dimethyl phthalate	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Di-n-butyl phthalate	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Di-n-octyl phthalate	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL



Reagent ID: 8270_LCS_Main_00020

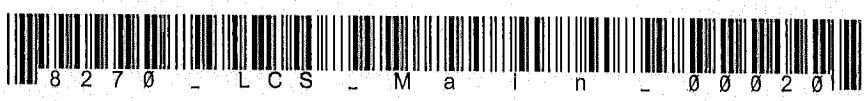
Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Diphenylamine	MS-569729_00015	05/31/2016	1713.40000	ug/mL	137.07200	ug/mL
Fluoranthene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Fluorene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorobutadiene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorocyclopentadiene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachloroethane	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexadecane	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Indeno[1,2,3-cd]pyrene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Isophorone	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Naphthalene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
n-Decane	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Nitrobenzene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodimethylamine	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodi-n-propylamine	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodiphenylamine	MS-569729_00015	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING



Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
n-Octadecane	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pentachlorophenol	MS-569729_00015	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
Phenanthrene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Phenol	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pyrene	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pyridine	MS-569729_00015	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,1'-Biphenyl	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2,4,5-Tetrachlorobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2,4-Trichlorobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,2-Dichlorobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,3-Dichlorobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,3-Dinitrobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,4-Dichlorobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1,4-Dioxane	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
1-Methylnaphthalene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,2'-oxybis[1-chloropropane]	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		
Comment:	Place all stocks in a sonication bath for 5 minutes before use. Dilute to 250mL in P&T MeOH 20mL of MS-569729 "8270 List 1 Std 1 MegaMix" 10mL of MS-569731 "Benzoic Acid" 1 year expiration Exchange /MeCl2 verification All Standards in fridge.		

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
2,3,4,6-Tetrachlorophenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4,5-Trichlorophenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4,6-Trichlorophenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dichlorophenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dimethylphenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,4-Dinitrophenol	MS-569729_00016	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
2,4-Dinitrotoluene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,6-Dichlorophenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2,6-Dinitrotoluene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Chloronaphthalene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Chlorophenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Methylnaphthalene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Methylphenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Nitroaniline	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
2-Nitrophenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3 & 4 Methylphenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

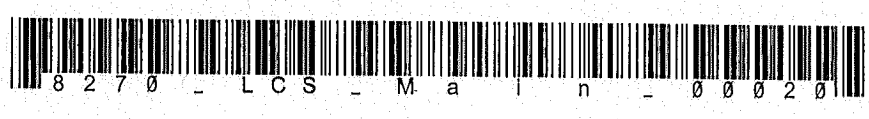
Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
3-Methylphenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
3-Nitroaniline	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4,6-Dinitro-2-methylphenol	MS-569729_00016	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
4-Bromophenyl phenyl ether	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chloro-3-methylphenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chloroaniline	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Chlorophenyl phenyl ether	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Methylphenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Nitroaniline	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
4-Nitrophenol	MS-569729_00016	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
Acenaphthene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Acenaphthylene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Acetophenone	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Aniline	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Anthracene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Azobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL



Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Benzo[a]anthracene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[a]pyrene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[b]fluoranthene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[g,h,i]perylene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzo[k]fluoranthene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzyl alcohol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-chloroethoxy)methane	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-chloroethyl)ether	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Bis(2-ethylhexyl) phthalate	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Butyl benzyl phthalate	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Carbazole	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Chrysene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dibenz(a,h)anthracene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dibenzofuran	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Diethyl phthalate	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Dimethyl phthalate	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		

Comment:
Place all stocks in a sonication bath for 5 minutes before use.
Dilute to 250mL in P&T MeOH
20mL of MS-569729 "8270 List 1 Std 1 MegaMix"
10mL of MS-569731 "Benzoic Acid"
1 year expiration
Exchange /MeCl2 verification
All Standards in fridge.

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
Di-n-butyl phthalate	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Di-n-octyl phthalate	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Diphenylamine	MS-569729_00016	05/31/2016	1713.40000	ug/mL	137.07200	ug/mL
Fluoranthene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Fluorene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorobutadiene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachlorocyclopentadiene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexachloroethane	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Hexadecane	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Indeno[1,2,3-cd]pyrene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Isophorone	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Naphthalene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
n-Decane	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Nitrobenzene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodimethylamine	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL

Reagent ID: 8270_LCS_Main_00020

Description:	USE IN CONJUCTION WITH 8270_LCS_S	Expiration Date:	03/27/2016
No. of Bottles:	2	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Isberg, Kyle
Reagent Volume:	250.000 mL	Solvent:	P&T Methanol
Creation Date:	03/27/2015	Solvent Lot:	MethanolP&T_00111
Open Date:			
Container(s):	3176000, 3176001		
Comment:	Place all stocks in a sonication bath for 5 minutes before use. Dilute to 250mL in P&T MeOH 20mL of MS-569729 "8270 List 1 Std 1 MegaMix" 10mL of MS-569731 "Benzoic Acid" 1 year expiration Exchange /MeCl2 verification All Standards in fridge.		

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
N-Nitrosodi-n-propylamine	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
N-Nitrosodiphenylamine	MS-569729_00016	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
n-Octadecane	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pentachlorophenol	MS-569729_00016	05/31/2016	2000.00000	ug/mL	160.00000	ug/mL
Phenanthrene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Phenol	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pyrene	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Pyridine	MS-569729_00016	05/31/2016	1000.00000	ug/mL	80.00000	ug/mL
Benzoic acid	MS-569731_00005	06/30/2016	2000.00000	ug/mL	80.00000	ug/mL
Indene	MS-569731_00005	06/30/2016	2000.00000	ug/mL	80.00000	ug/mL
Benzoic acid	MS-569731_00006	06/30/2016	2000.00000	ug/mL	80.00000	ug/mL
Indene	MS-569731_00006	06/30/2016	2000.00000	ug/mL	80.00000	ug/mL

Source Reagents

Reagent	Description	Type	Expiration	Vendor	Vendor Lot #	Vendor Cat Lot #	Volume Used	Volume Units
MS-569729_00013	HSLA Mega Mix (2015) 1000 ug/mL	ASTD	05/31/16	Restek	A0107399	569729	5.00000	mL
MS-569729_00014	HSLA Mega Mix (2015) 1000 ug/mL	ASTD	05/31/16	Restek	A0107399	569729	5.00000	mL
MS-569729_00015	HSLA Mega Mix (2015) 1000 ug/mL	ASTD	05/31/16	Restek	A0107399	569729	5.00000	mL
MS-569729_00016	HSLA Mega Mix (2015) 1000 ug/mL	ASTD	05/31/16	Restek	A0107399	569729	5.00000	mL
MS-569731_00005	HSLA Benzoic Acid (2015) 2000 ug/mL	ASTD	06/30/16	Restek	A0107943	569731	5.00000	mL
MS-569731_00006	HSLA Benzoic Acid (2015) 2000 ug/mL	ASTD	06/30/16	Restek	A0107943	569731	5.00000	mL

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

8270Surrogate_00078

Reagent ID: 8270Surrogate_00078

Description:	8270 Surrogate 100ug/ml	Expiration Date:	04/07/2016
No. of Bottles:	4	Laboratory:	TestAmerica Denver
Storage Location:	North Prep	Prepared By:	Knaub, Gentry L
Reagent Volume:	1000.000 mL	Solvent:	ACETONE
Creation Date:	04/07/2015	Solvent Lot:	Acetone_000115
Open Date:			
Container(s):	3189444, 3189445, 3189446, 3189447		
Comment:	Take 20mL of 8270SurHL and dilute to 1000mL. One year expiration date. Split into 4x250mL bottles. Requires solvent exchange to MeCl2 prior to submission for verification.		

Reagent Analyte Information

Analyte	Source ID	Source Exp. Date	Source Conc.	Source Conc. Units	Final Conc.	Final Conc. Units
2,4,6 - Tribromophenol	8270SurStkHL_00078	02/23/2018	5000.00000	ug/mL	100.00000	ug/mL
2,4,6-Tribromophenol	8270SurStkHL_00078	02/23/2018	5000.00000	ug/mL	100.00000	ug/mL
2-Fluorobiphenyl	8270SurStkHL_00078	02/23/2018	5000.00000	ug/mL	100.00000	ug/mL
2-Fluorophenol	8270SurStkHL_00078	02/23/2018	5000.00000	ug/mL	100.00000	ug/mL
Nitrobenzene-d5	8270SurStkHL_00078	02/23/2018	5000.00000	ug/mL	100.00000	ug/mL
Phenol-d5	8270SurStkHL_00078	02/23/2018	5000.00000	ug/mL	100.00000	ug/mL
Phenol-d6	8270SurStkHL_00078	02/23/2018	5000.00000	ug/mL	100.00000	ug/mL
Terphenyl-d14	8270SurStkHL_00078	02/23/2018	5000.00000	ug/mL	100.00000	ug/mL
2,4,6 - Tribromophenol	8270SurStkHL_00103	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
2,4,6-Tribromophenol	8270SurStkHL_00103	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
2-Fluorobiphenyl	8270SurStkHL_00103	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
2-Fluorophenol	8270SurStkHL_00103	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
Nitrobenzene-d5	8270SurStkHL_00103	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
Phenol-d5	8270SurStkHL_00103	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
Phenol-d6	8270SurStkHL_00103	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
Terphenyl-d14	8270SurStkHL_00103	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
2,4,6 - Tribromophenol	8270SurStkHL_00104	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL
2,4,6-Tribromophenol	8270SurStkHL_00104	05/31/2019	5000.00000	ug/mL	100.00000	ug/mL

Report Date: 09-Apr-2015 14:39:21 Chrom Revision: 2.2 13-Mar-2015 11:20:44

Preliminary Report

TestAmerica Denver
Recovery Report

Data File:	\\denchrom\chromdata\SMS_B2\20150409-33734.b\B2-9894.D				
Lims ID:	8270Surrogate_00078	Lab Sample ID:	Client 280-271876/28-A		
Client ID:					
Sample Type:	Client				
Inject. Date:	09-Apr-2015 13:16:30	ALS Bottle#:	6	Worklist Smp#:	28
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Sample Info:	8270Surrogate_00078				
Operator ID:	KIEKELD	Instrument ID:	SMS_B2		
Method:	\\denchrom\chromdata\SMS_B2\20150409-33734.b\SMSB2_8270C.m				
Limit Group:	MSSV - 8270C_625				
Method Label:	8270C / 625				
Last Update:	09-Apr-2015 14:39:10	Calib Date:	03-Apr-2015 11:31:30		
Integrator:	RTE	ID Type:	Deconvolution ID		
Quant Method:	Internal Standard	Quant By:	Initial Calibration		
Last ICal File:	\\denchrom\chromdata\SMS_B2\20150403-33531.b\B2-9841.D				
Column 1 :	VF-5ms (0.50 mm)		Det: MS SCAN		
Process Host:	XAWRK011				

Compound	Amount Added	Amount Recovered	% Rec.
$ 7 2-Fluorophenol	100.0	81.4	81.43
$ 8 Phenol-d5	100.0	80.2	80.21
$ 9 Nitrobenzene-d5	100.0	79.9	79.88
$ 11 2-Fluorobiphenyl	100.0	81.5	81.50
$ 12 2,4,6-Tribromophenol	100.0	66.4	66.35
$ 13 Terphenyl-d14	100.0	77.3	77.34

RESTEK

110 Benner Circle
Bellefonte, PA 16823-8812
Tel: (800)356-1688
Fax: (814)353-1309

www.restek.com

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ISO Guide 34 Accredited
Reference Material Producer
Certificate #3222.01

ACCREDITED
ISO/IEC 17025 Accredited
Testing Laboratory
Certificate #3222.02

Certificate of Composition

FOR LABORATORY USE ONLY-READ MSDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567685 **Lot No.:** A092712

Description : 8270 Surrogate Standard

8270 Surrogate Standard 5000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : January 2018 **Storage:** 10°C or colder

Handling: Sonicate prior to use.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	2-Fluorophenol	5,000.0 µg/mL	+/- 29.2761	µg/mL	Gravimetric
	CAS # 367-12-4		+/- 132.9947	µg/mL	Unstressed
	Purity 99%		+/- 163.4399	µg/mL	Stressed
2	Phenol-d5	5,000.0 µg/mL	+/- 29.2761	µg/mL	Gravimetric
	CAS # 4165-62-2		+/- 132.9947	µg/mL	Unstressed
	Purity 99%		+/- 163.4399	µg/mL	Stressed
3	Nitrobenzene-d5	5,000.0 µg/mL	+/- 29.2761	µg/mL	Gravimetric
	CAS # 4165-60-0		+/- 132.9947	µg/mL	Unstressed
	Purity 99%		+/- 163.4399	µg/mL	Stressed
4	2-Fluorobiphenyl	5,000.0 µg/mL	+/- 29.2761	µg/mL	Gravimetric
	CAS # 321-60-8		+/- 132.9947	µg/mL	Unstressed
	Purity 99%		+/- 163.4399	µg/mL	Stressed
5	2,4,6-Tribromophenol	5,000.0 µg/mL	+/- 29.2761	µg/mL	Gravimetric
	CAS # 118-79-6		+/- 132.9947	µg/mL	Unstressed
	Purity 99%		+/- 163.4399	µg/mL	Stressed
6	p-Terphenyl-d14	5,000.0 µg/mL	+/- 29.2761	µg/mL	Gravimetric
	CAS # 1718-51-0		+/- 132.9947	µg/mL	Unstressed
	Purity 99%		+/- 163.4399	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Tech Tips:
Due to the limited solubility of p-terphenyl-d14 in methanol, we do not recommend that this mixture be diluted in methanol.

General Certified Reference Material Notes

Expiration Notes:

- Expiration date of the unopened ampul stored at the recommended storage condition is the last day of the month listed in the expiration date field.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31840, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

Analytical Reference Materials

8270 Surrogate Standard

Catalog # 567685

Lot # A093638

110 Benner Circle Bellefonte, PA 16823-8812
(814) 353-1300

FOR LABORATORY USE ONLY. READ MSDS PRIOR TO USE.

RAW MATERIAL TEST INFORMATION AVAILABLE UPON REQUEST

MANUFACTURED UNDER RESTEK'S ISO 9001 REGISTERED QUALITY SYSTEM

RESTEK

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Certificate #3222.01

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ISO/IEC 17025 Accredited
Testing Laboratory
Certificate #3222.02

Certificate of Analysis

FOR LABORATORY USE ONLY-READ MSDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567685 **Lot No.:** A093638

Description : 8270 Surrogate Standard

8270 Surrogate Standard 5,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : February 2018 **Storage:** 10°C or colder

Handling: Sonicate prior to use.

C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	2-Fluorophenol	5,000.0 µg/mL	+/- 29.0689	µg/mL	Gravimetric
	CAS # 367-12-4		+/- 132.9492	µg/mL	Unstressed
	Purity 99%		+/- 163.4029	µg/mL	Stressed
2	Phenol-d5	5,000.0 µg/mL	+/- 29.0689	µg/mL	Gravimetric
	CAS # 4165-62-2		+/- 132.9492	µg/mL	Unstressed
	Purity 99%		+/- 163.4029	µg/mL	Stressed
3	Nitrobenzene-d5	5,000.0 µg/mL	+/- 29.0689	µg/mL	Gravimetric
	CAS # 4165-60-0		+/- 132.9492	µg/mL	Unstressed
	Purity 99%		+/- 163.4029	µg/mL	Stressed
4	2-Fluorobiphenyl	5,000.0 µg/mL	+/- 29.0689	µg/mL	Gravimetric
	CAS # 321-60-8		+/- 132.9492	µg/mL	Unstressed
	Purity 99%		+/- 163.4029	µg/mL	Stressed
5	2,4,6-Tribromophenol	5,000.0 µg/mL	+/- 29.0689	µg/mL	Gravimetric
	CAS # 118-79-6		+/- 132.9492	µg/mL	Unstressed
	Purity 99%		+/- 163.4029	µg/mL	Stressed
6	p-Terphenyl-d14	5,000.0 µg/mL	+/- 29.0689	µg/mL	Gravimetric
	CAS # 1718-51-0		+/- 132.9492	µg/mL	Unstressed
	Purity 99%		+/- 163.4029	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Tech Tips:

Due to the limited solubility of p-terphenyl-d14 in methanol, we do not recommend that this mixture be diluted in methanol.

Column:
30m x .25mm x .25um
Rtx-5 (cat.#10223)

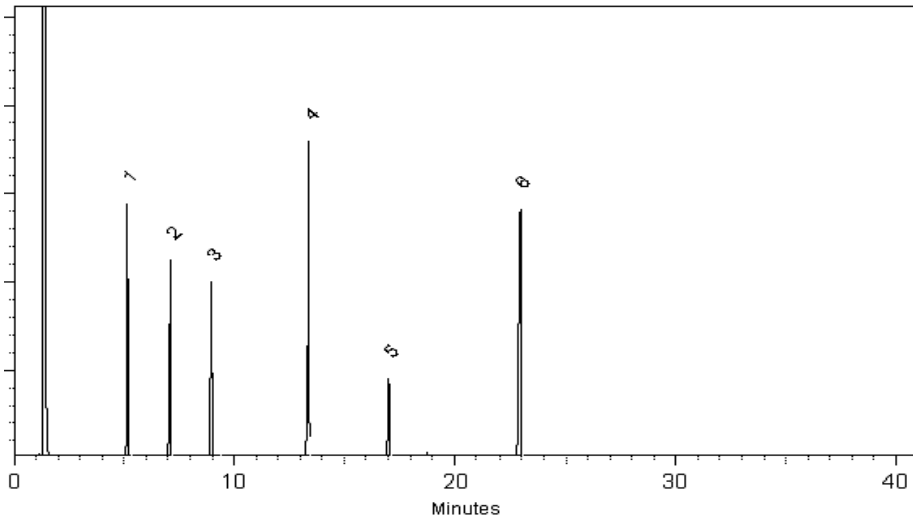
Carrier Gas:
hydrogen-constant pressure 10 psi.

Temp. Program:
40°C (hold 2 min.) to 330°C
@ 10°C/min. (hold 10 min.)

Inj. Temp:
250°C

Det. Temp:
330°C

Det. Type:
FID



Diane Shaffer - QA Analyst

Date Passed: 22-Feb-2013 **Balance:** 1128342313

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date of the unopened ampul stored at the recommended storage condition is the last day of the month listed in the expiration date field.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31840, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

RESTEK

Chemical Standard Batch Sheet
Lot #: A093638

Catalog #: 567685	**Target:** 5000 ug/mL	
Description: 8270 Surrogate Standard		
Solvent: Methylene Chloride	**Solvent Lot:** 127438	**Final Volume:** 6,000 ml

Made by: Matt Hepfer	**Date:** 2/19/2013 1:13:54PM			
Tested by: Jennifer Pollino	**Date:** 2/21/2013 11:49:35AM			
Pass	**By:** Diane Shaffer		**Date:** 2/22/2013 10:09:20A	
Packaged by: Alexandria Pavkovich / Kendra Swope	**Date:** 2/21/2013 10:36:16A	**No. Units:** 1,023	**Pkg Size:**	5 mL
Balance Used: BEDEARMBALPC1 XP205	**Serial #:** 1128342313			

Compound	CAS	Storage Location	Lot #	Purity	Target Conc(ug/mL)	Target	Actual	Calc Conc(ug/mL)
p-Terphenyl-d14	1718-51-0	R0504	PR-20577	0.99	5,000.00	30,000.00 mg	30,000.00 mg	5,000.0
Nitrobenzene-d5	4165-60-0	R0526	PR-20474	0.99	5,000.00	30,000.00 mg	30,000.00 mg	5,000.0
2-Fluorobiphenyl	321-60-8	R0539	E11Y047	0.99	5,000.00	30,000.00 mg	30,000.00 mg	5,000.0
2,4,6-Tribromophenol	118-79-6	R0619	1383140V	0.99	5,000.00	30,000.00 mg	30,000.00 mg	5,000.0
2-Fluorophenol	367-12-4	R0635	STBC5591V	0.99	5,000.00	30,000.00 mg	30,000.00 mg	5,000.0
Phenol-d5	4165-62-2	R1135	X479P1	0.99	5,000.00	30,000.00 mg	30,000.00 mg	5,000.0

QA Report: 8270 Surrogate Standard (Cat.#567685)

	Runs of Lot # A092712						Runs of Lot # A093638							
COMPONENT	Run #1	Run #2	Run #3	AVG	STD DEV	% RSD	Run #1	Run #2	Run #3	AVG	STD DEV	% RSD	%D MEAN	P/F
2-Fluorophenol	3567168	3686433	3610358	3621320	60383	1.67	3740895	3706516	3740206	3729206	19653	0.53	-2.98	PASS
Phenol-d5	4151283	4284919	4199657	4211953	67661	1.61	4299218	4265705	4301647	4288857	20087	0.47	-1.83	PASS
Nitrobenzene-d5	3474347	3583156	3514256	3523920	55044	1.56	3599035	3574157	3601170	3591454	15018	0.42	-1.92	PASS
2-Fluorobiphenyl	5477353	5637195	5541430	5551993	80443	1.45	5659153	5624786	5659681	5647873	19996	0.35	-1.73	PASS
2,4,6-Tribromophenol	1317679	1355641	1333550	1335623	19066	1.43	1356813	1353676	1360571	1357020	3452	0.25	-1.60	PASS
p-Terphenyl-d14	5916769	6051268	5989106	5985714	67314	1.12	6094261	6069529	6090064	6084618	13235	0.22	-1.65	PASS

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FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567685 **Lot No.:** A0103615

Description : 8270 Surrogate Standard

8270 Surrogate Standard 5,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : May 31, 2019 **Storage:** 10°C or colder

Handling: Sonicate prior to use.

CERTIFIED VALUES

Elution Order	Compound		Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	2-Fluorophenol		5,003.5 µg/mL	+/- 29.0892 µg/mL		Gravimetric
	CAS # 367-12-4	(Lot STBC5591V)		+/- 124.6713 µg/mL		Unstressed
	Purity 99%			+/- 156.7818 µg/mL		Stressed
2	Phenol-d5		5,002.9 µg/mL	+/- 29.0860 µg/mL		Gravimetric
	CAS # 4165-62-2	(Lot M387P4)		+/- 124.6575 µg/mL		Unstressed
	Purity 99%			+/- 156.7644 µg/mL		Stressed
3	Nitrobenzene-d5		5,001.4 µg/mL	+/- 29.0773 µg/mL		Gravimetric
	CAS # 4165-60-0	(Lot PR-20474)		+/- 124.6201 µg/mL		Unstressed
	Purity 99%			+/- 156.7174 µg/mL		Stressed
4	2-Fluorobiphenyl		5,004.4 µg/mL	+/- 29.0947 µg/mL		Gravimetric
	CAS # 321-60-8	(Lot E11Y047)		+/- 124.6949 µg/mL		Unstressed
	Purity 99%			+/- 156.8114 µg/mL		Stressed
5	2,4,6-Tribromophenol		5,003.9 µg/mL	+/- 29.0914 µg/mL		Gravimetric
	CAS # 118-79-6	(Lot 29699MJV)		+/- 124.6805 µg/mL		Unstressed
	Purity 99%			+/- 156.7934 µg/mL		Stressed
6	p-Terphenyl-d14		5,007.1 µg/mL	+/- 29.1100 µg/mL		Gravimetric
	CAS # 1718-51-0	(Lot PR-20577)		+/- 124.7604 µg/mL		Unstressed
	Purity 99%			+/- 156.8938 µg/mL		Stressed

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31840, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

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This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567685 **Lot No.:** A0103615

Description : 8270 Surrogate Standard

8270 Surrogate Standard 5,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : May 31, 2019 **Storage:** 10°C or colder

Handling: Sonicate prior to use.

CERTIFIED VALUES

Elution Order	Compound		Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	2-Fluorophenol		5,003.5 µg/mL	+/- 29.0892 µg/mL		Gravimetric
	CAS # 367-12-4	(Lot STBC5591V)		+/- 124.6713 µg/mL		Unstressed
	Purity 99%			+/- 156.7818 µg/mL		Stressed
2	Phenol-d5		5,002.9 µg/mL	+/- 29.0860 µg/mL		Gravimetric
	CAS # 4165-62-2	(Lot M387P4)		+/- 124.6575 µg/mL		Unstressed
	Purity 99%			+/- 156.7644 µg/mL		Stressed
3	Nitrobenzene-d5		5,001.4 µg/mL	+/- 29.0773 µg/mL		Gravimetric
	CAS # 4165-60-0	(Lot PR-20474)		+/- 124.6201 µg/mL		Unstressed
	Purity 99%			+/- 156.7174 µg/mL		Stressed
4	2-Fluorobiphenyl		5,004.4 µg/mL	+/- 29.0947 µg/mL		Gravimetric
	CAS # 321-60-8	(Lot E11Y047)		+/- 124.6949 µg/mL		Unstressed
	Purity 99%			+/- 156.8114 µg/mL		Stressed
5	2,4,6-Tribromophenol		5,003.9 µg/mL	+/- 29.0914 µg/mL		Gravimetric
	CAS # 118-79-6	(Lot 29699MJV)		+/- 124.6805 µg/mL		Unstressed
	Purity 99%			+/- 156.7934 µg/mL		Stressed
6	p-Terphenyl-d14		5,007.1 µg/mL	+/- 29.1100 µg/mL		Gravimetric
	CAS # 1718-51-0	(Lot PR-20577)		+/- 124.7604 µg/mL		Unstressed
	Purity 99%			+/- 156.8938 µg/mL		Stressed

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
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Purity Notes:

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0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

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Manufacturing Notes:

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Handling Notes:

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Catalog No. : 567685 **Lot No.:** A0103615

Description : 8270 Surrogate Standard

8270 Surrogate Standard 5,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : May 31, 2019 **Storage:** 10°C or colder

Handling: Sonicate prior to use.

CERTIFIED VALUES

Elution Order	Compound		Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	2-Fluorophenol		5,003.5 µg/mL	+/- 29.0892 µg/mL		Gravimetric
	CAS # 367-12-4	(Lot STBC5591V)		+/- 124.6713 µg/mL		Unstressed
	Purity 99%			+/- 156.7818 µg/mL		Stressed
2	Phenol-d5		5,002.9 µg/mL	+/- 29.0860 µg/mL		Gravimetric
	CAS # 4165-62-2	(Lot M387P4)		+/- 124.6575 µg/mL		Unstressed
	Purity 99%			+/- 156.7644 µg/mL		Stressed
3	Nitrobenzene-d5		5,001.4 µg/mL	+/- 29.0773 µg/mL		Gravimetric
	CAS # 4165-60-0	(Lot PR-20474)		+/- 124.6201 µg/mL		Unstressed
	Purity 99%			+/- 156.7174 µg/mL		Stressed
4	2-Fluorobiphenyl		5,004.4 µg/mL	+/- 29.0947 µg/mL		Gravimetric
	CAS # 321-60-8	(Lot E11Y047)		+/- 124.6949 µg/mL		Unstressed
	Purity 99%			+/- 156.8114 µg/mL		Stressed
5	2,4,6-Tribromophenol		5,003.9 µg/mL	+/- 29.0914 µg/mL		Gravimetric
	CAS # 118-79-6	(Lot 29699MJV)		+/- 124.6805 µg/mL		Unstressed
	Purity 99%			+/- 156.7934 µg/mL		Stressed
6	p-Terphenyl-d14		5,007.1 µg/mL	+/- 29.1100 µg/mL		Gravimetric
	CAS # 1718-51-0	(Lot PR-20577)		+/- 124.7604 µg/mL		Unstressed
	Purity 99%			+/- 156.8938 µg/mL		Stressed

General Certified Reference Material Notes

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0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

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Manufacturing Notes:

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Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31840, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

Standard Verification Form

Verification (New vendor or problematic Standard)	☐	Re-Verification	☒
TALS Reagent Record — ampules were not used therefore original CDs are un-changed			
New	☐	Copied	☐
COA Reviewed against formulary report			☒

Document instrument verification if need (Initial or re-verification):		
Department	**Acceptance Criteria**	
	Standard Analytes	**Poor Performers* and Esterified Analytes**
GC/HPLC	≤ 15 %D	≤ 35 %D or ≤ 50 %D for dinoseb
X GCMS/LCMS	≤ 35 %D	≤ 55 %D
MSVOA	≤ 25 %D	≤ 55 %D
Metals	≤ 8 %D	NA
Wet Chemistry	≤ 5 %D	NA

Standard Name	MS-567672	Standard ID	RES HSLA MajorMax	
Verified by	M. Hoffman	Instrument ID	Y	
Verification Date	8/11/14	Method Reference	8270C	
Reference Standard ID	567672SEC	Batch # A094002	Analytical Batch	
Analyte/Mix	Prepared Concentration	Verification Concentration	% Diff	Pass/Fail
See Attached Report.				All compounds Pass. pss 8/15/14
New Expiration Date:	5/2015	New TALS ID	★	

New expiration date can be no greater than ½ the designated standards shelf life from the date of re-verification. Standards can only be re-verified one time.

Comment:	test one ampule of lot A093709 & apply to all ampules of that lot

1st Level Review MH Date: 8/14/14

2nd Level Review [signature] Date: 8/14/14

QA Review (Re-verification only)	[signature]	Date: 8/15/14

Attach form, supporting documentation and original CoA to new verified or re-verified standard record in TALS.
*See analytical SOP for details on poor performing analytes.

★ Use same TALS ID with updated expiration date. Attach this paperwork to each record. pss 8/15/14

Analytical Reference Materials

8270 List 1 / Std #1 MegaMix

Catalog # 567672

Lot # A093709

110 Benner Circle Bellefonte, PA 16823-8812
(814) 353-1300

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RAW MATERIAL TEST INFORMATION AVAILABLE UPON REQUEST

MANUFACTURED UNDER RESTEK'S ISO 9001 REGISTERED QUALITY SYSTEM

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Certificate of Analysis

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This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567672 **Lot No.:** A093709

Description : 8270 List 1 / Std #1 MegaMix

8270 List 1 / Std #1 MegaMix 500-2000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : August 2014 **Storage:** 10°C or colder

Handling: Sonication required. Mix is photosensitive.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	1,4-Dioxane **CAS #** 123-91-1 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
2	Pyridine **CAS #** 110-86-1 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
3	N-Nitrosodimethylamine **CAS #** 62-75-9 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
4	Aniline **CAS #** 62-53-3 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
5	Phenol **CAS #** 108-95-2 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
6	Bis(2-chloroethyl)ether **CAS #** 111-44-4 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
7	2-Chlorophenol **CAS #** 95-57-8 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
8	1,3-Dichlorobenzene **CAS #** 541-73-1 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
9	1,4-Dichlorobenzene **CAS #** 106-46-7 **Purity** 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

10	1,2-Dichlorobenzene **CAS #** 95-50-1 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
11	Benzyl alcohol **CAS #** 100-51-6 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
12	Bis(2-chloroisopropyl)ether **CAS #** 108-60-1 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
13	2-Methylphenol (o-cresol) **CAS #** 95-48-7 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
14	Acetophenone **CAS #** 98-86-2 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
15	Hexachloroethane **CAS #** 67-72-1 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
16	N-Nitroso-di-n-propylamine **CAS #** 621-64-7 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
17	4-Methylphenol (p-cresol) **CAS #** 106-44-5 **Purity** 99%	500.0	µg/mL	+/- 2.9070 +/- 4.3978 +/- 8.6943	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
18	3-Methylphenol (m-cresol) **CAS #** 108-39-4 **Purity** 99%	500.0	µg/mL	+/- 2.9070 +/- 4.3978 +/- 8.6943	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
19	n-Decane (C10) **CAS #** 124-18-5 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
20	n-Octadecane (C18) **CAS #** 593-45-3 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
21	Nitrobenzene **CAS #** 98-95-3 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
22	Isophorone **CAS #** 78-59-1 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
23	2-Nitrophenol **CAS #** 88-75-5 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
24	2,4-Dimethylphenol **CAS #** 105-67-9 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
25	Bis(2-chlorocthoxy)methane **CAS #** 111-91-1 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
26	2,4-Dichlorophenol **CAS #** 120-83-2 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
27	1,2,4-Trichlorobenzene **CAS #** 120-82-1 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
28	Naphthalene **CAS #** 91-20-3 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

29	4-Chloroaniline	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 106-47-8			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
30	Hexachlorobutadiene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 87-68-3			+/-	8.7956	µg/mL	Unstressed
	Purity 97%			+/-	17.3886	µg/mL	Stressed
31	2-Methylnaphthalene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-57-6			+/-	8.7957	µg/mL	Unstressed
	Purity 96%			+/-	17.3887	µg/mL	Stressed
32	4-Chloro-3-methylphenol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 59-50-7			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
33	1-Methylnaphthalene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 90-12-0			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
34	1,2,4,5-Tetrachlorobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-94-3			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
35	Hexachlorocyclopentadiene	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 77-47-4			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
36	2,4,6-Trichlorophenol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 88-06-2			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
37	2,4,5-Trichlorophenol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-95-4			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
38	2-Chloronaphthalene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-58-7			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
39	Biphenyl	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 92-52-4			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
40	2-Nitroaniline	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 88-74-4			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
41	Acenaphthylene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 208-96-8			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
42	1,3-Dinitrobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 99-65-0			+/-	8.7956	µg/mL	Unstressed
	Purity 97%			+/-	17.3886	µg/mL	Stressed
43	Dimethylphthalate	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 131-11-3			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
44	2,6-Dinitrotoluene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 606-20-2			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
45	Acenaphthene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 83-32-9			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
46	3-Nitroaniline	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 99-09-2			+/-	8.7956	µg/mL	Unstressed
	Purity 97%			+/-	17.3886	µg/mL	Stressed
47	2,4-Dinitrophenol	2,000.0	µg/mL	+/-	11.6282	µg/mL	Gravimetric
	CAS # 51-28-5			+/-	17.5913	µg/mL	Unstressed
	Purity 99%			+/-	34.7772	µg/mL	Stressed

48	Dibenzofuran **CAS #** 132-64-9 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
49	2,4-Dinitrotoluene **CAS #** 121-14-2 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
50	4-Nitrophenol **CAS #** 100-02-7 **Purity** 99%	2,000.0	µg/mL	+/- 11.6282 +/- 17.5913 +/- 34.7772	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
51	2,3,4,6-Tetrachlorophenol **CAS #** 58-90-2 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
52	Fluorene **CAS #** 86-73-7 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
53	4-Chlorophenyl phenyl ether **CAS #** 7005-72-3 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
54	Diethylphthalate **CAS #** 84-66-2 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
55	4-Nitroaniline **CAS #** 100-01-6 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
56	4,6-Dinitro-2-methylphenol (Dinitro-o-cresol) **CAS #** 534-52-1 **Purity** 99%	2,000.0	µg/mL	+/- 11.6282 +/- 17.5913 +/- 34.7772	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
57	Azobenzene **CAS #** 103-33-3 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
58	4-Bromophenyl phenyl ether **CAS #** 101-55-3 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
59	Hexachlorobenzene **CAS #** 118-74-1 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
60	Pentachlorophenol **CAS #** 87-86-5 **Purity** 99%	2,000.0	µg/mL	+/- 11.6282 +/- 17.5913 +/- 34.7772	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
61	Phenanthrene **CAS #** 85-01-8 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
62	Anthracene **CAS #** 120-12-7 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
63	n-Hexadecane (C16) **CAS #** 544-76-3 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
64	Carbazole **CAS #** 86-74-8 **Purity** 98%	1,000.0	µg/mL	+/- 5.8140 +/- 8.7956 +/- 17.3885	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
65	Di-n-butylphthalate **CAS #** 84-74-2 **Purity** 99%	1,000.0	µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
66	Fluoranthene **CAS #** 206-44-0 **Purity** 98%	1,000.0	µg/mL	+/- 5.8140 +/- 8.7956 +/- 17.3885	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

67	Pyrene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 129-00-0			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
68	Benzyl butyl phthalate	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 85-68-7			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
69	Benz(a)anthracene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 56-55-3			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
70	Chrysene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 218-01-9			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
71	Bis(2-ethylhexyl)phthalate	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 117-81-7			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
72	Di-n-octyl phthalate	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 117-84-0			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
73	Benzo(b)fluoranthene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 205-99-2			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
74	Benzo(k)fluoranthene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 207-08-9			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
75	Benzo(a)pyrene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 50-32-8			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
76	Indeno(1,2,3-cd)pyrene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 193-39-5			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
77	Dibenz(a,h)anthracene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 53-70-3			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
78	Benzo(g,h,i)perylene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 191-24-2			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Column:
30m x .25mm x .25um
Rtx-5 (cat.#10223)

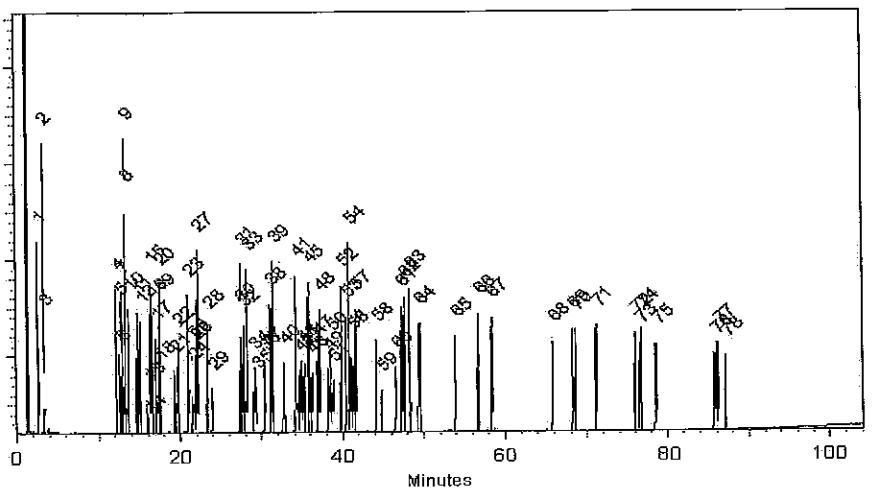
Carrier Gas:
hydrogen-constant pressure 10 psi

Temp. Program:
35°C (hold 3 min.) to 330°C
@ 3°C/min. (hold 3 min.)

Inj. Temp:
250°C

Det. Temp:
300°C

Det. Type:
FID



Jodi E. Breon - QA Analyst

Date Passed: 11-Mar-2013 **Balance:** 1128342313

Manufactured under Restek's ISO 9001:2008
Registered Quality System
Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date of the unopened ampul stored at the recommended storage condition is the last day of the month listed in the expiration date field.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31840, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

RESTEK

Chemical Standard Batch Sheet
Lot #: A093709

Catalog #: 567672	**Target:** 500-2000 ug/mL	
Description: 8270 List 1 / Std #1 MegaMix		
Solvent: Methylene Chloride	**Solvent Lot:** 127788	**Final Volume:** 4,000 ml

Made by: Matt Hepfer	**Date:** 3/8/2013 12:21:16PM	
Tested by: Jodi Breon	**Date:** 3/11/2013 3:28:30PM	
Pass	**By:** Jodi Breon	**Date:** 3/11/2013 3:28:33PM
Packaged by: Brandon Reish / Matt Hepfer	**Date:** 3/10/2013 5:08:52PM **No. Units:** 666	**Pkg Size:** 5 mL
Balance Used: BEDEARMBALPC1 XP205	**Serial #:** 1128342313	

Compound	CAS	Storage Location	Lot #	Purity	Target Conc(ug/mL)	Target	Actual	Calc Conc(ug/mL)
N-Nitrosodimethylamine	62-75-9	CAL012	1260200	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Pyridine	110-86-1	F0049	02718MW	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Bis(2-chloroethoxy)methane	111-91-1	R0253	317200	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
4-Bromophenyl phenyl ether	101-55-3	R0254	STBB9729V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
4-Chlorophenyl phenyl ether	7005-72-3	R0255	MKBL1347V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2,4-Dimethylphenol	105-67-9	R0256	10165155	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2-Nitrophenol	88-75-5	R0263	02611AH	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Benz(a)anthracene	56-55-3	R0469	ER121707-01	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Benzo(b)fluoranthene	205-99-2	R0470	ER022008-02	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Benzo(k)fluoranthene	207-08-9	R0471	ER061608-02	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Benzo(g,h,i)perylene	191-24-2	R0473	ER020708-08	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Benzo(a)pyrene	50-32-8	R0474	ER071309-02	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Benzyl butyl phthalate	85-68-7	R0475	03027HV	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Bis(2-ethylhexyl)phthalate	117-81-7	R0479	MKBH9511V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Carbazole	86-74-8	R0480	S42950-417	0.98	1,000.00	4,081.63 mg	4,081.60 mg	1,000.0
1,2-Dichlorobenzene	95-50-1	R0484	68996CMV	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
1,3-Dichlorobenzene	541-73-1	R0485	BCBC1891V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
1,4-Dichlorobenzene	106-46-7	R0486	MKBG7690V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2,4-Dinitrotoluene	121-14-2	R0488	MKAA0690V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2,6-Dinitrotoluene	606-20-2	R0489	1437483V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2-Methylnaphthalene	91-57-6	R0490	19399MJV	0.96	1,000.00	4,166.67 mg	4,166.70 mg	1,000.0
2-Nitroaniline	88-74-4	R0491	MKBF9132V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
3-Nitroaniline	99-09-2	R0492	MKBH5131V	0.97	1,000.00	4,123.71 mg	4,123.70 mg	1,000.0
4-Nitroaniline	100-01-6	R0493	BCBG4702V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
4-Nitrophenol	100-02-7	R0494	MKBF5564V	0.99	2,000.00	8,000.00 mg	8,000.00 mg	2,000.0
1,2,4-Trichlorobenzene	120-82-1	R0496	26896BM	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Acenaphthylene	208-96-8	R0497	ER030707-01	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Anthracene	120-12-7	R0499	MKBK5208V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Pentachlorophenol	87-86-5	R0501	130201JLM	0.99	2,000.00	8,000.00 mg	8,000.00 mg	2,000.0
Phenanthrene	85-01-8	R0502	MKBJ4205V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Pyrene	129-00-0	R0503	S22012V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Chrysene	218-01-9	R0505	ER120810-02	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Dibenz(a,h)anthracene	53-70-3	R0507	ER032211-01	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Di-n-butylphthalate	84-74-2	R0508	MKBG1851V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Fluoranthene	206-44-0	R0514	00828AJ	0.98	1,000.00	4,081.63 mg	4,081.60 mg	1,000.0
Fluorene	86-73-7	R0515	1326187	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Hexachlorocyclopentadiene	77-47-4	R0517	14752	0.98	1,000.00	4,081.63 mg	4,081.60 mg	1,000.0
Hexachlorobenzene	118-74-1	R0518	LB93343V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Hexachlorobutadiene	87-68-3	R0519	K22W009	0.97	1,000.00	4,123.71 mg	4,123.70 mg	1,000.0
Hexachloroethane	67-72-1	R0520	4H3SF	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Indeno(1,2,3-cd)pyrene	193-39-5	R0521	ER082107-02	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Naphthalene	91-20-3	R0524	MKBF8633V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Nitrobenzene	98-95-3	R0525	65096APV	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
N-Nitroso-di-n-propylamine	621-64-7	R0527	OPAGF-LS	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2,3,4,6-Tetrachlorophenol	58-90-2	R0528	ER100206-01	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0

Compound	CAS	Storage Location	Lot #	Purity	Target Conc(ug/mL)	Target	Actual	Calc Conc(ug/mL)
n-Decane (C10)	124-18-5	R0531	SHBB8486V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Acenaphthene	83-32-9	R0537	MKBH3748V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
n-Octadecane (C18)	593-45-3	R0549	FGG01-YPFG	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
1,2,4,5-Tetrachlorobenzene	95-94-3	R0555	06024AIV	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Acetophenone	98-86-2	R0561	MKBH7920V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Biphenyl	92-52-4	R0562	1277976	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
n-Hexadecane (C16)	544-76-3	R0563	SHBC3991V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
1-Methylnaphthalene	90-12-0	R0571	5250.00-10	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
1,4-Dioxane	123-91-1	R0581	SHBC3639V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Benzyl alcohol	100-51-6	R0606	02236CC	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Bis(2-chloroethyl)ether	111-44-4	R0607	45296HKV	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Bis(2-chloroisopropyl)ether	108-60-1	R0608	LB95678V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
4-Chloroaniline	106-47-8	R0609	12528PH	0.98	1,000.00	4,081.63 mg	4,081.60 mg	1,000.0
4-Chloro-3-methylphenol	59-50-7	R0610	STBC0769V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2-Chloronaphthalene	91-58-7	R0611	FIJ01	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2-Chlorophenol	95-57-8	R0612	MKBD3900V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2,4-Dichlorophenol	120-83-2	R0614	BCBH1617V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
4,6-Dinitro-2-methylphenol (Dinitro-o-cr	534-52-1	R0615	13212DW	0.99	2,000.00	8,000.00 mg	8,000.00 mg	2,000.0
2,4-Dinitrophenol	51-28-5	R0616	MKBH0709V	0.99	2,000.00	8,000.00 mg	8,000.00 mg	2,000.0
2-Methylphenol (o-cresol)	95-48-7	R0617	29996EK	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
4-Methylphenol (p-cresol)	106-44-5	R0618	49396APV	0.99	500.00	2,000.00 mg	2,000.00 mg	500.0
2,4,5-Trichlorophenol	95-95-4	R0620	FHM01	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
2,4,6-Trichlorophenol	88-06-2	R0621	MKBH7393V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Phenol	108-95-2	R0627	BCBC0688	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Dibenzofuran	132-64-9	R0630	MKBK2375V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Diethylphthalate	84-66-2	R0631	MKBG0599V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Dimethylphthalate	131-11-3	R0632	10117699	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Di-n-octyl phthalate	117-84-0	R0633	36100	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Isophorone	78-59-1	R0636	06705DE	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Azobenzene	103-33-3	R0645	130305JLM	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
Aniline	62-53-3	R0647	BCBG3682V	0.99	1,000.00	4,000.00 mg	4,000.00 mg	1,000.0
3-Methylphenol (m-cresol)	108-39-4	R0648	04627KA	0.99	500.00	2,000.00 mg	2,000.00 mg	500.0
1,3-Dinitrobenzene	99-65-0	R0690	0001338960	0.97	1,000.00	4,123.71 mg	4,123.70 mg	1,000.0

QA Report: 8270 List 1/ Std #1 MegaMix (Cat.#567672)

COMPONENT	Runs of Lot # A094002 Run #1	Run #2	Run #3	AVG	STD DEV	% RSD	Runs of Lot # A093709 Run #1	Run #2	Run #3	AVG	STD DEV	% RSD	%D MEAN	P/F
1,4-Dioxane	188646	192821	191880	191116	2190	1.15	192336	189509	189679	190508	1585	0.83	0.32	PASS
pyridine	415451	428642	427105	423733	7213	1.70	415345	408462	408591	410799	3937	0.96	3.05	PASS
n-nitrosodimethylamine	223317	225514	225683	224838	1320	0.59	230349	230199	229641	230063	373	0.16	-2.32	PASS
aniline	493725	503628	503624	500326	5716	1.14	500907	496113	495103	497374	3101	0.62	0.59	PASS
phenol	817868	839195	837639	831567	11889	1.43	835201	826269	822169	827880	6664	0.80	0.44	PASS
Bis(2-chlorotheyl) ether	149557	150074	151312	150314	902	0.60	151577	151758	151779	151705	111	0.07	-0.92	PASS
2-chlorophenol	555903	568972	568346	564407	7371	1.31	569567	561806	562793	564722	4225	0.75	-0.06	PASS
1,3-dichlorobenzene	315458	322043	321021	319507	3544	1.11	322065	319592	317894	319850	2097	0.66	-0.11	PASS
1,4-dichlorobenzene	330944	338234	337988	335722	4140	1.23	334527	331280	330934	332247	1982	0.60	1.04	PASS
1,2-dichlorobenzene	330894	338189	337997	335693	4157	1.24	336250	333171	332806	334076	1892	0.57	0.48	PASS
benzyl alcohol	495115	507587	507493	503398	7174	1.43	497826	493630	493743	495066	2391	0.48	1.66	PASS
bis(2-chloroisopropyl) ether	172397	176239	175001	174546	1961	1.12	234012	231770	230920	232234	1597	0.69	-33.05	PASSES per customer ap...
2-methylphenol	571992	584082	581689	579254	6402	1.11	508833	504219	502597	505216	3235	0.64	12.78	PASSES per customer ap...
Acetophenone & 'hexachloroethane	491568	505669	495927	497721	7220	1.45	493188	488122	492221	491177	2690	0.55	1.31	PASS
N-Nitroso-di-n-propylamine	361179	368406	366589	365391	3759	1.03	368628	365316	363726	365890	2501	0.68	-0.14	PASS
4 & 3-methylphenol	483616	493737	491316	489556	5285	1.08	486478	482308	480327	483038	3140	0.65	1.33	PASS
nitrobenzene	386839	393863	391993	390898	3638	0.93	390916	387583	385700	388066	2641	0.68	0.72	PASS
isophorone	480255	491568	489148	486990	5957	1.22	469615	467260	464590	467155	2514	0.54	4.07	PASS
2-nitrophenol	301545	309268	306582	305798	3921	1.28	307220	305842	303776	305613	1733	0.57	0.06	PASS
2,4-dimethylphenol	497531	508856	507673	504687	6225	1.23	508912	504842	502705	505486	3153	0.62	-0.16	PASS
bis(2-chloroethoxy) methane	155792	159441	158707	157980	1930	1.22	163541	162149	161852	162514	902	0.55	-2.87	PASS
2,4-dichlorophenol	272471	278784	279120	276792	3746	1.35	277729	275396	274832	275986	1536	0.56	0.29	PASS
1,2,4-trichlorobenzene	274752	280745	280628	278708	3427	1.23	275719	273362	271330	273470	2197	0.80	1.88	PASS
naphthalene	632713	647089	644183	641328	7601	1.19	645901	640789	638935	641875	3608	0.56	-0.09	PASS
4-chloroaniline	364745	372864	372918	370176	4703	1.27	378204	377330	375425	376986	1421	0.38	-1.84	PASS
hexachlorobutadiene	130937	133892	133787	132872	1677	1.26	134118	131409	131706	132411	1486	1.12	0.35	PASS
2-methylnaphthalene	627310	643366	643371	638016	9271	1.45	650934	645347	644288	646856	3571	0.55	-1.39	PASS
4-chloro-3-methylphenol	363509	372177	373094	369593	5289	1.43	370470	367322	366997	368263	1918	0.52	0.36	PASS
1-methylnaphthalene	605609	618871	620998	615159	8339	1.36	629048	623146	622223	624806	3703	0.59	-1.57	PASS
1,2,4,5-Tetrachlorobenzene	241238	245964	245232	244145	2544	1.04	243630	242510	241372	242504	1129	0.47	0.67	PASS
hexachlorocyclopentadiene	116601	120591	119462	118885	2057	1.73	116395	116227	115501	116041	475	0.41	2.39	PASS
2,4,6-trichlorophenol	234196	239699	239132	237676	3027	1.27	234520	233876	232573	233656	992	0.42	1.69	PASS
2,4,5-trichlorophenol	228868	234109	233717	232231	2919	1.26	232088	231980	230147	231405	1091	0.47	0.36	PASS
2-chloronaphthalene	512484	523485	523317	519762	6303	1.21	521500	518502	516177	518726	2669	0.51	0.20	PASS
Biphenyl	649641	663720	663441	658934	8049	1.22	660072	656120	653319	656504	3393	0.52	0.37	PASS
2-nitroaniline	340217	347361	346641	344740	3933	1.14	343896	340835	339983	341571	2058	0.60	0.92	PASS
1,4-dinitrobenzene	653793	665831	665803	661809	6942	1.05	654825	649138	646348	650104	4320	0.66	1.77	PASS
acenaphthylene	275967	282442	280456	279622	3317	1.19	291597	289220	288137	289651	1770	0.61	-3.59	PASS
1,3-dinitrobenzene	334868	340560	340949	338792	3404	1.00	342476	339715	337599	339930	2446	0.72	-0.34	PASS
dimethylphthalate	300923	306038	306893	304618	3228	1.06	306806	304219	302295	304440	2264	0.74	0.06	PASS
2,6-dinitrotoluene	665538	678708	677791	674012	7353	1.09	667927	662763	659031	663240	4467	0.67	1.60	PASS

acenaphthene	328807	335342	334538	332896	3564	1.07	343085	340825	338826	340912	2131	0.63	-2.41	PASS
3-nitroaniline	403212	420082	423890	415726	11005	2.65	431022	430954	428441	430139	1471	0.34	-3.47	PASS
2,4-dinitrophenol	554539	566141	565341	562007	6480	1.15	565280	561042	558344	561555	3496	0.62	0.08	PASS
dibenzofuran	483565	494357	486748	488223	5545	1.14	487703	490962	482122	486929	4471	0.92	0.27	PASS
2,4-dinitrotoluene	411655	419196	423575	418142	6029	1.44	422350	415647	418494	418830	3364	0.80	-0.16	PASS
4-nitrophenol	196668	202431	201931	200343	3193	1.59	199923	197906	197053	198294	1474	0.74	1.02	PASS
2,3,4,6-tetrachlorophenol	657087	672353	669979	666473	8215	1.23	656335	653469	650119	653308	3111	0.48	1.98	PASS
Fluorene	450667	459779	458548	456331	4944	1.08	457658	454118	451948	454575	2882	0.63	0.38	PASS
4-chlorophenyl phenyl ether	951920	972861	972060	965614	11866	1.23	971107	961700	961700	964836	5431	0.56	0.08	PASS
Diethylphthalate	333909	345775	348870	342851	7897	2.30	353496	350700	342723	348973	5590	1.60	-1.79	PASS
4-nitroaniline	435113	442519	436284	437972	3981	0.91	430065	430765	427725	429518	1592	0.37	1.93	PASS
4,6-dinitro-2-methylphenol	528898	539833	538390	535707	5941	1.11	534553	531043	528139	531245	3212	0.60	0.83	PASS
azobenzene	367392	374621	372888	371634	3774	1.02	373338	370643	367579	370520	2881	0.78	0.30	PASS
4-bromophenyl phenyl ester	180139	182859	182855	181951	1569	0.86	181000	181096	179710	180602	774	0.43	0.74	PASS
hexachlorobenzene	349415	356979	355983	354126	4110	1.16	342216	338859	339126	340067	1866	0.55	3.97	PASS
pentachlorophenol	645355	657962	655899	653072	6762	1.04	655020	650361	645902	650428	4559	0.70	0.40	PASS
phenanthrene	662496	675758	672344	670199	6886	1.03	670067	665965	662600	666211	3740	0.56	0.60	PASS
anthracene	585500	598538	597645	593894	7283	1.23	597987	593007	590910	593968	3635	0.61	-0.01	PASS
carbazole	586303	599793	595925	594007	6947	1.17	603232	599000	595798	599343	3729	0.62	-0.90	PASS
di-n-butyl phthalate	410246	419359	416547	415384	4666	1.12	418935	416137	414843	416638	2092	0.50	-0.30	PASS
fluoranthene	654471	666002	664254	661576	6215	0.94	668549	666073	661948	665523	3335	0.50	-0.60	PASS
pyrene	650967	665685	661682	659445	7610	1.15	656710	654800	650047	653852	3431	0.52	0.85	PASS
benzyl butyl phthalate	450042	459622	457648	455771	5058	1.11	452073	448100	447363	449179	2534	0.56	1.45	PASS
benzo(a) anthracene	635549	647464	650178	644397	7782	1.21	645231	642038	635647	640972	4880	0.76	0.53	PASS
chrysene	657355	670946	671586	666629	8038	1.21	642404	638698	633711	638271	4362	0.68	4.25	PASS
bis(2-ethylhexyl) phthalate	463772	473038	471516	469442	4969	1.06	472680	468287	465366	468778	3682	0.79	0.14	PASS
di-n-octyl phthalate	462265	470733	469244	467414	4521	0.97	458612	456339	453427	456126	2599	0.57	2.41	PASS
benzo(b) fluoroanthene	601072	611440	610799	607770	5810	0.96	612725	608324	609303	610117	2311	0.38	-0.39	PASS
benzo(k) fluoroanthene	637653	651057	648645	645785	7145	1.11	667310	660680	657980	661990	4801	0.73	-2.51	PASS
benzo(a) pyrene	598863	608918	610923	606235	6462	1.07	618065	613693	610326	614028	3880	0.63	-1.29	PASS
indeno(1,2,3-cd) pyrene	463837	479874	471869	471860	8019	1.70	504977	499671	499953	501534	2985	0.60	-6.29	PASS
dibenz(a,h) anthracene	578298	595218	585868	586461	8476	1.45	596894	596078	589414	594129	4103	0.69	-1.31	PASS
benzo(g,h,i) perylene	532334	548831	541334	540833	8260	1.53	543819	533453	532600	536624	6246	1.16	0.78	PASS

TestAmerica Laboratories
ICV, ICal Verification Report

Data Path: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0576.D
Worklist Name: 080814-HSL Worklist Num: 26258
Instrument: SMS_Y Method: SMSY_8270C
Limit Group: MSSV - 8270C_625
Analysis Type: SemiVOA
Inj Volume: 0.50 Inj Vol Units: ul

Detector 1: MS SCAN

Compound	Amount Added	Amount Recovered	%Drift
27 1,4-Dioxane	100.0	97.3	-2.66
28 N-Nitrosodimethylamine	100.0	96.2	-3.83
29 Pyridine	100.0	104.4	4.40
40 Phenol	100.0	97.4	-2.58
43 Aniline	100.0	94.0	-5.96
44 Alpha Methyl Styrene	100.0	101.2	1.20
45 Bis(2-chloroethyl)ether	100.0	105.2	5.20
47 n-Decane	100.0	97.4	-2.63
48 2-Chlorophenol	100.0	100.0	0.00
49 1,3-Dichlorobenzene	100.0	99.2	-0.80
50 1,4-Dichlorobenzene	100.0	102.4	2.40
51 Benzyl alcohol	100.0	103.8	3.80
52 1,2-Dichlorobenzene	100.0	97.4	-2.65
54 2-Methylphenol	100.0	101.0	1.00
55 2,2'-oxybis[1-chloropropane]	72.0	72.6	0.82
56 Indene	100.0	100.5	0.50
58 3-Methylphenol	100.0	101.1	1.10
59 4-Methylphenol	100.0	101.1	1.10
60 3 & 4 Methylphenol	100.0	101.1	1.10
61 N-Nitrosodi-n-propylamine	100.0	104.3	4.30
63 Acetophenone	100.0	100.2	0.20
66 Hexachloroethane	100.0	98.8	-1.21
67 Nitrobenzene	100.0	97.4	-2.59
71 Isophorone	100.0	108.5	8.50
73 2,4-Dimethylphenol	100.0	93.0	-7.01
74 2-Nitrophenol	100.0	103.5	3.50
77 Benzoic acid	200.0	203.2	1.60
78 Bis(2-chloroethoxy)methane	100.0	95.0	-5.00
79 3,5-Dimethylphenol	100.0	98.9	-1.10

Compound	Amount Added	Amount Recovered	%Drift
82 2,4-Dichlorophenol	100.0	98.2	-1.79
84 1,2,4-Trichlorobenzene	100.0	99.5	-0.47
87 Naphthalene	100.0	98.3	-1.68
88 4-Chloroaniline	100.0	95.9	-4.12
91 Hexachlorobutadiene	100.0	99.4	-0.63
97 4-Chloro-3-methylphenol	100.0	100.1	0.10
101 2-Methylnaphthalene	100.0	98.7	-1.26
103 1-Methylnaphthalene	100.0	97.7	-2.29
105 Hexachlorocyclopentadiene	100.0	112.7	12.70
106 1,2,4,5-Tetrachlorobenzene	100.0	98.1	-1.89
108 2,4,6-Trichlorophenol	100.0	103.6	3.60
109 2,3-Dichlorobenzenamine	100.0	96.5	-3.50
110 2,4,5-Trichlorophenol	100.0	101.0	1.00
112 1,1'-Biphenyl	100.0	96.7	-3.32
114 2-Chloronaphthalene	100.0	96.0	-4.05
116 2-Nitroaniline	100.0	101.1	1.10
119 Dimethyl phthalate	100.0	97.3	-2.66
120 1,3-Dinitrobenzene	100.0	99.3	-0.67
121 2,6-Dinitrotoluene	100.0	105.5	5.50
122 Acenaphthylene	100.0	104.0	4.00
123 3-Nitroaniline	100.0	97.2	-2.82
124 Acenaphthene	100.0	97.8	-2.21
126 2,4-Dinitrophenol	200.0	188.8	-5.60
127 4-Nitrophenol	200.0	207.0	3.50
130 2,4-Dinitrotoluene	100.0	104.9	4.90
132 Dibenzofuran	100.0	94.6	-5.43
134 2,3,4,6-Tetrachlorophenol	100.0	103.9	3.90
136 Hexadecane	100.0	97.3	-2.70
139 Diethyl phthalate	100.0	96.2	-3.78
142 4-Chlorophenyl phenyl ether	100.0	97.2	-2.80
143 Fluorene	100.0	100.0	0.00
146 4-Nitroaniline	100.0	101.2	1.20
147 4,6-Dinitro-2-methylphenol	200.0	202.7	1.35
150 N-Nitrosodiphenylamine	100.0	94.5	-5.52
151 1,2-Diphenylhydrazine	101.1	100.2	-0.89
152 Azobenzene	100.0	99.1	-0.90

Compound	Amount Added	Amount Recovered	%Drift
164 4-Bromophenyl phenyl ether	100.0	98.4	-1.61
165 Hexachlorobenzene	100.0	94.7	-5.27
169 n-Octadecane	100.0	103.0	3.00
171 Pentachlorophenol	200.0	202.5	1.25
176 Phenanthrene	100.0	95.9	-4.11
177 Anthracene	100.0	100.2	0.20
178 Carbazole	100.0	96.6	-3.39
179 Alachlor	100.0	102.5	2.50
181 Di-n-butyl phthalate	100.0	105.1	5.10
188 Fluoranthene	100.0	102.3	2.30
191 Pyrene	100.0	99.8	-0.19
198 Butyl benzyl phthalate	100.0	94.9	-5.10
203 Bis(2-ethylhexyl) phthalate	100.0	91.4	-8.62
206 Benzo[a]anthracene	100.0	100.9	0.90
207 Chrysene	100.0	103.5	3.50
209 Di-n-octyl phthalate	100.0	95.0	-5.04
211 Benzo[b]fluoranthene	100.0	96.7	-3.31
212 Benzo[k]fluoranthene	100.0	105.4	5.40
214 Benzo[e]pyrene	100.0	118.7	18.70
215 Benzo[a]pyrene	100.0	98.1	-1.94
221 Dibenz(a,h)anthracene	100.0	102.0	2.00
222 Indeno[1,2,3-cd]pyrene	100.0	94.8	-5.20
223 Benzo[g,h,i]perylene	100.0	107.0	7.00

Mega Mix ICV

Standard Verification Form

Verification (New vendor or problematic Standard)	☐	Re-Verification	☑
TALS Reagent Record			
New	☐	Copied	☐
COA Reviewed against formulary report		☑	

Document instrument verification if need (Initial or re-verification):		
Department	**Acceptance Criteria**	
	Standard Analytes	**Poor Performers* and Esterified Analytes**
GC/HPLC	≤ 15 %D	≤ 35 %D or ≤ 50 %D for dinoseb
GCMS/LCMS ✓	≤ 35 %D	≤ 55 %D
MSVOA	≤ 25 %D	≤ 55 %D
Metals	≤ 8 %D	NA
Wet Chemistry	≤ 5 %D	NA

Standard Name	MS-567672 SEC	Standard ID		
Verified by	M. Hoffman	Instrument ID	Y	
Verification Date	8/14/14	Method Reference	8270C	
Reference Standard ID	MS-567672	Batch #	A093709	
Analyte/Mix	Prepared Concentration	Verification Concentration	% Diff	Pass/Fail
See attached report				All compounds pass.
	2/2015 (ER)	08/28/14		
New Expiration Date:	~~May 2015~~	New TALS ID	Same w/ new exp. date.	

New expiration date can be no greater than ½ the designated standards shelf life from the date of re-verification. Standards can only be re-verified one time.

Comment:	

1st Level Review: MH Date: 8/28/14

2nd Level Review: (ER) Date: 08/28/14

QA Review (Re-verification only)		Date: 8/28/14

Attach form, supporting documentation and original CoA to new verified or re-verified standard record in TALS.

*See analytical SOP for details on poor performing analytes.

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Certificate #3282.02

FOR LABORATORY USE ONLY-READ MSDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567672.sec **Lot No.:** A094002

Description : 8270 List 1 / Std #1 MegaMix

8270 List 1 / Std #1 MegaMix 500-2000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : September 2014 **Storage:** 10°C or colder

Handling: Sonication required. Mix is photosensitive.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	1,4-Dioxane CAS # 123-91-1.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
2	Pyridine CAS # 110-86-1.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
3	N-Nitrosodimethylamine CAS # 62-75-9.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
4	Aniline CAS # 62-53-3.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
5	Phenol CAS # 108-95-2.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
6	Bis(2-chloroethyl)ether CAS # 111-44-4.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
7	2-Chlorophenol CAS # 95-57-8.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
8	1,3-Dichlorobenzene CAS # 541-73-1.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed
9	1,4-Dichlorobenzene CAS # 106-46-7.SEC Purity 99%	1,000.0 µg/mL	+/- 5.8141 +/- 8.7957 +/- 17.3886	µg/mL µg/mL µg/mL	Gravimetric Unstressed Stressed

10	1,2-Dichlorobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-50-1.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
11	Benzyl alcohol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 100-51-6.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
12	Bis(2-chloroisopropyl)ether	720.0	µg/mL	+/-	4.1861	µg/mL	Gravimetric
	CAS # 108-60-1.SEC			+/-	6.3329	µg/mL	Unstressed
	Purity 72%			+/-	12.5198	µg/mL	Stressed
13	2-Methylphenol (o-cresol)	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 95-48-7.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
14	Acetophenone	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 98-86-2.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
15	Hexachloroethane	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 67-72-1.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
16	N-Nitroso-di-n-propylamine	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 621-64-7.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
17	4-Methylphenol (p-cresol)	500.0	µg/mL	+/-	2.9138	µg/mL	Gravimetric
	CAS # 106-44-5.SEC			+/-	4.4023	µg/mL	Unstressed
	Purity 99%			+/-	8.6966	µg/mL	Stressed
18	3-Methylphenol (m-cresol)	500.0	µg/mL	+/-	2.9138	µg/mL	Gravimetric
	CAS # 108-39-4.SEC			+/-	4.4023	µg/mL	Unstressed
	Purity 99%			+/-	8.6966	µg/mL	Stressed
19	n-Decane (C10)	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 124-18-5.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
20	n-Octadecane (C18)	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 593-45-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
21	Nitrobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 98-95-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
22	Isophorone	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 78-59-1.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
23	2-Nitrophenol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 88-75-5.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
24	2,4-Dimethylphenol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 105-67-9.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
25	Bis(2-chloroethoxy)methane	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 111-91-1			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
26	2,4-Dichlorophenol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-83-2.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
27	1,2,4-Trichlorobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-82-1.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
28	Naphthalene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-20-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed

29	4-Chloroaniline	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 106-47-8.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
30	Hexachlorobutadiene	1,000.0	µg/mL	+/-	5.8139	µg/mL	Gravimetric
	CAS # 87-68-3.SEC			+/-	8.7954	µg/mL	Unstressed
	Purity 97%			+/-	17.3881	µg/mL	Stressed
31	2-Methylnaphthalene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-57-6.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
32	4-Chloro-3-methylphenol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 59-50-7.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
33	1-Methylnaphthalene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 90-12-0.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
34	1,2,4,5-Tetrachlorobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-94-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
35	Hexachlorocyclopentadiene	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 77-47-4.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
36	2,4,6-Trichlorophenol	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 88-06-2.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
37	2,4,5-Trichlorophenol	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-95-4.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
38	2-Chloronaphthalene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-58-7.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
39	Biphenyl	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 92-52-4.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
40	2-Nitroaniline	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 88-74-4.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
41	Acenaphthylene	1,000.0	µg/mL	+/-	5.8139	µg/mL	Gravimetric
	CAS # 208-96-8.SEC			+/-	8.7954	µg/mL	Unstressed
	Purity 97%			+/-	17.3881	µg/mL	Stressed
42	1,3-Dinitrobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 99-65-0.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
43	Dimethylphthalate	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 131-11-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
44	2,6-Dinitrotoluene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 606-20-2.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
45	Acenaphthene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 83-32-9.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
46	3-Nitroaniline	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 99-09-2.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
47	2,4-Dinitrophenol	2,000.0	µg/mL	+/-	11.6282	µg/mL	Gravimetric
	CAS # 51-28-5.SEC			+/-	17.5913	µg/mL	Unstressed
	Purity 99%			+/-	34.7772	µg/mL	Stressed

48	Dibenzofuran	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 132-64-9.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
49	2,4-Dinitrotoluene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 121-14-2.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
50	4-Nitrophenol	2,000.0	µg/mL	+/-	11.6282	µg/mL	Gravimetric
	CAS # 100-02-7.SEC			+/-	17.5913	µg/mL	Unstressed
	Purity 99%			+/-	34.7772	µg/mL	Stressed
51	2,3,4,6-Tetrachlorophenol	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 58-90-2.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
52	Fluorene	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 86-73-7.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
53	4-Chlorophenyl phenyl ether	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 7005-72-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
54	Diethylphthalate	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 84-66-2.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
55	4-Nitroaniline	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 100-01-6.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
56	4,6-Dinitro-2-methylphenol (Dinitro-o-cresol)	2,000.0	µg/mL	+/-	11.6281	µg/mL	Gravimetric
	CAS # 534-52-1.SEC			+/-	17.5912	µg/mL	Unstressed
	Purity 98%			+/-	34.7769	µg/mL	Stressed
57	Azobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 103-33-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
58	4-Bromophenyl phenyl ether	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 101-55-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
59	Hexachlorobenzene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 118-74-1.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
60	Pentachlorophenol	2,000.0	µg/mL	+/-	11.6282	µg/mL	Gravimetric
	CAS # 87-86-5.SEC			+/-	17.5913	µg/mL	Unstressed
	Purity 99%			+/-	34.7772	µg/mL	Stressed
61	Phenanthrene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 85-01-8.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
62	Anthracene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-12-7.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
63	n-Hexadecane (C16)	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 544-76-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
64	Carbazole	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 86-74-8.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
65	Di-n-butylphthalate	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 84-74-2.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
66	Fluoranthene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 206-44-0.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed

67	Pyrene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 129-00-0.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
68	Benzyl butyl phthalate	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 85-68-7.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
69	Benz(a)anthracene	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 56-55-3.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
70	chrysene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 218-01-9.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
71	Bis(2-ethylhexyl)phthalate	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 117-81-7.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
72	Di-n-octyl phthalate	1,000.0	µg/mL	+/-	5.8140	µg/mL	Gravimetric
	CAS # 117-84-0.SEC			+/-	8.7956	µg/mL	Unstressed
	Purity 98%			+/-	17.3885	µg/mL	Stressed
73	Benzo(b)fluoranthene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 205-99-2.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
74	Benzo(k)fluoranthene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 207-08-9.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
75	Benzo(a)pyrene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 50-32-8.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
76	Indeno(1,2,3-cd)pyrene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 193-39-5.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
77	Dibenz(a,h)anthracene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 53-70-3.SEC			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed
78	Benzo(g,h,i)perylene	1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 191-24-2			+/-	8.7957	µg/mL	Unstressed
	Purity 99%			+/-	17.3886	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Specific Reference Material Notes:
The Bis(2-chloroisopropyl)ether contains a 28% impurity of Propane, 1,1'oxybis,, 3-chloro.

Report Date: 14-Aug-2014 06:29:22 Chrom Revision: 2.2 24-Jun-2014 07:21:42

Preliminary Report

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0571.D
Lims ID: ICIS HSLA080
Client ID:
Sample Type: ICIS Calib Level: 5
Inject. Date: 08-Aug-2014 09:55:30 ALS Bottle#: 5 Worklist Smp#: 22
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: ICIS HSLA080
Operator ID: hoffmanm Instrument ID: SMS_Y
Sublist: chrom-SMSY_8270C*sub20

Method: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\SMSY_8270C.m
Limit Group: MSSV - 8270C_625
Method Label: 8270C / 625
Last Update: 12-Aug-2014 08:24:06 Calib Date: 08-Aug-2014 11:47:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0575.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK049

First Level Reviewer: hoffmanm Date: 11-Aug-2014 11:55:22

	Compound	Sig	RT (min.)	Adj RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
*	1 1,4-Dichlorobenzene-d4	152	4.357	4.357	0.000	96	208751	40.0	40.0	
*	2 Naphthalene-d8	136	5.391	5.391	0.000	100	830145	40.0	40.0	
*	3 Acenaphthene-d10	164	6.865	6.865	0.000	92	463953	40.0	40.0	
*	4 Phenanthrene-d10	188	8.128	8.128	0.000	97	764844	40.0	40.0	
*	5 Chrysene-d12	240	11.870	11.870	0.000	96	722208	40.0	40.0	
*	6 Perylene-d12	264	15.964	15.964	0.000	95	563515	40.0	40.0	
$	7 2-Fluorophenol	112	3.335	3.335	0.000	93	574536	80.0	82.3	
$	8 Phenol-d5	99	4.016	4.016	0.000	93	752107	80.0	83.1	
$	9 Nitrobenzene-d5	82	4.797	4.797	0.000	91	675640	80.0	83.2	
$	11 2-Fluorobiphenyl	172	6.266	6.266	0.000	99	1196587	80.0	79.0	
$	12 2,4,6-Tribromophenol	330	7.535	7.535	0.000	91	135458	80.0	83.7	p
$	13 Terphenyl-d14	244	9.826	9.826	0.000	98	1140380	80.0	82.6	
	27 1,4-Dioxane	88	2.177	2.170	0.000	96	261983	80.0	77.7	
	28 N-Nitrosodimethylamine	74	2.389	2.389	0.000	91	395689	80.0	82.1	
	29 Pyridine	79	2.436	2.436	0.000	95	689782	80.0	82.8	
	40 Phenol	94	4.028	4.028	0.000	99	776997	80.0	81.7	
	43 Aniline	93	4.092	4.092	0.000	97	998990	80.0	84.4	
	44 Alpha Methyl Styrene	118	4.122	4.122	0.000	94	625968	80.0	82.6	
	45 Bis(2-chloroethyl)ether	93	4.116	4.116	0.000	91	581149	80.0	78.9	
	47 n-Decane	43	4.163	4.163	0.000	90	618794	80.0	80.6	
	48 2-Chlorophenol	128	4.186	4.186	0.000	97	602527	80.0	81.8	
	49 1,3-Dichlorobenzene	146	4.316	4.316	0.000	97	660171	80.0	81.3	
	50 1,4-Dichlorobenzene	146	4.368	4.368	0.000	93	667498	80.0	81.7	
	51 Benzyl alcohol	108	4.439	4.439	0.000	92	393559	80.0	84.3	
	52 1,2-Dichlorobenzene	146	4.492	4.492	0.000	97	637403	80.0	79.8	
	54 2-Methylphenol	108	4.509	4.509	0.000	96	593564	80.0	83.4	
	55 2,2'-oxybis[1-chloropropan	45	4.533	4.534	0.000	93	850367	80.0	81.9	
	56 Indene	116	4.562	4.562	0.000	90	994182	80.0	80.8	
	58 3-Methylphenol	108	4.627	4.627	0.000	96	606414	80.0	82.2	p

Preliminary Report

Data File: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0571.D

Compound	Sig	RT (min.)	Adj RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
59 4-Methylphenol	108	4.627	4.627	0.000	96	606414	80.0	82.2	p
60 3 & 4 Methylphenol	108	4.627	4.627	0.000	96	606414	80.0	82.2	p
61 N-Nitrosodi-n-propylamine	70	4.639	4.639	0.000	91	436567	80.0	83.8	
63 Acetophenone	105	4.662	4.662	0.000	95	834675	80.0	80.1	
66 Hexachloroethane	117	4.768	4.768	0.000	97	250173	80.0	78.4	
67 Nitrobenzene	77	4.815	4.815	0.000	89	659580	80.0	80.6	
71 Isophorone	82	4.991	4.991	0.000	99	1133316	80.0	84.1	
73 2,4-Dimethylphenol	107	5.062	5.062	0.000	96	606986	80.0	82.4	
74 2-Nitrophenol	139	5.073	5.073	0.000	94	324183	80.0	85.2	
77 Benzoic acid	105	5.109	5.109	0.000	88	885035	160.0	164.5	
78 Bis(2-chloroethoxy)methane	93	5.144	5.144	0.000	99	689223	80.0	79.8	
79 3,5-Dimethylphenol	107	5.167	5.167	0.000	94	631749	80.0	81.4	
82 2,4-Dichlorophenol	162	5.255	5.255	0.000	96	468005	80.0	80.3	
84 1,2,4-Trichlorobenzene	180	5.332	5.332	0.000	94	512810	80.0	79.6	
87 Naphthalene	128	5.408	5.408	0.000	98	1733746	80.0	79.8	
88 4-Chloroaniline	127	5.438	5.438	0.000	95	762624	80.0	81.7	
91 Hexachlorobutadiene	225	5.485	5.485	0.000	97	266105	80.0	80.2	
94 Caprolactam	55	5.720	5.723	0.000	79	290670	80.0	84.0	
97 4-Chloro-3-methylphenol	107	5.808	5.808	0.000	97	519526	80.0	82.8	
101 2-Methylnaphthalene	142	5.978	5.978	0.000	93	1158573	80.0	80.9	
103 1-Methylnaphthalene	142	6.066	6.066	0.000	94	1088630	80.0	80.2	
105 Hexachlorocyclopentadiene	237	6.096	6.096	0.000	95	260664	80.0	83.7	
106 1,2,4,5-Tetrachlorobenzene	216	6.119	6.119	0.000	98	439813	80.0	79.2	
108 2,4,6-Trichlorophenol	196	6.207	6.207	0.000	94	311636	80.0	84.4	
109 2,3-Dichlorobenzenamine	161	6.219	6.219	0.000	96	581729	80.0	79.1	
110 2,4,5-Trichlorophenol	196	6.236	6.236	0.000	92	335663	80.0	83.5	
112 1,1'-Biphenyl	154	6.360	6.360	0.000	95	1358617	80.0	79.4	
114 2-Chloronaphthalene	162	6.395	6.395	0.000	98	1028753	80.0	79.2	
116 2-Nitroaniline	65	6.477	6.477	0.000	83	369274	80.0	86.8	
119 Dimethyl phthalate	163	6.589	6.589	0.000	98	1151460	80.0	81.4	
120 1,3-Dinitrobenzene	168	6.648	6.648	0.000	83	192745	80.0	86.1	
121 2,6-Dinitrotoluene	165	6.665	6.665	0.000	94	273407	80.0	86.7	
122 Acenaphthylene	152	6.753	6.753	0.000	99	1638049	80.0	82.9	
123 3-Nitroaniline	138	6.818	6.818	0.000	93	354630	80.0	85.5	
124 Acenaphthene	153	6.894	6.894	0.000	94	1073421	80.0	80.3	
126 2,4-Dinitrophenol	184	6.900	6.900	0.000	80	322324	160.0	158.6	
127 4-Nitrophenol	109	6.930	6.930	0.000	88	400576	160.0	176.3	
130 2,4-Dinitrotoluene	165	7.006	7.006	0.000	91	356996	80.0	85.6	
132 Dibenzofuran	168	7.035	7.035	0.000	96	1460457	80.0	78.5	
134 2,3,4,6-Tetrachlorophenol	232	7.129	7.129	0.000	81	258810	80.0	85.8	
136 Hexadecane	57	7.135	7.135	0.000	94	784967	80.0	82.3	
139 Diethyl phthalate	149	7.171	7.171	0.000	98	1175652	80.0	80.8	
142 4-Chlorophenyl phenyl ethe	204	7.300	7.300	0.000	97	513056	80.0	79.0	
143 Fluorene	166	7.323	7.323	0.000	95	1200615	80.0	80.2	
146 4-Nitroaniline	138	7.341	7.341	0.000	85	356465	80.0	86.8	
147 4,6-Dinitro-2-methylphenol	198	7.353	7.353	0.000	84	405535	160.0	162.6	
150 N-Nitrosodiphenylamine	169	7.400	7.400	0.000	62	878554	80.0	82.0	
151 1,2-Diphenylhydrazine	77	7.435	7.435	0.000	99	1346090	80.9	82.1	p
152 Azobenzene	77	7.435	7.435	0.000	100	1346090	80.0	81.2	p
155 2,4,6 - Tribromophenol	330	7.535	7.535	0.000	91	136633	80.0	84.4	p
164 4-Bromophenyl phenyl ether	248	7.717	7.717	0.000	72	278581	80.0	80.8	

Preliminary Report

Data File: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0571.D

Compound	Sig	RT (min.)	Adj RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
165 Hexachlorobenzene	284	7.793	7.793	0.000	90	279985	80.0	78.6	
167 Atrazine	200	7.823	7.823	0.000	90	257443	80.0	85.9	
169 n-Octadecane	85	7.905	7.905	0.000	94	328683	80.0	85.9	
171 Pentachlorophenol	266	7.952	7.952	0.000	89	372171	160.0	162.4	
176 Phenanthrene	178	8.146	8.146	0.000	98	1721546	80.0	78.5	
177 Anthracene	178	8.193	8.193	0.000	98	1777401	80.0	82.8	
178 Carbazole	167	8.322	8.322	0.000	96	1689104	80.0	83.5	
179 Alachlor	188	8.387	8.387	0.000	97	211290	80.0	86.4	
181 Di-n-butyl phthalate	149	8.551	8.551	0.000	100	1921614	80.0	87.2	
188 Fluoranthene	202	9.356	9.356	0.000	99	1712847	80.0	85.3	
191 Pyrene	202	9.667	9.667	0.000	97	1801637	80.0	83.0	
197 Famphur	218	10.513	10.514	0.000	87	553352	80.0	83.8	
198 Butyl benzyl phthalate	149	10.595	10.595	0.000	98	823791	80.0	82.3	
203 Bis(2-ethylhexyl) phthalat	149	11.782	11.782	0.000	98	1007930	80.0	79.4	
205 3,3'-Dichlorobenzidine	252	11.794	11.794	0.000	76	546183	80.0	82.0	
206 Benzo[a]anthracene	228	11.847	11.847	0.000	99	1575300	80.0	83.4	
207 Chrysene	228	11.935	11.935	0.000	98	1485182	80.0	79.7	
209 Di-n-octyl phthalate	149	13.574	13.574	0.000	99	1436762	80.0	78.6	
211 Benzo[b]fluoranthene	252	14.760	14.760	0.000	99	1286907	80.0	80.6	
212 Benzo[k]fluoranthene	252	14.842	14.842	0.000	98	1490888	80.0	89.1	
214 Benzo[e]pyrene	252	15.612	15.612	0.000	97	1160559	80.0	86.7	
215 Benzo[a]pyrene	252	15.782	15.782	0.000	80	1332291	80.0	81.8	
221 Dibenz(a,h)anthracene	278	19.372	19.372	0.000	94	1030012	80.0	79.6	
222 Indeno[1,2,3-cd]pyrene	276	19.307	19.306	0.000	96	958528	80.0	76.2	M
223 Benzo[g,h,i]perylene	276	19.842	19.842	0.000	95	1208264	80.0	87.7	
S 225 Methyl Phenols,Total	108				0			165.6	
S 226 Total Cresols	108				0			165.6	

QC Flag Legend

Processing Flags
p - Peak ID'ed as Multiple Compounds

Review Flags
M - Manually Integrated

Reagents:

MS-HSLA080_00009 Amount Added: 200.00 Units: uL

Preliminary Report

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0576.D
Lims ID: ICV HSLB1B3
Client ID:
Sample Type: ICV
Inject. Date: 08-Aug-2014 12:15:30 ALS Bottle#: 10 Worklist Smp#: 27
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: ICV HSLB1B3
Operator ID: hoffmanm Instrument ID: SMS_Y
Sublist:

Method: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\SMSY_8270C.m
Limit Group: MSSV - 8270C_625
Method Label: 8270C / 625
Last Update: 12-Aug-2014 08:24:06 Calib Date: 08-Aug-2014 11:47:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0575.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK049

First Level Reviewer: hoffmanm Date: 11-Aug-2014 12:37:52

	Compound	Sig	RT (min.)	Adj RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
*	1 1,4-Dichlorobenzene-d4	152	4.358	4.357	0.001	96	205437	40.0	40.0	
*	2 Naphthalene-d8	136	5.392	5.391	0.001	100	816575	40.0	40.0	
*	3 Acenaphthene-d10	164	6.867	6.865	0.002	91	466874	40.0	40.0	
*	4 Phenanthrene-d10	188	8.130	8.128	0.002	97	771135	40.0	40.0	
*	5 Chrysene-d12	240	11.872	11.870	0.002	96	718207	40.0	40.0	
*	6 Perylene-d12	264	15.960	15.964	-0.004	95	568555	40.0	40.0	
	27 1,4-Dioxane	88	2.179	2.170	0.002	97	322864	100.0	97.3	
	28 N-Nitrosodimethylamine	74	2.384	2.389	-0.005	89	456191	100.0	96.2	
	29 Pyridine	79	2.431	2.436	-0.005	93	856044	100.0	104.4	
	40 Phenol	94	4.029	4.028	0.001	99	912317	100.0	97.4	
	43 Aniline	93	4.094	4.092	0.002	97	1095396	100.0	94.0	
	44 Alpha Methyl Styrene	118	4.117	4.122	-0.005	97	754840	100.0	101.2	
	45 Bis(2-chloroethyl)ether	93	4.117	4.116	0.001	92	762976	100.0	105.2	
	47 n-Decane	43	4.164	4.163	0.001	89	735348	100.0	97.4	
	48 2-Chlorophenol	128	4.188	4.186	0.002	97	725564	100.0	100.0	
	49 1,3-Dichlorobenzene	146	4.317	4.316	0.001	98	793068	100.0	99.2	
	50 1,4-Dichlorobenzene	146	4.370	4.368	0.002	93	822895	100.0	102.4	
	51 Benzyl alcohol	108	4.440	4.439	0.001	92	476820	100.0	103.8	
	52 1,2-Dichlorobenzene	146	4.493	4.492	0.001	96	764893	100.0	97.4	
	54 2-Methylphenol	108	4.511	4.509	0.002	95	707591	100.0	101.0	
	55 2,2'-oxybis[1-chloropropan	45	4.534	4.534	0.001	94	741917	72.0	72.6	M
	56 Indene	116	4.564	4.562	0.002	90	1217387	100.0	100.5	
	58 3-Methylphenol	108	4.628	4.627	0.001	96	733385	100.0	101.1	p
	59 4-Methylphenol	108	4.628	4.627	0.001	96	733385	100.0	101.1	p
	60 3 & 4 Methylphenol	108	4.628	4.627	0.001	96	733385	100.0	101.1	p
	61 N-Nitrosodi-n-propylamine	70	4.640	4.639	0.001	90	535145	100.0	104.3	
	63 Acetophenone	105	4.664	4.662	0.002	96	1027514	100.0	100.2	
	66 Hexachloroethane	117	4.769	4.768	0.001	97	310425	100.0	98.8	
	67 Nitrobenzene	77	4.816	4.815	0.001	90	784602	100.0	97.4	

Preliminary Report

Data File: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0576.D

Compound	Sig	RT (min.)	Adj RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
71 Isophorone	82	4.987	4.991	-0.004	99	1437028	100.0	108.5	
73 2,4-Dimethylphenol	107	5.063	5.062	0.001	95	673746	100.0	93.0	
74 2-Nitrophenol	139	5.075	5.073	0.002	87	387526	100.0	103.5	
77 Benzoic acid	105	5.116	5.109	0.007	89	1082193	200.0	203.2	
78 Bis(2-chloroethoxy)methane	93	5.145	5.144	0.001	98	807079	100.0	95.0	
79 3,5-Dimethylphenol	107	5.169	5.167	0.002	93	755199	100.0	98.9	
82 2,4-Dichlorophenol	162	5.257	5.255	0.002	96	563277	100.0	98.2	
84 1,2,4-Trichlorobenzene	180	5.333	5.332	0.001	94	630381	100.0	99.5	
87 Naphthalene	128	5.410	5.408	0.002	98	2100466	100.0	98.3	
88 4-Chloroaniline	127	5.439	5.438	0.001	96	880373	100.0	95.9	
91 Hexachlorobutadiene	225	5.486	5.485	0.001	97	324239	100.0	99.4	
97 4-Chloro-3-methylphenol	107	5.809	5.808	0.001	97	617768	100.0	100.1	
101 2-Methylnaphthalene	142	5.980	5.978	0.002	93	1391634	100.0	98.7	
103 1-Methylnaphthalene	142	6.068	6.066	0.002	94	1305488	100.0	97.7	
105 Hexachlorocyclopentadiene	237	6.097	6.096	0.001	95	353292	100.0	112.7	
106 1,2,4,5-Tetrachlorobenzene	216	6.115	6.119	-0.004	97	536150	100.0	98.1	
108 2,4,6-Trichlorophenol	196	6.203	6.207	-0.004	95	384640	100.0	103.6	
109 2,3-Dichlorobenzenamine	161	6.220	6.219	0.001	96	714148	100.0	96.5	
110 2,4,5-Trichlorophenol	196	6.238	6.236	0.002	95	408739	100.0	101.0	
112 1,1'-Biphenyl	154	6.361	6.360	0.001	95	1663971	100.0	96.7	
114 2-Chloronaphthalene	162	6.397	6.395	0.002	98	1254267	100.0	95.9	
116 2-Nitroaniline	65	6.479	6.477	0.002	85	432654	100.0	101.1	
119 Dimethyl phthalate	163	6.590	6.589	0.001	98	1385368	100.0	97.3	
120 1,3-Dinitrobenzene	168	6.649	6.648	0.001	83	223654	100.0	99.3	
121 2,6-Dinitrotoluene	165	6.667	6.665	0.002	94	334796	100.0	105.5	
122 Acenaphthylene	152	6.749	6.753	-0.004	99	2068095	100.0	104.0	
123 3-Nitroaniline	138	6.820	6.818	0.002	94	405573	100.0	97.2	
124 Acenaphthene	153	6.896	6.894	0.002	93	1315482	100.0	97.8	
126 2,4-Dinitrophenol	184	6.902	6.900	0.002	82	389678	200.0	188.8	
127 4-Nitrophenol	109	6.925	6.930	-0.005	89	473433	200.0	207.0	
130 2,4-Dinitrotoluene	165	7.008	7.006	0.002	91	440180	100.0	104.9	
132 Dibenzofuran	168	7.037	7.035	0.002	96	1770687	100.0	94.6	
134 2,3,4,6-Tetrachlorophenol	232	7.131	7.129	0.002	75	315432	100.0	103.9	
136 Hexadecane	57	7.137	7.135	0.002	95	934416	100.0	97.3	
139 Diethyl phthalate	149	7.172	7.171	0.001	97	1408488	100.0	96.2	
142 4-Chlorophenyl phenyl ethe	204	7.301	7.300	0.001	96	635053	100.0	97.2	
143 Fluorene	166	7.325	7.323	0.002	95	1506037	100.0	100.0	
146 4-Nitroaniline	138	7.342	7.341	0.001	83	418243	100.0	101.2	
147 4,6-Dinitro-2-methylphenol	198	7.354	7.353	0.001	84	512742	200.0	202.7	
150 N-Nitrosodiphenylamine	169	7.401	7.400	0.001	63	1020409	100.0	94.5	
151 1,2-Diphenylhydrazine	77	7.436	7.435	0.001	99	1653504	101.1	100.2	p
152 Azobenzene	77	7.436	7.435	0.001	100	1653504	100.0	99.1	p
164 4-Bromophenyl phenyl ether	248	7.718	7.717	0.001	72	341920	100.0	98.4	
165 Hexachlorobenzene	284	7.795	7.793	0.002	89	340065	100.0	94.7	
169 n-Octadecane	85	7.906	7.905	0.001	94	397489	100.0	103.0	
171 Pentachlorophenol	266	7.953	7.952	0.001	89	470038	200.0	202.5	
176 Phenanthrene	178	8.147	8.146	0.001	98	2121286	100.0	95.9	
177 Anthracene	178	8.194	8.193	0.001	98	2166945	100.0	100.2	
178 Carbazole	167	8.323	8.322	0.001	96	1970178	100.0	96.6	
179 Alachlor	188	8.388	8.387	0.001	97	252689	100.0	102.5	
181 Di-n-butyl phthalate	149	8.553	8.551	0.001	100	2335645	100.0	105.1	

Preliminary Report

Data File: \\Denchrom\ChromData\SMS_Y\20140811-26258.b\Y0576.D

Compound	Sig	RT (min.)	Adj RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
188 Fluoranthene	202	9.357	9.356	0.001	99	2072557	100.0	102.3	
191 Pyrene	202	9.669	9.667	0.002	96	2155864	100.0	99.8	
198 Butyl benzyl phthalate	149	10.591	10.595	-0.004	97	948010	100.0	94.9	
203 Bis(2-ethylhexyl) phthalat	149	11.783	11.782	0.001	98	1161370	100.0	91.4	
206 Benzo[a]anthracene	228	11.842	11.847	-0.005	99	1896853	100.0	100.9	
207 Chrysene	228	11.936	11.935	0.001	98	1917451	100.0	103.5	
209 Di-n-octyl phthalate	149	13.581	13.574	0.007	99	1779825	100.0	95.0	
211 Benzo[b]fluoranthene	252	14.762	14.760	0.002	98	1563440	100.0	96.7	
212 Benzo[k]fluoranthene	252	14.844	14.842	0.002	99	1780662	100.0	105.4	
214 Benzo[e]pyrene	252	15.619	15.612	0.007	96	1603512	100.0	118.7	
215 Benzo[a]pyrene	252	15.778	15.782	-0.004	80	1618192	100.0	98.1	
221 Dibenz(a,h)anthracene	278	19.373	19.372	0.001	94	1339673	100.0	102.0	
222 Indeno[1,2,3-cd]pyrene	276	19.309	19.306	0.002	96	1196951	100.0	94.8	
223 Benzo[g,h,i]perylene	276	19.843	19.842	0.001	94	1486786	100.0	107.0	

QC Flag Legend

Processing Flags
p - Peak ID'ed as Multiple Compounds

Review Flags
M - Manually Integrated

Reagents:

MS-HSLB1B3SSV_00011 Amount Added: 200.00 Units: uL

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FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567673.SEC **Lot No.:** A0101015

Description : 8270 List 1 / Std #2 Amines

8270 List 1 / Std #2 Amines 2,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 10 mL **Pkg Amt:** > 5 mL

Expiration Date : July 31, 2015 **Storage:** 10°C or colder

Handling: Contains carcinogen

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	epsilon-Caprolactam	2,010.0 µg/mL	+/- 11.7958	µg/mL	Gravimetric
	CAS # 105-60-2.SEC (Lot BLJTB)		+/- 22.0652	µg/mL	Unstressed
	Purity 99%		+/- 37.3617	µg/mL	Stressed
2	Atrazine	2,009.3 µg/mL	+/- 11.7919	µg/mL	Gravimetric
	CAS # 1912-24-9.SEC (Lot 1132400)		+/- 22.0579	µg/mL	Unstressed
	Purity 99%		+/- 37.3493	µg/mL	Stressed
3	Benzidine	2,018.0 µg/mL	+/- 11.8428	µg/mL	Gravimetric
	CAS # 92-87-5.SEC (Lot 1301900)		+/- 22.1530	µg/mL	Unstressed
	Purity 99%		+/- 37.5104	µg/mL	Stressed
4	3,3'-Dichlorobenzidine	2,016.7 µg/mL	+/- 11.8349	µg/mL	Gravimetric
	CAS # 91-94-1.SEC (Lot 2010900)		+/- 22.1384	µg/mL	Unstressed
	Purity 99%		+/- 37.4856	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Column:
30m x 0.25mm x 0.25µm
Rtx-5 (cat.#10223)

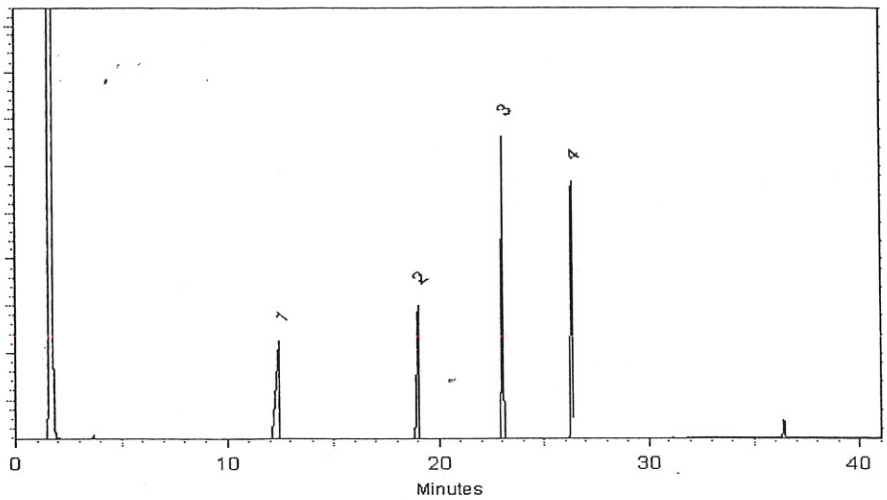
Carrier Gas:
hydrogen-constant pressure 10 psi.

Temp. Program:
40°C (hold 2 min.) to 330°C
@ 10°C/min. (hold 10 min.)

Inj. Temp:
250°C

Det. Temp:
330°C

Det. Type:
FID



This chromatogram represents a general set of testing conditions chosen to guarantee product quality. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

Brandon Reish - Mix Technician

Date Mixed: 30-Jan-2014 **Balance:** B345965662

Jodi E. Breon - QA Analyst

Date Passed: 03-Feb-2014

Manufactured under Restek's ISO 9001:2008
Registered Quality System
Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31840, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

Analytical Reference Materials

8270 List 1 / Std #3 Benzoic Acid

Catalog # 567674.sec

Lot # A093654 & A093441

110 Benner Circle Bellefonte, PA 16823-8812
(814) 353-1300

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RAW MATERIAL TEST INFORMATION AVAILABLE UPON REQUEST

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This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567674.sec **Lot No.:** A093654

Description : 8270 List 1 / Std #3 Benzoic Acid

8270 List 1 / Std #3 Benzoic Acid 2,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : February 2016 **Storage:** 10°C or colder

C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Benzoic acid	2,000.0 µg/mL	+/- 11.6284	µg/mL	Gravimetric
	CAS # 65-85-0.SEC		+/- 96.5270	µg/mL	Unstressed
	Purity 97%		+/- 96.6098	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Column:
30m x .25mm x .25um
Rtx-5 (cat.#10223)

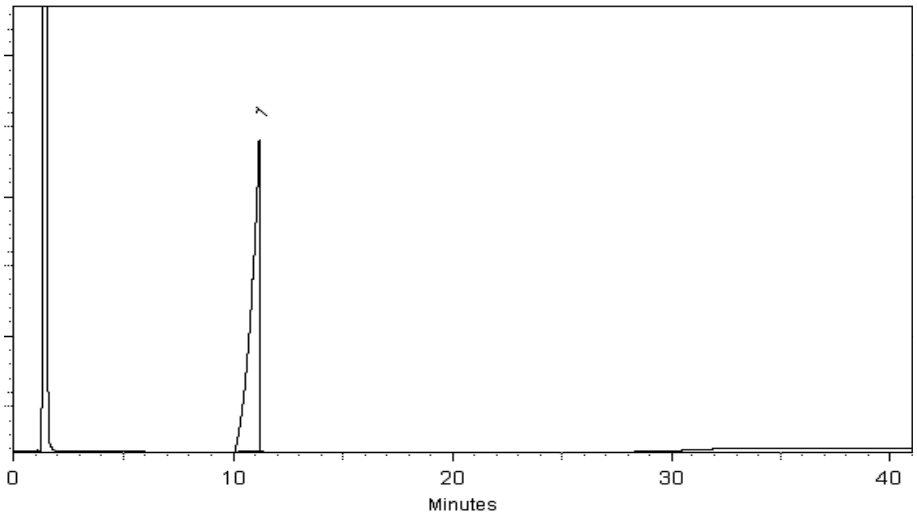
Carrier Gas:
hydrogen-constant pressure 10 psi.

Temp. Program:
40°C (hold 2 min.) to 330°C
@ 10°C/min. (hold 10 min.)

Inj. Temp:
250°C

Det. Temp:
330°C

Det. Type:
FID



Jodi E. Breon - QA Analyst

Date Passed: 22-Feb-2013 **Balance:** 1128353505

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date of the unopened ampul stored at the recommended storage condition is the last day of the month listed in the expiration date field.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31840, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

RESTEK

Chemical Standard Batch Sheet
Lot #: A093654

Catalog #: 567674.sec	**Target:** 2000 ug/mL	
Description: 8270 List 1 / Std #3 Benzoic Acid		
Solvent: Methylene Chloride	**Solvent Lot:** 126244	**Final Volume:** 1,000 ml

Made by: Mary Ellen Wood	**Date:** 2/19/2013 3:12:24PM		
Tested by: Jennifer Pollino	**Date:** 2/21/2013 11:48:51AM		
Pass	**By:** Jodi Breon	**Date:** 2/22/2013 12:34:09PM	
Packaged by: Alexandria Pavkovich / Alexandria Pavkov	**Date:** 2/21/2013 8:30:06AM	**No. Units:** 161	**Pkg Size:** 5 mL
Balance Used: BEDEARMBALPC2 XP205	**Serial #:** 1128353505		

Compound	CAS	Storage Location	Lot #	Purity	Target Conc(ug/mL)	Target	Actual	Calc Conc(ug/mL)
Benzoic acid	65-85-0.SEC	RS063	QD3UO	0.97	2,000.00	2,061.86 mg	2,061.90 mg	2,000.0

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This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 567674 **Lot No.:** A093441

Description : 8270 List 1 / Std #3 Benzoic Acid

8270 List 1 / Std #3 Benzoic Acid 2,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : February 2016 **Storage:** 10°C or colder

C E R T I F I E D V A L U E S

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Benzoic acid	2,000.0 µg/mL	+/- 11.6282	µg/mL	Gravimetric
	CAS # 65-85-0		+/- 96.5249	µg/mL	Unstressed
	Purity 99%		+/- 96.6077	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Column:
30m x .25mm x .25um
Rtx-5 (cat.#10223)

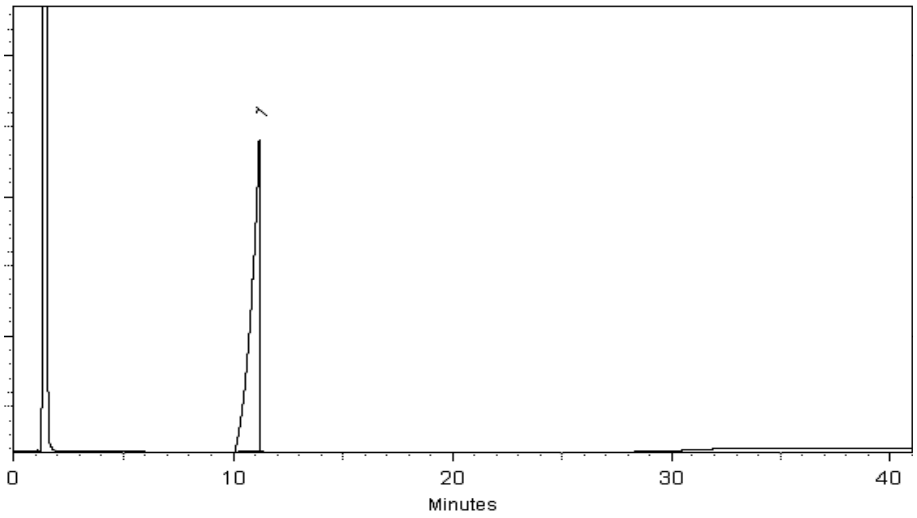
Carrier Gas:
hydrogen-constant pressure 10 psi.

Temp. Program:
40°C (hold 2 min.) to 330°C
@ 10°C/min. (hold 10 min.)

Inj. Temp:
250°C

Det. Temp:
330°C

Det. Type:
FID



Jodi E. Breon - QA Analyst

Date Passed: 22-Feb-2013 **Balance:** 1128342313

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date of the unopened ampul stored at the recommended storage condition is the last day of the month listed in the expiration date field.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31840, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

RESTEK

Chemical Standard Batch Sheet
Lot #: A093441

Catalog #: 567674	**Target:** 2000 ug/mL	
Description: 8270 List 1 / Std #3 Benzoic Acid		
Solvent: Methylene Chloride	**Solvent Lot:** 127438	**Final Volume:** 4,000 ml

Made by: Matt Hepfer	**Date:** 2/7/2013 1:17:18PM		
Tested by: Jodi Breon	**Date:** 2/8/2013 3:42:28PM		
Pass	**By:** Jodi Breon	**Date:** 2/22/2013 12:33:55PM	
Packaged by: Kendra Swope / Alexandria Pavkovich	**Date:** 2/8/2013 10:02:22AM	**No. Units:** 615	**Pkg Size:** 5 mL
Balance Used: BEDEARMBALPC1 XP205	**Serial #:** 1128342313		

Compound	CAS	Storage Location	Lot #	Purity	Target Conc(ug/mL)	Target	Actual	Calc Conc(ug/mL)
Benzoic acid	65-85-0	R0472	MKBG9391V	0.99	2,000.00	8,000.00 mg	8,000.00 mg	2,000.0

QA Report: 8270 List 1/ Std. #3 Benzoic Acid (Cat.#567674)

	Runs of Lot # A093441						Runs of Lot # A093654							
COMPONENT	Run #1	Run #2	Run #3	AVG	STD DEV	% RSD	Run #1	Run #2	Run #3	AVG	STD DEV	% RSD	%D MEAN	**P/F**
Benzoic acid	5123644	5100875	5144322	5122947	21732	0.42	5288104	5295526	5370045	5317892	45318	0.85	-3.81	PASS

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Gravimetric Certificate

FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 568023 **Lot No.:** A0101191

Description : 8270 Famphur Standard

8270 Famphur Standard 2,000 µg/ml, Methylene Chloride, 1 ml/ampul

Container Size : 2 mL **Pkg Amt:** > 1 mL

Expiration Date : February 29, 2016 **Storage:** 10°C or colder

Handling: This product is photosensitive.

CERTIFIED VALUES

Component #	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Famphur	2,011.6 µg/mL	+/- 20.2643	µg/mL	Gravimetric
	CAS # 52-85-7 (Lot 1269600)		+/- 74.2962	µg/mL	Unstressed
	Purity 94%		+/- 74.2987	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Michael J Maye

Date Mixed: 06-Feb-2014 **Balance:** 1128353505

Manufactured under Restek's ISO 9001:2008
Registered Quality System
Certificate #FM 80397

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FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 568038 **Lot No.:** A0101189

Description : Diphenylamine Standard

Diphenylamine Standard 1708 µg/ml, Methylene Chloride, 1 ml/ampul

Container Size : 2 mL **Pkg Amt:** > 1 mL

Expiration Date : February 28, 2017 **Storage:** 10°C or colder

CERTIFIED VALUES

Component #	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Diphenylamine	1,712.0 µg/mL	+/- 17.2462	µg/mL	Gravimetric
	CAS # 122-39-4 (Lot 07525MF)		+/- 23.4457	µg/mL	Unstressed
	Purity 99%		+/- 34.7730	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Michael J Maye

Date Mixed: 06-Feb-2014 **Balance:** 1128353505

Manufactured under Restek's ISO 9001:2008
Registered Quality System
Certificate #FM 80397

Tech Tips:

N-Nitrosodiphenylamine is prone to breakdown in the injection port and will be converted to diphenylamine.
N-Nitrosodiphenylamine is also a reactive species that can initiate premature decomposition of other compounds in the mix. For these reasons diphenylamine is used in the preparation of this mixture. When comparing the response of this compound to mixtures manufactured using N-nitrosodiphenylamine, a difference in response will be observed.

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CERTIFIED REFERENCE MATERIAL

Certificate of Analysis

ACCREDITED
ISO Guide 34 Accredited
Reference Material Producer
Certificate #3222.01

ACCREDITED
ISO/IEC 17025 Accredited
Testing Laboratory
Certificate #3222.02

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Catalog No. : 569729 **Lot No.:** A0107399

Description : 8270 List 1 / Std #1 MegaMix (2015)

8270 List 1 / Std #1 MegaMix (2015) 500-2000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 10 mL **Pkg Amt:** > 5 mL

Expiration Date : May 31, 2016 **Storage:** 10°C or colder

Handling: Carcinogen/reproductive toxin. Photosensitive. Sonicate.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)	
1	1,4-Dioxane **CAS #** 123-91-1 (Lot SHBD8744V) **Purity** 99%	1,004.4 µg/mL	+/- 5.8397 µg/mL +/- 10.9969 µg/mL +/- 18.6525 µg/mL	Gravimetric Unstressed Stressed
2	Pyridine **CAS #** 110-86-1 (Lot SHBC7174V) **Purity** 99%	1,001.0 µg/mL	+/- 5.8199 µg/mL +/- 10.9596 µg/mL +/- 18.5894 µg/mL	Gravimetric Unstressed Stressed
3	N-Nitrosodimethylamine **CAS #** 62-75-9 (Lot 3213100) **Purity** 99%	1,000.2 µg/mL	+/- 5.8152 µg/mL +/- 10.9509 µg/mL +/- 18.5745 µg/mL	Gravimetric Unstressed Stressed
4	Aniline **CAS #** 62-53-3 (Lot K22Z462) **Purity** 99%	1,002.3 µg/mL	+/- 5.8275 µg/mL +/- 10.9739 µg/mL +/- 18.6135 µg/mL	Gravimetric Unstressed Stressed
5	Bis(2-chloroethyl)ether **CAS #** 111-44-4 (Lot 45296HKV) **Purity** 99%	1,001.4 µg/mL	+/- 5.8222 µg/mL +/- 10.9640 µg/mL +/- 18.5968 µg/mL	Gravimetric Unstressed Stressed
6	2-Chlorophenol **CAS #** 95-57-8 (Lot MKBD3900V) **Purity** 99%	1,000.8 µg/mL	+/- 5.8187 µg/mL +/- 10.9575 µg/mL +/- 18.5856 µg/mL	Gravimetric Unstressed Stressed
7	Phenol **CAS #** 108-95-2 (Lot SHBC6998V) **Purity** 99%	1,006.9 µg/mL	+/- 5.8542 µg/mL +/- 11.0242 µg/mL +/- 18.6989 µg/mL	Gravimetric Unstressed Stressed

8	n-Decane (C10)		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 124-18-5	(Lot SHBF1587V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
9	1,4-Dichlorobenzene		1,005.3	µg/mL	+/-	5.8449	µg/mL	Gravimetric
	CAS # 106-46-7	(Lot MKBL3891V)			+/-	11.0067	µg/mL	Unstressed
	Purity 99%				+/-	18.6692	µg/mL	Stressed
10	1,3-Dichlorobenzene		1,002.0	µg/mL	+/-	5.8257	µg/mL	Gravimetric
	CAS # 541-73-1	(Lot BCBC1891V)			+/-	10.9706	µg/mL	Unstressed
	Purity 99%				+/-	18.6079	µg/mL	Stressed
11	1,2-Dichlorobenzene		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 95-50-1	(Lot 68996CMV)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
12	Benzyl alcohol		1,000.4	µg/mL	+/-	5.8164	µg/mL	Gravimetric
	CAS # 100-51-6	(Lot SHBC1850V)			+/-	10.9531	µg/mL	Unstressed
	Purity 99%				+/-	18.5782	µg/mL	Stressed
13	2,2'-oxybis(1-chloropropane)		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 108-60-1	(Lot 2-EAW-18-3)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
14	2-Methylphenol (o-cresol)		1,003.6	µg/mL	+/-	5.8350	µg/mL	Gravimetric
	CAS # 95-48-7	(Lot SHBC1479V)			+/-	10.9881	µg/mL	Unstressed
	Purity 99%				+/-	18.6376	µg/mL	Stressed
15	Hexachloroethane		1,005.9	µg/mL	+/-	5.8484	µg/mL	Gravimetric
	CAS # 67-72-1	(Lot 4H3SF)			+/-	11.0133	µg/mL	Unstressed
	Purity 99%				+/-	18.6804	µg/mL	Stressed
16	Acetophenone		1,002.7	µg/mL	+/-	5.8298	µg/mL	Gravimetric
	CAS # 98-86-2	(Lot MKBR7156V)			+/-	10.9783	µg/mL	Unstressed
	Purity 99%				+/-	18.6209	µg/mL	Stressed
17	N-Nitroso-di-n-propylamine		1,003.9	µg/mL	+/-	5.8368	µg/mL	Gravimetric
	CAS # 621-64-7	(Lot OPAGF)			+/-	10.9914	µg/mL	Unstressed
	Purity 99%				+/-	18.6432	µg/mL	Stressed
18	4-Methylphenol (p-cresol)		500.4	µg/mL	+/-	2.9161	µg/mL	Gravimetric
	CAS # 106-44-5	(Lot 49396APV)			+/-	5.4823	µg/mL	Unstressed
	Purity 99%				+/-	9.2949	µg/mL	Stressed
19	3-Methylphenol (m-cresol)		500.2	µg/mL	+/-	2.9149	µg/mL	Gravimetric
	CAS # 108-39-4	(Lot SHBD0627V)			+/-	5.4801	µg/mL	Unstressed
	Purity 99%				+/-	9.2912	µg/mL	Stressed
20	Nitrobenzene		1,001.1	µg/mL	+/-	5.8205	µg/mL	Gravimetric
	CAS # 98-95-3	(Lot 65096APV)			+/-	10.9607	µg/mL	Unstressed
	Purity 99%				+/-	18.5912	µg/mL	Stressed
21	Isophorone		999.3	µg/mL	+/-	5.8100	µg/mL	Gravimetric
	CAS # 78-59-1	(Lot 06705DE)			+/-	10.9410	µg/mL	Unstressed
	Purity 97%				+/-	18.5577	µg/mL	Stressed
22	2-Nitrophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 88-75-5	(Lot BCBH7602V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
23	2,4-Dimethylphenol		1,003.2	µg/mL	+/-	5.8327	µg/mL	Gravimetric
	CAS # 105-67-9	(Lot 10165155)			+/-	10.9837	µg/mL	Unstressed
	Purity 99%				+/-	18.6302	µg/mL	Stressed

24	Bis(2-chloroethoxy)methane		1,004.5	µg/mL	+/-	5.8402	µg/mL	Gravimetric
	CAS # 111-91-1	(Lot 317200)			+/-	10.9980	µg/mL	Unstressed
	Purity 99%				+/-	18.6544	µg/mL	Stressed
25	2,4-Dichlorophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-83-2	(Lot BCBH1617V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
26	1,2,4-Trichlorobenzene		1,000.6	µg/mL	+/-	5.8176	µg/mL	Gravimetric
	CAS # 120-82-1	(Lot 26896BM)			+/-	10.9553	µg/mL	Unstressed
	Purity 99%				+/-	18.5819	µg/mL	Stressed
27	Naphthalene		1,002.3	µg/mL	+/-	5.8275	µg/mL	Gravimetric
	CAS # 91-20-3	(Lot MKBH4351V)			+/-	10.9739	µg/mL	Unstressed
	Purity 99%				+/-	18.6135	µg/mL	Stressed
28	2,6-Dichlorophenol		1,000.1	µg/mL	+/-	5.8147	µg/mL	Gravimetric
	CAS # 87-65-0	(Lot MKBN2776V)			+/-	10.9498	µg/mL	Unstressed
	Purity 99%				+/-	18.5726	µg/mL	Stressed
29	4-Chloroaniline		1,000.3	µg/mL	+/-	5.8157	µg/mL	Gravimetric
	CAS # 106-47-8	(Lot 12528PH)			+/-	10.9518	µg/mL	Unstressed
	Purity 98%				+/-	18.5761	µg/mL	Stressed
30	Hexachlorobutadiene		999.9	µg/mL	+/-	5.8135	µg/mL	Gravimetric
	CAS # 87-68-3	(Lot K22W009)			+/-	10.9475	µg/mL	Unstressed
	Purity 98%				+/-	18.5688	µg/mL	Stressed
31	2-Methylnaphthalene		998.6	µg/mL	+/-	5.8059	µg/mL	Gravimetric
	CAS # 91-57-6	(Lot 19399MJV)			+/-	10.9333	µg/mL	Unstressed
	Purity 96%				+/-	18.5446	µg/mL	Stressed
32	4-Chloro-3-methylphenol		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 59-50-7	(Lot STBC0769V)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
33	1-Methylnaphthalene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 90-12-0	(Lot 525000-10)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
34	1,2,4,5-Tetrachlorobenzene		1,000.2	µg/mL	+/-	5.8152	µg/mL	Gravimetric
	CAS # 95-94-3	(Lot 06024AIV)			+/-	10.9509	µg/mL	Unstressed
	Purity 99%				+/-	18.5745	µg/mL	Stressed
35	Hexachlorocyclopentadiene		1,002.4	µg/mL	+/-	5.8280	µg/mL	Gravimetric
	CAS # 77-47-4	(Lot 3140300)			+/-	10.9750	µg/mL	Unstressed
	Purity 99%				+/-	18.6154	µg/mL	Stressed
36	2,4,6-Trichlorophenol		1,001.4	µg/mL	+/-	5.8222	µg/mL	Gravimetric
	CAS # 88-06-2	(Lot MKBH7393V)			+/-	10.9640	µg/mL	Unstressed
	Purity 99%				+/-	18.5968	µg/mL	Stressed
37	2,4,5-Trichlorophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-95-4	(Lot FHM01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
38	2-Chloronaphthalene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-58-7	(Lot FIJ01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
39	Biphenyl		1,006.1	µg/mL	+/-	5.8496	µg/mL	Gravimetric
	CAS # 92-52-4	(Lot 1277976)			+/-	11.0155	µg/mL	Unstressed
	Purity 99%				+/-	18.6841	µg/mL	Stressed

40	2-Nitroaniline		1,004.6	µg/mL	+/-	5.8408	µg/mL	Gravimetric
	CAS # 88-74-4	(Lot MKBK7597V)			+/-	10.9991	µg/mL	Unstressed
	Purity 99%				+/-	18.6562	µg/mL	Stressed
41	Acenaphthylene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 208-96-8	(Lot ER030707-01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
42	1,3-Dinitrobenzene		1,006.4	µg/mL	+/-	5.8513	µg/mL	Gravimetric
	CAS # 99-65-0	(Lot BCBB1436V)			+/-	11.0188	µg/mL	Unstressed
	Purity 99%				+/-	18.6896	µg/mL	Stressed
43	Dimethylphthalate		1,001.0	µg/mL	+/-	5.8199	µg/mL	Gravimetric
	CAS # 131-11-3	(Lot 10117699)			+/-	10.9596	µg/mL	Unstressed
	Purity 99%				+/-	18.5894	µg/mL	Stressed
44	2,6-Dinitrotoluene		1,001.6	µg/mL	+/-	5.8234	µg/mL	Gravimetric
	CAS # 606-20-2	(Lot 1437483V)			+/-	10.9662	µg/mL	Unstressed
	Purity 99%				+/-	18.6005	µg/mL	Stressed
45	Acenaphthene		1,001.8	µg/mL	+/-	5.8246	µg/mL	Gravimetric
	CAS # 83-32-9	(Lot MKBP0384V)			+/-	10.9684	µg/mL	Unstressed
	Purity 99%				+/-	18.6042	µg/mL	Stressed
46	2,4-Dinitrophenol		2,008.7	µg/mL	+/-	11.6788	µg/mL	Gravimetric
	CAS # 51-28-5	(Lot MKBP5833V)			+/-	21.9927	µg/mL	Unstressed
	Purity 99%				+/-	37.3031	µg/mL	Stressed
47	Dibenzofuran		1,001.8	µg/mL	+/-	5.8246	µg/mL	Gravimetric
	CAS # 132-64-9	(Lot MKBK2375V)			+/-	10.9684	µg/mL	Unstressed
	Purity 99%				+/-	18.6042	µg/mL	Stressed
48	3-Nitroaniline		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 99-09-2	(Lot MKBH5131V)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
49	2,4-Dinitrotoluene		1,001.7	µg/mL	+/-	5.8240	µg/mL	Gravimetric
	CAS # 121-14-2	(Lot MKAA0690V)			+/-	10.9673	µg/mL	Unstressed
	Purity 99%				+/-	18.6024	µg/mL	Stressed
50	4-Nitrophenol		2,001.0	µg/mL	+/-	11.6340	µg/mL	Gravimetric
	CAS # 100-02-7	(Lot MKBK1842V)			+/-	21.9083	µg/mL	Unstressed
	Purity 99%				+/-	37.1601	µg/mL	Stressed
51	2,3,4,6-Tetrachlorophenol		1,002.1	µg/mL	+/-	5.8263	µg/mL	Gravimetric
	CAS # 58-90-2	(Lot FN10221307)			+/-	10.9717	µg/mL	Unstressed
	Purity 99%				+/-	18.6098	µg/mL	Stressed
52	Fluorene		1,000.5	µg/mL	+/-	5.8169	µg/mL	Gravimetric
	CAS # 86-73-7	(Lot 10174662)			+/-	10.9540	µg/mL	Unstressed
	Purity 98%				+/-	18.5797	µg/mL	Stressed
53	4-Chlorophenyl phenyl ether		1,000.4	µg/mL	+/-	5.8164	µg/mL	Gravimetric
	CAS # 7005-72-3	(Lot MKBL1347V)			+/-	10.9531	µg/mL	Unstressed
	Purity 99%				+/-	18.5782	µg/mL	Stressed
54	n-Hexadecane (C16)		1,002.2	µg/mL	+/-	5.8269	µg/mL	Gravimetric
	CAS # 544-76-3	(Lot SHBD4570V)			+/-	10.9728	µg/mL	Unstressed
	Purity 99%				+/-	18.6116	µg/mL	Stressed
55	Diethylphthalate		1,001.1	µg/mL	+/-	5.8205	µg/mL	Gravimetric
	CAS # 84-66-2	(Lot MKBJ3578V)			+/-	10.9607	µg/mL	Unstressed
	Purity 99%				+/-	18.5912	µg/mL	Stressed

56	Azobenzene		1,002.3	µg/mL	+/-	5.8275	µg/mL	Gravimetric
	CAS # 103-33-3	(Lot MKBS2559V)			+/-	10.9739	µg/mL	Unstressed
	Purity 99%				+/-	18.6135	µg/mL	Stressed
57	Diphenylamine		1,713.4	µg/mL	+/-	9.9619	µg/mL	Gravimetric
	CAS # 122-39-4	(Lot 07525MF)			+/-	18.7595	µg/mL	Unstressed
	Purity 99%				+/-	31.8192	µg/mL	Stressed
58	4-Nitroaniline		1,002.8	µg/mL	+/-	5.8304	µg/mL	Gravimetric
	CAS # 100-01-6	(Lot BCBG4702V)			+/-	10.9794	µg/mL	Unstressed
	Purity 99%				+/-	18.6228	µg/mL	Stressed
59	4,6-Dinitro-2-methylphenol (Dinitro-o-cresol)		2,002.0	µg/mL	+/-	11.6398	µg/mL	Gravimetric
	CAS # 534-52-1	(Lot LC06195V)			+/-	21.9193	µg/mL	Unstressed
	Purity 99%				+/-	37.1787	µg/mL	Stressed
60	4-Bromophenyl phenyl ether		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 101-55-3	(Lot STBB9729V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
61	Hexachlorobenzene		1,002.1	µg/mL	+/-	5.8260	µg/mL	Gravimetric
	CAS # 118-74-1	(Lot LC04221V)			+/-	10.9711	µg/mL	Unstressed
	Purity 98%				+/-	18.6089	µg/mL	Stressed
62	Pentachlorophenol		2,000.3	µg/mL	+/-	11.6299	µg/mL	Gravimetric
	CAS # 87-86-5	(Lot 140626JLM)			+/-	21.9007	µg/mL	Unstressed
	Purity 99%				+/-	37.1471	µg/mL	Stressed
63	Phenanthrene		999.0	µg/mL	+/-	5.8083	µg/mL	Gravimetric
	CAS # 85-01-8	(Lot MKBL6906V)			+/-	10.9379	µg/mL	Unstressed
	Purity 98%				+/-	18.5524	µg/mL	Stressed
64	n-Octadecane (C18)		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 593-45-3	(Lot OGCDK)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
65	Anthracene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-12-7	(Lot MKBK5208V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
66	Carbazole		1,000.1	µg/mL	+/-	5.8146	µg/mL	Gravimetric
	CAS # 86-74-8	(Lot S42950-417)			+/-	10.9497	µg/mL	Unstressed
	Purity 98%				+/-	18.5725	µg/mL	Stressed
67	Di-n-butylphthalate		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 84-74-2	(Lot MKBL8501V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
68	Fluoranthene		999.7	µg/mL	+/-	5.8123	µg/mL	Gravimetric
	CAS # 206-44-0	(Lot MKBQ6360V)			+/-	10.9454	µg/mL	Unstressed
	Purity 98%				+/-	18.5652	µg/mL	Stressed
69	Pyrene		999.1	µg/mL	+/-	5.8089	µg/mL	Gravimetric
	CAS # 129-00-0	(Lot BCBJ0984V)			+/-	10.9390	µg/mL	Unstressed
	Purity 98%				+/-	18.5543	µg/mL	Stressed
70	Benzyl butyl phthalate		1,001.2	µg/mL	+/-	5.8211	µg/mL	Gravimetric
	CAS # 85-68-7	(Lot 03027HV)			+/-	10.9618	µg/mL	Unstressed
	Purity 99%				+/-	18.5931	µg/mL	Stressed
71	Benz(a)anthracene		1,001.4	µg/mL	+/-	5.8222	µg/mL	Gravimetric
	CAS # 56-55-3	(Lot ER031412-01)			+/-	10.9640	µg/mL	Unstressed
	Purity 99%				+/-	18.5968	µg/mL	Stressed

72	Chrysene		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 218-01-9	(Lot ER120810-02)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
73	Bis(2-ethylhexyl)phthalate		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 117-81-7	(Lot MKBK2695V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
74	Di-n-octyl phthalate		1,002.1	µg/mL	+/-	5.8263	µg/mL	Gravimetric
	CAS # 117-84-0	(Lot 3128600)			+/-	10.9717	µg/mL	Unstressed
	Purity 99%				+/-	18.6098	µg/mL	Stressed
75	Benzo(b)fluoranthene		1,000.1	µg/mL	+/-	5.8147	µg/mL	Gravimetric
	CAS # 205-99-2	(Lot ER03101401)			+/-	10.9498	µg/mL	Unstressed
	Purity 99%				+/-	18.5726	µg/mL	Stressed
76	Benzo(k)fluoranthene		1,003.3	µg/mL	+/-	5.8333	µg/mL	Gravimetric
	CAS # 207-08-9	(Lot ER041513-01)			+/-	10.9848	µg/mL	Unstressed
	Purity 99%				+/-	18.6321	µg/mL	Stressed
77	Benzo(a)pyrene		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 50-32-8	(Lot ER071309-02)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
78	Indeno(1,2,3-cd)pyrene		1,000.6	µg/mL	+/-	5.8176	µg/mL	Gravimetric
	CAS # 193-39-5	(Lot ER082107-02)			+/-	10.9553	µg/mL	Unstressed
	Purity 99%				+/-	18.5819	µg/mL	Stressed
79	Dibenz(a,h)anthracene		1,004.4	µg/mL	+/-	5.8397	µg/mL	Gravimetric
	CAS # 53-70-3	(Lot ER032211-01)			+/-	10.9969	µg/mL	Unstressed
	Purity 99%				+/-	18.6525	µg/mL	Stressed
80	Benzo(g,h,i)perylene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 191-24-2	(Lot ER020708-08)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Specific Reference Material Notes:

N-nitrosodiphenylamine 2000 ug/mL equivalent when used for GC analysis. Actual formulation is diphenylamine 1710 ug/mL. N-Nitrosodiphenylamine is prone to breakdown in the injection port and will be converted to diphenylamine. N-Nitrosodiphenylamine is also a reactive species that can initiate premature decomposition of other compounds in the mix. For these reasons diphenylamine is used in the preparation of this mixture. When comparing the response of this compound to mixtures manufactured using N-nitrosodiphenylamine, a difference in response will be observed.

Column:
30m x 0.25mm x 0.25µm
Rtx-5 (cat.#10223)

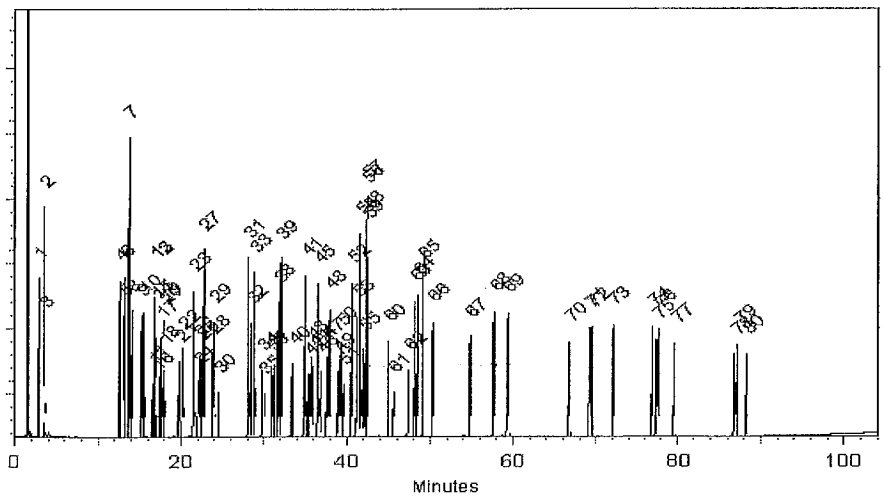
Carrier Gas:
hydrogen-constant pressure 10 psi

Temp. Program:
35°C (hold 3 min.) to 330°C
@ 3°C/min. (hold 3 min.)

Inj. Temp:
250°C

Det. Temp:
300°C

Det. Type:
FID



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

F. Joseph Tallon - Mix Technician **Date Mixed:** 24-Nov-2014 **Balance:** 1128360905

Jodi E. Breon - QA Analyst **Date Passed:** 05-Dec-2014

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

k is a coverage factor of 2, which gives a level of confidence of approximately 95%.

- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

RESTEK®
110 Benner Circle
Bellefonte, PA 16823-8812
Tel: (800)356-1688
Fax: (814)353-1309

www.restek.com

CERTIFIED REFERENCE MATERIAL

Certificate of Analysis

ACCREDITED
ISO Guide 34 Accredited
Reference Material Producer
Certificate #3222.01

ilac-MRA

ACCREDITED
ISO/IEC 17025 Accredited
Testing Laboratory
Certificate #3222.02

FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 569729 **Lot No.:** A0107399

Description : 8270 List 1 / Std #1 MegaMix (2015)

8270 List 1 / Std #1 MegaMix (2015) 500-2000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 10 mL **Pkg Amt:** > 5 mL

Expiration Date : May 31, 2016 **Storage:** 10°C or colder

Handling: Carcinogen/reproductive toxin. Photosensitive. Sonicate.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)	
1	1,4-Dioxane **CAS #** 123-91-1 (Lot SHBD8744V) **Purity** 99%	1,004.4 µg/mL	+/- 5.8397 µg/mL +/- 10.9969 µg/mL +/- 18.6525 µg/mL	Gravimetric Unstressed Stressed
2	Pyridine **CAS #** 110-86-1 (Lot SHBC7174V) **Purity** 99%	1,001.0 µg/mL	+/- 5.8199 µg/mL +/- 10.9596 µg/mL +/- 18.5894 µg/mL	Gravimetric Unstressed Stressed
3	N-Nitrosodimethylamine **CAS #** 62-75-9 (Lot 3213100) **Purity** 99%	1,000.2 µg/mL	+/- 5.8152 µg/mL +/- 10.9509 µg/mL +/- 18.5745 µg/mL	Gravimetric Unstressed Stressed
4	Aniline **CAS #** 62-53-3 (Lot K22Z462) **Purity** 99%	1,002.3 µg/mL	+/- 5.8275 µg/mL +/- 10.9739 µg/mL +/- 18.6135 µg/mL	Gravimetric Unstressed Stressed
5	Bis(2-chloroethyl)ether **CAS #** 111-44-4 (Lot 45296HKV) **Purity** 99%	1,001.4 µg/mL	+/- 5.8222 µg/mL +/- 10.9640 µg/mL +/- 18.5968 µg/mL	Gravimetric Unstressed Stressed
6	2-Chlorophenol **CAS #** 95-57-8 (Lot MKBD3900V) **Purity** 99%	1,000.8 µg/mL	+/- 5.8187 µg/mL +/- 10.9575 µg/mL +/- 18.5856 µg/mL	Gravimetric Unstressed Stressed
7	Phenol **CAS #** 108-95-2 (Lot SHBC6998V) **Purity** 99%	1,006.9 µg/mL	+/- 5.8542 µg/mL +/- 11.0242 µg/mL +/- 18.6989 µg/mL	Gravimetric Unstressed Stressed

8	n-Decane (C10)		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 124-18-5	(Lot SHBF1587V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
9	1,4-Dichlorobenzene		1,005.3	µg/mL	+/-	5.8449	µg/mL	Gravimetric
	CAS # 106-46-7	(Lot MKBL3891V)			+/-	11.0067	µg/mL	Unstressed
	Purity 99%				+/-	18.6692	µg/mL	Stressed
10	1,3-Dichlorobenzene		1,002.0	µg/mL	+/-	5.8257	µg/mL	Gravimetric
	CAS # 541-73-1	(Lot BCBC1891V)			+/-	10.9706	µg/mL	Unstressed
	Purity 99%				+/-	18.6079	µg/mL	Stressed
11	1,2-Dichlorobenzene		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 95-50-1	(Lot 68996CMV)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
12	Benzyl alcohol		1,000.4	µg/mL	+/-	5.8164	µg/mL	Gravimetric
	CAS # 100-51-6	(Lot SHBC1850V)			+/-	10.9531	µg/mL	Unstressed
	Purity 99%				+/-	18.5782	µg/mL	Stressed
13	2,2'-oxybis(1-chloropropane)		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 108-60-1	(Lot 2-EAW-18-3)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
14	2-Methylphenol (o-cresol)		1,003.6	µg/mL	+/-	5.8350	µg/mL	Gravimetric
	CAS # 95-48-7	(Lot SHBC1479V)			+/-	10.9881	µg/mL	Unstressed
	Purity 99%				+/-	18.6376	µg/mL	Stressed
15	Hexachloroethane		1,005.9	µg/mL	+/-	5.8484	µg/mL	Gravimetric
	CAS # 67-72-1	(Lot 4H3SF)			+/-	11.0133	µg/mL	Unstressed
	Purity 99%				+/-	18.6804	µg/mL	Stressed
16	Acetophenone		1,002.7	µg/mL	+/-	5.8298	µg/mL	Gravimetric
	CAS # 98-86-2	(Lot MKBR7156V)			+/-	10.9783	µg/mL	Unstressed
	Purity 99%				+/-	18.6209	µg/mL	Stressed
17	N-Nitroso-di-n-propylamine		1,003.9	µg/mL	+/-	5.8368	µg/mL	Gravimetric
	CAS # 621-64-7	(Lot OPAGF)			+/-	10.9914	µg/mL	Unstressed
	Purity 99%				+/-	18.6432	µg/mL	Stressed
18	4-Methylphenol (p-cresol)		500.4	µg/mL	+/-	2.9161	µg/mL	Gravimetric
	CAS # 106-44-5	(Lot 49396APV)			+/-	5.4823	µg/mL	Unstressed
	Purity 99%				+/-	9.2949	µg/mL	Stressed
19	3-Methylphenol (m-cresol)		500.2	µg/mL	+/-	2.9149	µg/mL	Gravimetric
	CAS # 108-39-4	(Lot SHBD0627V)			+/-	5.4801	µg/mL	Unstressed
	Purity 99%				+/-	9.2912	µg/mL	Stressed
20	Nitrobenzene		1,001.1	µg/mL	+/-	5.8205	µg/mL	Gravimetric
	CAS # 98-95-3	(Lot 65096APV)			+/-	10.9607	µg/mL	Unstressed
	Purity 99%				+/-	18.5912	µg/mL	Stressed
21	Isophorone		999.3	µg/mL	+/-	5.8100	µg/mL	Gravimetric
	CAS # 78-59-1	(Lot 06705DE)			+/-	10.9410	µg/mL	Unstressed
	Purity 97%				+/-	18.5577	µg/mL	Stressed
22	2-Nitrophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 88-75-5	(Lot BCBH7602V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
23	2,4-Dimethylphenol		1,003.2	µg/mL	+/-	5.8327	µg/mL	Gravimetric
	CAS # 105-67-9	(Lot 10165155)			+/-	10.9837	µg/mL	Unstressed
	Purity 99%				+/-	18.6302	µg/mL	Stressed

24	Bis(2-chloroethoxy)methane		1,004.5	µg/mL	+/-	5.8402	µg/mL	Gravimetric
	CAS # 111-91-1	(Lot 317200)			+/-	10.9980	µg/mL	Unstressed
	Purity 99%				+/-	18.6544	µg/mL	Stressed
25	2,4-Dichlorophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-83-2	(Lot BCBH1617V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
26	1,2,4-Trichlorobenzene		1,000.6	µg/mL	+/-	5.8176	µg/mL	Gravimetric
	CAS # 120-82-1	(Lot 26896BM)			+/-	10.9553	µg/mL	Unstressed
	Purity 99%				+/-	18.5819	µg/mL	Stressed
27	Naphthalene		1,002.3	µg/mL	+/-	5.8275	µg/mL	Gravimetric
	CAS # 91-20-3	(Lot MKBH4351V)			+/-	10.9739	µg/mL	Unstressed
	Purity 99%				+/-	18.6135	µg/mL	Stressed
28	2,6-Dichlorophenol		1,000.1	µg/mL	+/-	5.8147	µg/mL	Gravimetric
	CAS # 87-65-0	(Lot MKBN2776V)			+/-	10.9498	µg/mL	Unstressed
	Purity 99%				+/-	18.5726	µg/mL	Stressed
29	4-Chloroaniline		1,000.3	µg/mL	+/-	5.8157	µg/mL	Gravimetric
	CAS # 106-47-8	(Lot 12528PH)			+/-	10.9518	µg/mL	Unstressed
	Purity 98%				+/-	18.5761	µg/mL	Stressed
30	Hexachlorobutadiene		999.9	µg/mL	+/-	5.8135	µg/mL	Gravimetric
	CAS # 87-68-3	(Lot K22W009)			+/-	10.9475	µg/mL	Unstressed
	Purity 98%				+/-	18.5688	µg/mL	Stressed
31	2-Methylnaphthalene		998.6	µg/mL	+/-	5.8059	µg/mL	Gravimetric
	CAS # 91-57-6	(Lot 19399MJV)			+/-	10.9333	µg/mL	Unstressed
	Purity 96%				+/-	18.5446	µg/mL	Stressed
32	4-Chloro-3-methylphenol		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 59-50-7	(Lot STBC0769V)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
33	1-Methylnaphthalene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 90-12-0	(Lot 525000-10)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
34	1,2,4,5-Tetrachlorobenzene		1,000.2	µg/mL	+/-	5.8152	µg/mL	Gravimetric
	CAS # 95-94-3	(Lot 06024AIV)			+/-	10.9509	µg/mL	Unstressed
	Purity 99%				+/-	18.5745	µg/mL	Stressed
35	Hexachlorocyclopentadiene		1,002.4	µg/mL	+/-	5.8280	µg/mL	Gravimetric
	CAS # 77-47-4	(Lot 3140300)			+/-	10.9750	µg/mL	Unstressed
	Purity 99%				+/-	18.6154	µg/mL	Stressed
36	2,4,6-Trichlorophenol		1,001.4	µg/mL	+/-	5.8222	µg/mL	Gravimetric
	CAS # 88-06-2	(Lot MKBH7393V)			+/-	10.9640	µg/mL	Unstressed
	Purity 99%				+/-	18.5968	µg/mL	Stressed
37	2,4,5-Trichlorophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-95-4	(Lot FHM01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
38	2-Chloronaphthalene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-58-7	(Lot FIJ01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
39	Biphenyl		1,006.1	µg/mL	+/-	5.8496	µg/mL	Gravimetric
	CAS # 92-52-4	(Lot 1277976)			+/-	11.0155	µg/mL	Unstressed
	Purity 99%				+/-	18.6841	µg/mL	Stressed

40	2-Nitroaniline		1,004.6	µg/mL	+/-	5.8408	µg/mL	Gravimetric
	CAS # 88-74-4	(Lot MKBK7597V)			+/-	10.9991	µg/mL	Unstressed
	Purity 99%				+/-	18.6562	µg/mL	Stressed
41	Acenaphthylene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 208-96-8	(Lot ER030707-01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
42	1,3-Dinitrobenzene		1,006.4	µg/mL	+/-	5.8513	µg/mL	Gravimetric
	CAS # 99-65-0	(Lot BCBB1436V)			+/-	11.0188	µg/mL	Unstressed
	Purity 99%				+/-	18.6896	µg/mL	Stressed
43	Dimethylphthalate		1,001.0	µg/mL	+/-	5.8199	µg/mL	Gravimetric
	CAS # 131-11-3	(Lot 10117699)			+/-	10.9596	µg/mL	Unstressed
	Purity 99%				+/-	18.5894	µg/mL	Stressed
44	2,6-Dinitrotoluene		1,001.6	µg/mL	+/-	5.8234	µg/mL	Gravimetric
	CAS # 606-20-2	(Lot 1437483V)			+/-	10.9662	µg/mL	Unstressed
	Purity 99%				+/-	18.6005	µg/mL	Stressed
45	Acenaphthene		1,001.8	µg/mL	+/-	5.8246	µg/mL	Gravimetric
	CAS # 83-32-9	(Lot MKBP0384V)			+/-	10.9684	µg/mL	Unstressed
	Purity 99%				+/-	18.6042	µg/mL	Stressed
46	2,4-Dinitrophenol		2,008.7	µg/mL	+/-	11.6788	µg/mL	Gravimetric
	CAS # 51-28-5	(Lot MKBP5833V)			+/-	21.9927	µg/mL	Unstressed
	Purity 99%				+/-	37.3031	µg/mL	Stressed
47	Dibenzofuran		1,001.8	µg/mL	+/-	5.8246	µg/mL	Gravimetric
	CAS # 132-64-9	(Lot MKBK2375V)			+/-	10.9684	µg/mL	Unstressed
	Purity 99%				+/-	18.6042	µg/mL	Stressed
48	3-Nitroaniline		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 99-09-2	(Lot MKBH5131V)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
49	2,4-Dinitrotoluene		1,001.7	µg/mL	+/-	5.8240	µg/mL	Gravimetric
	CAS # 121-14-2	(Lot MKAA0690V)			+/-	10.9673	µg/mL	Unstressed
	Purity 99%				+/-	18.6024	µg/mL	Stressed
50	4-Nitrophenol		2,001.0	µg/mL	+/-	11.6340	µg/mL	Gravimetric
	CAS # 100-02-7	(Lot MKBK1842V)			+/-	21.9083	µg/mL	Unstressed
	Purity 99%				+/-	37.1601	µg/mL	Stressed
51	2,3,4,6-Tetrachlorophenol		1,002.1	µg/mL	+/-	5.8263	µg/mL	Gravimetric
	CAS # 58-90-2	(Lot FN10221307)			+/-	10.9717	µg/mL	Unstressed
	Purity 99%				+/-	18.6098	µg/mL	Stressed
52	Fluorene		1,000.5	µg/mL	+/-	5.8169	µg/mL	Gravimetric
	CAS # 86-73-7	(Lot 10174662)			+/-	10.9540	µg/mL	Unstressed
	Purity 98%				+/-	18.5797	µg/mL	Stressed
53	4-Chlorophenyl phenyl ether		1,000.4	µg/mL	+/-	5.8164	µg/mL	Gravimetric
	CAS # 7005-72-3	(Lot MKBL1347V)			+/-	10.9531	µg/mL	Unstressed
	Purity 99%				+/-	18.5782	µg/mL	Stressed
54	n-Hexadecane (C16)		1,002.2	µg/mL	+/-	5.8269	µg/mL	Gravimetric
	CAS # 544-76-3	(Lot SHBD4570V)			+/-	10.9728	µg/mL	Unstressed
	Purity 99%				+/-	18.6116	µg/mL	Stressed
55	Diethylphthalate		1,001.1	µg/mL	+/-	5.8205	µg/mL	Gravimetric
	CAS # 84-66-2	(Lot MKBJ3578V)			+/-	10.9607	µg/mL	Unstressed
	Purity 99%				+/-	18.5912	µg/mL	Stressed

56	Azobenzene		1,002.3	µg/mL	+/-	5.8275	µg/mL	Gravimetric
	CAS # 103-33-3	(Lot MKBS2559V)			+/-	10.9739	µg/mL	Unstressed
	Purity 99%				+/-	18.6135	µg/mL	Stressed
57	Diphenylamine		1,713.4	µg/mL	+/-	9.9619	µg/mL	Gravimetric
	CAS # 122-39-4	(Lot 07525MF)			+/-	18.7595	µg/mL	Unstressed
	Purity 99%				+/-	31.8192	µg/mL	Stressed
58	4-Nitroaniline		1,002.8	µg/mL	+/-	5.8304	µg/mL	Gravimetric
	CAS # 100-01-6	(Lot BCBG4702V)			+/-	10.9794	µg/mL	Unstressed
	Purity 99%				+/-	18.6228	µg/mL	Stressed
59	4,6-Dinitro-2-methylphenol (Dinitro-o-cresol)		2,002.0	µg/mL	+/-	11.6398	µg/mL	Gravimetric
	CAS # 534-52-1	(Lot LC06195V)			+/-	21.9193	µg/mL	Unstressed
	Purity 99%				+/-	37.1787	µg/mL	Stressed
60	4-Bromophenyl phenyl ether		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 101-55-3	(Lot STBB9729V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
61	Hexachlorobenzene		1,002.1	µg/mL	+/-	5.8260	µg/mL	Gravimetric
	CAS # 118-74-1	(Lot LC04221V)			+/-	10.9711	µg/mL	Unstressed
	Purity 98%				+/-	18.6089	µg/mL	Stressed
62	Pentachlorophenol		2,000.3	µg/mL	+/-	11.6299	µg/mL	Gravimetric
	CAS # 87-86-5	(Lot 140626JLM)			+/-	21.9007	µg/mL	Unstressed
	Purity 99%				+/-	37.1471	µg/mL	Stressed
63	Phenanthrene		999.0	µg/mL	+/-	5.8083	µg/mL	Gravimetric
	CAS # 85-01-8	(Lot MKBL6906V)			+/-	10.9379	µg/mL	Unstressed
	Purity 98%				+/-	18.5524	µg/mL	Stressed
64	n-Octadecane (C18)		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 593-45-3	(Lot OGCDK)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
65	Anthracene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-12-7	(Lot MKBK5208V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
66	Carbazole		1,000.1	µg/mL	+/-	5.8146	µg/mL	Gravimetric
	CAS # 86-74-8	(Lot S42950-417)			+/-	10.9497	µg/mL	Unstressed
	Purity 98%				+/-	18.5725	µg/mL	Stressed
67	Di-n-butylphthalate		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 84-74-2	(Lot MKBL8501V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
68	Fluoranthene		999.7	µg/mL	+/-	5.8123	µg/mL	Gravimetric
	CAS # 206-44-0	(Lot MKBQ6360V)			+/-	10.9454	µg/mL	Unstressed
	Purity 98%				+/-	18.5652	µg/mL	Stressed
69	Pyrene		999.1	µg/mL	+/-	5.8089	µg/mL	Gravimetric
	CAS # 129-00-0	(Lot BCBJ0984V)			+/-	10.9390	µg/mL	Unstressed
	Purity 98%				+/-	18.5543	µg/mL	Stressed
70	Benzyl butyl phthalate		1,001.2	µg/mL	+/-	5.8211	µg/mL	Gravimetric
	CAS # 85-68-7	(Lot 03027HV)			+/-	10.9618	µg/mL	Unstressed
	Purity 99%				+/-	18.5931	µg/mL	Stressed
71	Benz(a)anthracene		1,001.4	µg/mL	+/-	5.8222	µg/mL	Gravimetric
	CAS # 56-55-3	(Lot ER031412-01)			+/-	10.9640	µg/mL	Unstressed
	Purity 99%				+/-	18.5968	µg/mL	Stressed

72	Chrysene		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 218-01-9	(Lot ER120810-02)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
73	Bis(2-ethylhexyl)phthalate		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 117-81-7	(Lot MKBK2695V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
74	Di-n-octyl phthalate		1,002.1	µg/mL	+/-	5.8263	µg/mL	Gravimetric
	CAS # 117-84-0	(Lot 3128600)			+/-	10.9717	µg/mL	Unstressed
	Purity 99%				+/-	18.6098	µg/mL	Stressed
75	Benzo(b)fluoranthene		1,000.1	µg/mL	+/-	5.8147	µg/mL	Gravimetric
	CAS # 205-99-2	(Lot ER03101401)			+/-	10.9498	µg/mL	Unstressed
	Purity 99%				+/-	18.5726	µg/mL	Stressed
76	Benzo(k)fluoranthene		1,003.3	µg/mL	+/-	5.8333	µg/mL	Gravimetric
	CAS # 207-08-9	(Lot ER041513-01)			+/-	10.9848	µg/mL	Unstressed
	Purity 99%				+/-	18.6321	µg/mL	Stressed
77	Benzo(a)pyrene		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 50-32-8	(Lot ER071309-02)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
78	Indeno(1,2,3-cd)pyrene		1,000.6	µg/mL	+/-	5.8176	µg/mL	Gravimetric
	CAS # 193-39-5	(Lot ER082107-02)			+/-	10.9553	µg/mL	Unstressed
	Purity 99%				+/-	18.5819	µg/mL	Stressed
79	Dibenz(a,h)anthracene		1,004.4	µg/mL	+/-	5.8397	µg/mL	Gravimetric
	CAS # 53-70-3	(Lot ER032211-01)			+/-	10.9969	µg/mL	Unstressed
	Purity 99%				+/-	18.6525	µg/mL	Stressed
80	Benzo(g,h,i)perylene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 191-24-2	(Lot ER020708-08)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Specific Reference Material Notes:

N-nitrosodiphenylamine 2000 ug/mL equivalent when used for GC analysis. Actual formulation is diphenylamine 1710 ug/mL. N-Nitrosodiphenylamine is prone to breakdown in the injection port and will be converted to diphenylamine. N-Nitrosodiphenylamine is also a reactive species that can initiate premature decomposition of other compounds in the mix. For these reasons diphenylamine is used in the preparation of this mixture. When comparing the response of this compound to mixtures manufactured using N-nitrosodiphenylamine, a difference in response will be observed.

Column:
30m x 0.25mm x 0.25µm
Rtx-5 (cat.#10223)

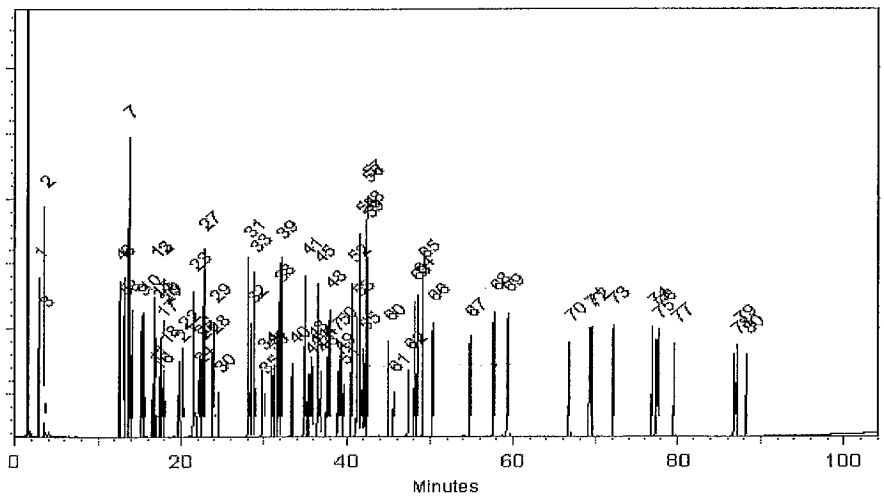
Carrier Gas:
hydrogen-constant pressure 10 psi

Temp. Program:
35°C (hold 3 min.) to 330°C
@ 3°C/min. (hold 3 min.)

Inj. Temp:
250°C

Det. Temp:
300°C

Det. Type:
FID



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

F. Joseph Tallon - Mix Technician **Date Mixed:** 24-Nov-2014 **Balance:** 1128360905

Jodi E. Breon - QA Analyst **Date Passed:** 05-Dec-2014

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

k is a coverage factor of 2, which gives a level of confidence of approximately 95%.

- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

RESTEK®
110 Benner Circle
Bellefonte, PA 16823-8812
Tel: (800)356-1688
Fax: (814)353-1309

www.restek.com

CERTIFIED REFERENCE MATERIAL

Certificate of Analysis

ACCREDITED
ISO Guide 34 Accredited
Reference Material Producer
Certificate #3222.01

ilac-MRA

ACCREDITED
ISO/IEC 17025 Accredited
Testing Laboratory
Certificate #3222.02

FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 569729 **Lot No.:** A0107399

Description : 8270 List 1 / Std #1 MegaMix (2015)

8270 List 1 / Std #1 MegaMix (2015) 500-2000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 10 mL **Pkg Amt:** > 5 mL

Expiration Date : May 31, 2016 **Storage:** 10°C or colder

Handling: Carcinogen/reproductive toxin. Photosensitive. Sonicate.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)	
1	1,4-Dioxane **CAS #** 123-91-1 (Lot SHBD8744V) **Purity** 99%	1,004.4 µg/mL	+/- 5.8397 µg/mL +/- 10.9969 µg/mL +/- 18.6525 µg/mL	Gravimetric Unstressed Stressed
2	Pyridine **CAS #** 110-86-1 (Lot SHBC7174V) **Purity** 99%	1,001.0 µg/mL	+/- 5.8199 µg/mL +/- 10.9596 µg/mL +/- 18.5894 µg/mL	Gravimetric Unstressed Stressed
3	N-Nitrosodimethylamine **CAS #** 62-75-9 (Lot 3213100) **Purity** 99%	1,000.2 µg/mL	+/- 5.8152 µg/mL +/- 10.9509 µg/mL +/- 18.5745 µg/mL	Gravimetric Unstressed Stressed
4	Aniline **CAS #** 62-53-3 (Lot K22Z462) **Purity** 99%	1,002.3 µg/mL	+/- 5.8275 µg/mL +/- 10.9739 µg/mL +/- 18.6135 µg/mL	Gravimetric Unstressed Stressed
5	Bis(2-chloroethyl)ether **CAS #** 111-44-4 (Lot 45296HKV) **Purity** 99%	1,001.4 µg/mL	+/- 5.8222 µg/mL +/- 10.9640 µg/mL +/- 18.5968 µg/mL	Gravimetric Unstressed Stressed
6	2-Chlorophenol **CAS #** 95-57-8 (Lot MKBD3900V) **Purity** 99%	1,000.8 µg/mL	+/- 5.8187 µg/mL +/- 10.9575 µg/mL +/- 18.5856 µg/mL	Gravimetric Unstressed Stressed
7	Phenol **CAS #** 108-95-2 (Lot SHBC6998V) **Purity** 99%	1,006.9 µg/mL	+/- 5.8542 µg/mL +/- 11.0242 µg/mL +/- 18.6989 µg/mL	Gravimetric Unstressed Stressed

8	n-Decane (C10)		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 124-18-5	(Lot SHBF1587V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
9	1,4-Dichlorobenzene		1,005.3	µg/mL	+/-	5.8449	µg/mL	Gravimetric
	CAS # 106-46-7	(Lot MKBL3891V)			+/-	11.0067	µg/mL	Unstressed
	Purity 99%				+/-	18.6692	µg/mL	Stressed
10	1,3-Dichlorobenzene		1,002.0	µg/mL	+/-	5.8257	µg/mL	Gravimetric
	CAS # 541-73-1	(Lot BCBC1891V)			+/-	10.9706	µg/mL	Unstressed
	Purity 99%				+/-	18.6079	µg/mL	Stressed
11	1,2-Dichlorobenzene		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 95-50-1	(Lot 68996CMV)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
12	Benzyl alcohol		1,000.4	µg/mL	+/-	5.8164	µg/mL	Gravimetric
	CAS # 100-51-6	(Lot SHBC1850V)			+/-	10.9531	µg/mL	Unstressed
	Purity 99%				+/-	18.5782	µg/mL	Stressed
13	2,2'-oxybis(1-chloropropane)		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 108-60-1	(Lot 2-EAW-18-3)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
14	2-Methylphenol (o-cresol)		1,003.6	µg/mL	+/-	5.8350	µg/mL	Gravimetric
	CAS # 95-48-7	(Lot SHBC1479V)			+/-	10.9881	µg/mL	Unstressed
	Purity 99%				+/-	18.6376	µg/mL	Stressed
15	Hexachloroethane		1,005.9	µg/mL	+/-	5.8484	µg/mL	Gravimetric
	CAS # 67-72-1	(Lot 4H3SF)			+/-	11.0133	µg/mL	Unstressed
	Purity 99%				+/-	18.6804	µg/mL	Stressed
16	Acetophenone		1,002.7	µg/mL	+/-	5.8298	µg/mL	Gravimetric
	CAS # 98-86-2	(Lot MKBR7156V)			+/-	10.9783	µg/mL	Unstressed
	Purity 99%				+/-	18.6209	µg/mL	Stressed
17	N-Nitroso-di-n-propylamine		1,003.9	µg/mL	+/-	5.8368	µg/mL	Gravimetric
	CAS # 621-64-7	(Lot OPAGF)			+/-	10.9914	µg/mL	Unstressed
	Purity 99%				+/-	18.6432	µg/mL	Stressed
18	4-Methylphenol (p-cresol)		500.4	µg/mL	+/-	2.9161	µg/mL	Gravimetric
	CAS # 106-44-5	(Lot 49396APV)			+/-	5.4823	µg/mL	Unstressed
	Purity 99%				+/-	9.2949	µg/mL	Stressed
19	3-Methylphenol (m-cresol)		500.2	µg/mL	+/-	2.9149	µg/mL	Gravimetric
	CAS # 108-39-4	(Lot SHBD0627V)			+/-	5.4801	µg/mL	Unstressed
	Purity 99%				+/-	9.2912	µg/mL	Stressed
20	Nitrobenzene		1,001.1	µg/mL	+/-	5.8205	µg/mL	Gravimetric
	CAS # 98-95-3	(Lot 65096APV)			+/-	10.9607	µg/mL	Unstressed
	Purity 99%				+/-	18.5912	µg/mL	Stressed
21	Isophorone		999.3	µg/mL	+/-	5.8100	µg/mL	Gravimetric
	CAS # 78-59-1	(Lot 06705DE)			+/-	10.9410	µg/mL	Unstressed
	Purity 97%				+/-	18.5577	µg/mL	Stressed
22	2-Nitrophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 88-75-5	(Lot BCBH7602V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
23	2,4-Dimethylphenol		1,003.2	µg/mL	+/-	5.8327	µg/mL	Gravimetric
	CAS # 105-67-9	(Lot 10165155)			+/-	10.9837	µg/mL	Unstressed
	Purity 99%				+/-	18.6302	µg/mL	Stressed

24	Bis(2-chloroethoxy)methane		1,004.5	µg/mL	+/-	5.8402	µg/mL	Gravimetric
	CAS # 111-91-1	(Lot 317200)			+/-	10.9980	µg/mL	Unstressed
	Purity 99%				+/-	18.6544	µg/mL	Stressed
25	2,4-Dichlorophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-83-2	(Lot BCBH1617V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
26	1,2,4-Trichlorobenzene		1,000.6	µg/mL	+/-	5.8176	µg/mL	Gravimetric
	CAS # 120-82-1	(Lot 26896BM)			+/-	10.9553	µg/mL	Unstressed
	Purity 99%				+/-	18.5819	µg/mL	Stressed
27	Naphthalene		1,002.3	µg/mL	+/-	5.8275	µg/mL	Gravimetric
	CAS # 91-20-3	(Lot MKBH4351V)			+/-	10.9739	µg/mL	Unstressed
	Purity 99%				+/-	18.6135	µg/mL	Stressed
28	2,6-Dichlorophenol		1,000.1	µg/mL	+/-	5.8147	µg/mL	Gravimetric
	CAS # 87-65-0	(Lot MKBN2776V)			+/-	10.9498	µg/mL	Unstressed
	Purity 99%				+/-	18.5726	µg/mL	Stressed
29	4-Chloroaniline		1,000.3	µg/mL	+/-	5.8157	µg/mL	Gravimetric
	CAS # 106-47-8	(Lot 12528PH)			+/-	10.9518	µg/mL	Unstressed
	Purity 98%				+/-	18.5761	µg/mL	Stressed
30	Hexachlorobutadiene		999.9	µg/mL	+/-	5.8135	µg/mL	Gravimetric
	CAS # 87-68-3	(Lot K22W009)			+/-	10.9475	µg/mL	Unstressed
	Purity 98%				+/-	18.5688	µg/mL	Stressed
31	2-Methylnaphthalene		998.6	µg/mL	+/-	5.8059	µg/mL	Gravimetric
	CAS # 91-57-6	(Lot 19399MJV)			+/-	10.9333	µg/mL	Unstressed
	Purity 96%				+/-	18.5446	µg/mL	Stressed
32	4-Chloro-3-methylphenol		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 59-50-7	(Lot STBC0769V)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
33	1-Methylnaphthalene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 90-12-0	(Lot 525000-10)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
34	1,2,4,5-Tetrachlorobenzene		1,000.2	µg/mL	+/-	5.8152	µg/mL	Gravimetric
	CAS # 95-94-3	(Lot 06024AIV)			+/-	10.9509	µg/mL	Unstressed
	Purity 99%				+/-	18.5745	µg/mL	Stressed
35	Hexachlorocyclopentadiene		1,002.4	µg/mL	+/-	5.8280	µg/mL	Gravimetric
	CAS # 77-47-4	(Lot 3140300)			+/-	10.9750	µg/mL	Unstressed
	Purity 99%				+/-	18.6154	µg/mL	Stressed
36	2,4,6-Trichlorophenol		1,001.4	µg/mL	+/-	5.8222	µg/mL	Gravimetric
	CAS # 88-06-2	(Lot MKBH7393V)			+/-	10.9640	µg/mL	Unstressed
	Purity 99%				+/-	18.5968	µg/mL	Stressed
37	2,4,5-Trichlorophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-95-4	(Lot FHM01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
38	2-Chloronaphthalene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-58-7	(Lot FIJ01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
39	Biphenyl		1,006.1	µg/mL	+/-	5.8496	µg/mL	Gravimetric
	CAS # 92-52-4	(Lot 1277976)			+/-	11.0155	µg/mL	Unstressed
	Purity 99%				+/-	18.6841	µg/mL	Stressed

40	2-Nitroaniline		1,004.6	µg/mL	+/-	5.8408	µg/mL	Gravimetric
	CAS # 88-74-4	(Lot MKBK7597V)			+/-	10.9991	µg/mL	Unstressed
	Purity 99%				+/-	18.6562	µg/mL	Stressed
41	Acenaphthylene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 208-96-8	(Lot ER030707-01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
42	1,3-Dinitrobenzene		1,006.4	µg/mL	+/-	5.8513	µg/mL	Gravimetric
	CAS # 99-65-0	(Lot BCBB1436V)			+/-	11.0188	µg/mL	Unstressed
	Purity 99%				+/-	18.6896	µg/mL	Stressed
43	Dimethylphthalate		1,001.0	µg/mL	+/-	5.8199	µg/mL	Gravimetric
	CAS # 131-11-3	(Lot 10117699)			+/-	10.9596	µg/mL	Unstressed
	Purity 99%				+/-	18.5894	µg/mL	Stressed
44	2,6-Dinitrotoluene		1,001.6	µg/mL	+/-	5.8234	µg/mL	Gravimetric
	CAS # 606-20-2	(Lot 1437483V)			+/-	10.9662	µg/mL	Unstressed
	Purity 99%				+/-	18.6005	µg/mL	Stressed
45	Acenaphthene		1,001.8	µg/mL	+/-	5.8246	µg/mL	Gravimetric
	CAS # 83-32-9	(Lot MKBP0384V)			+/-	10.9684	µg/mL	Unstressed
	Purity 99%				+/-	18.6042	µg/mL	Stressed
46	2,4-Dinitrophenol		2,008.7	µg/mL	+/-	11.6788	µg/mL	Gravimetric
	CAS # 51-28-5	(Lot MKBP5833V)			+/-	21.9927	µg/mL	Unstressed
	Purity 99%				+/-	37.3031	µg/mL	Stressed
47	Dibenzofuran		1,001.8	µg/mL	+/-	5.8246	µg/mL	Gravimetric
	CAS # 132-64-9	(Lot MKBK2375V)			+/-	10.9684	µg/mL	Unstressed
	Purity 99%				+/-	18.6042	µg/mL	Stressed
48	3-Nitroaniline		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 99-09-2	(Lot MKBH5131V)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
49	2,4-Dinitrotoluene		1,001.7	µg/mL	+/-	5.8240	µg/mL	Gravimetric
	CAS # 121-14-2	(Lot MKAA0690V)			+/-	10.9673	µg/mL	Unstressed
	Purity 99%				+/-	18.6024	µg/mL	Stressed
50	4-Nitrophenol		2,001.0	µg/mL	+/-	11.6340	µg/mL	Gravimetric
	CAS # 100-02-7	(Lot MKBK1842V)			+/-	21.9083	µg/mL	Unstressed
	Purity 99%				+/-	37.1601	µg/mL	Stressed
51	2,3,4,6-Tetrachlorophenol		1,002.1	µg/mL	+/-	5.8263	µg/mL	Gravimetric
	CAS # 58-90-2	(Lot FN10221307)			+/-	10.9717	µg/mL	Unstressed
	Purity 99%				+/-	18.6098	µg/mL	Stressed
52	Fluorene		1,000.5	µg/mL	+/-	5.8169	µg/mL	Gravimetric
	CAS # 86-73-7	(Lot 10174662)			+/-	10.9540	µg/mL	Unstressed
	Purity 98%				+/-	18.5797	µg/mL	Stressed
53	4-Chlorophenyl phenyl ether		1,000.4	µg/mL	+/-	5.8164	µg/mL	Gravimetric
	CAS # 7005-72-3	(Lot MKBL1347V)			+/-	10.9531	µg/mL	Unstressed
	Purity 99%				+/-	18.5782	µg/mL	Stressed
54	n-Hexadecane (C16)		1,002.2	µg/mL	+/-	5.8269	µg/mL	Gravimetric
	CAS # 544-76-3	(Lot SHBD4570V)			+/-	10.9728	µg/mL	Unstressed
	Purity 99%				+/-	18.6116	µg/mL	Stressed
55	Diethylphthalate		1,001.1	µg/mL	+/-	5.8205	µg/mL	Gravimetric
	CAS # 84-66-2	(Lot MKBJ3578V)			+/-	10.9607	µg/mL	Unstressed
	Purity 99%				+/-	18.5912	µg/mL	Stressed

56	Azobenzene		1,002.3	µg/mL	+/-	5.8275	µg/mL	Gravimetric
	CAS # 103-33-3	(Lot MKBS2559V)			+/-	10.9739	µg/mL	Unstressed
	Purity 99%				+/-	18.6135	µg/mL	Stressed
57	Diphenylamine		1,713.4	µg/mL	+/-	9.9619	µg/mL	Gravimetric
	CAS # 122-39-4	(Lot 07525MF)			+/-	18.7595	µg/mL	Unstressed
	Purity 99%				+/-	31.8192	µg/mL	Stressed
58	4-Nitroaniline		1,002.8	µg/mL	+/-	5.8304	µg/mL	Gravimetric
	CAS # 100-01-6	(Lot BCBG4702V)			+/-	10.9794	µg/mL	Unstressed
	Purity 99%				+/-	18.6228	µg/mL	Stressed
59	4,6-Dinitro-2-methylphenol (Dinitro-o-cresol)		2,002.0	µg/mL	+/-	11.6398	µg/mL	Gravimetric
	CAS # 534-52-1	(Lot LC06195V)			+/-	21.9193	µg/mL	Unstressed
	Purity 99%				+/-	37.1787	µg/mL	Stressed
60	4-Bromophenyl phenyl ether		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 101-55-3	(Lot STBB9729V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
61	Hexachlorobenzene		1,002.1	µg/mL	+/-	5.8260	µg/mL	Gravimetric
	CAS # 118-74-1	(Lot LC04221V)			+/-	10.9711	µg/mL	Unstressed
	Purity 98%				+/-	18.6089	µg/mL	Stressed
62	Pentachlorophenol		2,000.3	µg/mL	+/-	11.6299	µg/mL	Gravimetric
	CAS # 87-86-5	(Lot 140626JLM)			+/-	21.9007	µg/mL	Unstressed
	Purity 99%				+/-	37.1471	µg/mL	Stressed
63	Phenanthrene		999.0	µg/mL	+/-	5.8083	µg/mL	Gravimetric
	CAS # 85-01-8	(Lot MKBL6906V)			+/-	10.9379	µg/mL	Unstressed
	Purity 98%				+/-	18.5524	µg/mL	Stressed
64	n-Octadecane (C18)		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 593-45-3	(Lot OGCDK)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
65	Anthracene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-12-7	(Lot MKBK5208V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
66	Carbazole		1,000.1	µg/mL	+/-	5.8146	µg/mL	Gravimetric
	CAS # 86-74-8	(Lot S42950-417)			+/-	10.9497	µg/mL	Unstressed
	Purity 98%				+/-	18.5725	µg/mL	Stressed
67	Di-n-butylphthalate		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 84-74-2	(Lot MKBL8501V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
68	Fluoranthene		999.7	µg/mL	+/-	5.8123	µg/mL	Gravimetric
	CAS # 206-44-0	(Lot MKBQ6360V)			+/-	10.9454	µg/mL	Unstressed
	Purity 98%				+/-	18.5652	µg/mL	Stressed
69	Pyrene		999.1	µg/mL	+/-	5.8089	µg/mL	Gravimetric
	CAS # 129-00-0	(Lot BCBJ0984V)			+/-	10.9390	µg/mL	Unstressed
	Purity 98%				+/-	18.5543	µg/mL	Stressed
70	Benzyl butyl phthalate		1,001.2	µg/mL	+/-	5.8211	µg/mL	Gravimetric
	CAS # 85-68-7	(Lot 03027HV)			+/-	10.9618	µg/mL	Unstressed
	Purity 99%				+/-	18.5931	µg/mL	Stressed
71	Benz(a)anthracene		1,001.4	µg/mL	+/-	5.8222	µg/mL	Gravimetric
	CAS # 56-55-3	(Lot ER031412-01)			+/-	10.9640	µg/mL	Unstressed
	Purity 99%				+/-	18.5968	µg/mL	Stressed

72	Chrysene		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 218-01-9	(Lot ER120810-02)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
73	Bis(2-ethylhexyl)phthalate		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 117-81-7	(Lot MKBK2695V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
74	Di-n-octyl phthalate		1,002.1	µg/mL	+/-	5.8263	µg/mL	Gravimetric
	CAS # 117-84-0	(Lot 3128600)			+/-	10.9717	µg/mL	Unstressed
	Purity 99%				+/-	18.6098	µg/mL	Stressed
75	Benzo(b)fluoranthene		1,000.1	µg/mL	+/-	5.8147	µg/mL	Gravimetric
	CAS # 205-99-2	(Lot ER03101401)			+/-	10.9498	µg/mL	Unstressed
	Purity 99%				+/-	18.5726	µg/mL	Stressed
76	Benzo(k)fluoranthene		1,003.3	µg/mL	+/-	5.8333	µg/mL	Gravimetric
	CAS # 207-08-9	(Lot ER041513-01)			+/-	10.9848	µg/mL	Unstressed
	Purity 99%				+/-	18.6321	µg/mL	Stressed
77	Benzo(a)pyrene		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 50-32-8	(Lot ER071309-02)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
78	Indeno(1,2,3-cd)pyrene		1,000.6	µg/mL	+/-	5.8176	µg/mL	Gravimetric
	CAS # 193-39-5	(Lot ER082107-02)			+/-	10.9553	µg/mL	Unstressed
	Purity 99%				+/-	18.5819	µg/mL	Stressed
79	Dibenz(a,h)anthracene		1,004.4	µg/mL	+/-	5.8397	µg/mL	Gravimetric
	CAS # 53-70-3	(Lot ER032211-01)			+/-	10.9969	µg/mL	Unstressed
	Purity 99%				+/-	18.6525	µg/mL	Stressed
80	Benzo(g,h,i)perylene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 191-24-2	(Lot ER020708-08)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Specific Reference Material Notes:

N-nitrosodiphenylamine 2000 ug/mL equivalent when used for GC analysis. Actual formulation is diphenylamine 1710 ug/mL. N-Nitrosodiphenylamine is prone to breakdown in the injection port and will be converted to diphenylamine. N-Nitrosodiphenylamine is also a reactive species that can initiate premature decomposition of other compounds in the mix. For these reasons diphenylamine is used in the preparation of this mixture. When comparing the response of this compound to mixtures manufactured using N-nitrosodiphenylamine, a difference in response will be observed.

Column:
30m x 0.25mm x 0.25µm
Rtx-5 (cat.#10223)

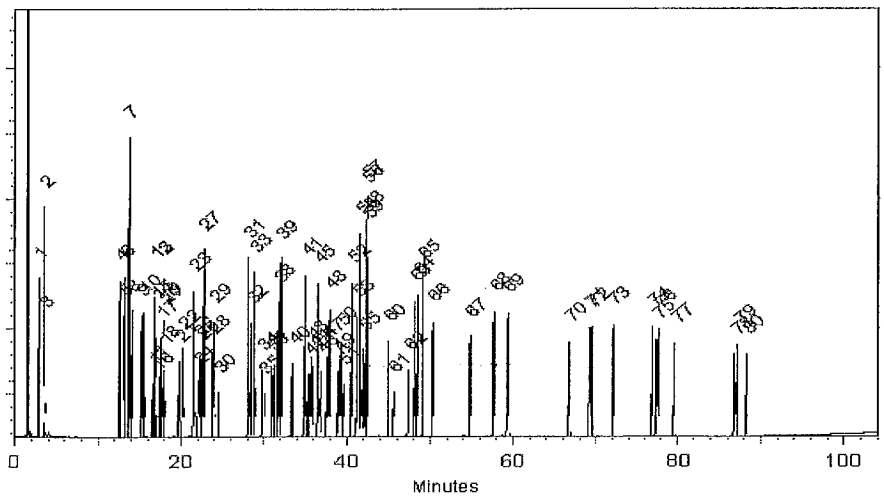
Carrier Gas:
hydrogen-constant pressure 10 psi

Temp. Program:
35°C (hold 3 min.) to 330°C
@ 3°C/min. (hold 3 min.)

Inj. Temp:
250°C

Det. Temp:
300°C

Det. Type:
FID



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

F. Joseph Tallon - Mix Technician **Date Mixed:** 24-Nov-2014 **Balance:** 1128360905

Jodi E. Breon - QA Analyst **Date Passed:** 05-Dec-2014

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

k is a coverage factor of 2, which gives a level of confidence of approximately 95%.

- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

RESTEK®
110 Benner Circle
Bellefonte, PA 16823-8812
Tel: (800)356-1688
Fax: (814)353-1309

www.restek.com

CERTIFIED REFERENCE MATERIAL

Certificate of Analysis

ACCREDITED
ISO Guide 34 Accredited
Reference Material Producer
Certificate #3222.01

ACCREDITED
ISO/IEC 17025 Accredited
Testing Laboratory
Certificate #3222.02

FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 569729 **Lot No.:** A0107399

Description : 8270 List 1 / Std #1 MegaMix (2015)

8270 List 1 / Std #1 MegaMix (2015) 500-2000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 10 mL **Pkg Amt:** > 5 mL

Expiration Date : May 31, 2016 **Storage:** 10°C or colder

Handling: Carcinogen/reproductive toxin. Photosensitive. Sonicate.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)	
1	1,4-Dioxane **CAS #** 123-91-1 (Lot SHBD8744V) **Purity** 99%	1,004.4 µg/mL	+/- 5.8397 µg/mL +/- 10.9969 µg/mL +/- 18.6525 µg/mL	Gravimetric Unstressed Stressed
2	Pyridine **CAS #** 110-86-1 (Lot SHBC7174V) **Purity** 99%	1,001.0 µg/mL	+/- 5.8199 µg/mL +/- 10.9596 µg/mL +/- 18.5894 µg/mL	Gravimetric Unstressed Stressed
3	N-Nitrosodimethylamine **CAS #** 62-75-9 (Lot 3213100) **Purity** 99%	1,000.2 µg/mL	+/- 5.8152 µg/mL +/- 10.9509 µg/mL +/- 18.5745 µg/mL	Gravimetric Unstressed Stressed
4	Aniline **CAS #** 62-53-3 (Lot K22Z462) **Purity** 99%	1,002.3 µg/mL	+/- 5.8275 µg/mL +/- 10.9739 µg/mL +/- 18.6135 µg/mL	Gravimetric Unstressed Stressed
5	Bis(2-chloroethyl)ether **CAS #** 111-44-4 (Lot 45296HKV) **Purity** 99%	1,001.4 µg/mL	+/- 5.8222 µg/mL +/- 10.9640 µg/mL +/- 18.5968 µg/mL	Gravimetric Unstressed Stressed
6	2-Chlorophenol **CAS #** 95-57-8 (Lot MKBD3900V) **Purity** 99%	1,000.8 µg/mL	+/- 5.8187 µg/mL +/- 10.9575 µg/mL +/- 18.5856 µg/mL	Gravimetric Unstressed Stressed
7	Phenol **CAS #** 108-95-2 (Lot SHBC6998V) **Purity** 99%	1,006.9 µg/mL	+/- 5.8542 µg/mL +/- 11.0242 µg/mL +/- 18.6989 µg/mL	Gravimetric Unstressed Stressed

8	n-Decane (C10)		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 124-18-5	(Lot SHBF1587V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
9	1,4-Dichlorobenzene		1,005.3	µg/mL	+/-	5.8449	µg/mL	Gravimetric
	CAS # 106-46-7	(Lot MKBL3891V)			+/-	11.0067	µg/mL	Unstressed
	Purity 99%				+/-	18.6692	µg/mL	Stressed
10	1,3-Dichlorobenzene		1,002.0	µg/mL	+/-	5.8257	µg/mL	Gravimetric
	CAS # 541-73-1	(Lot BCBC1891V)			+/-	10.9706	µg/mL	Unstressed
	Purity 99%				+/-	18.6079	µg/mL	Stressed
11	1,2-Dichlorobenzene		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 95-50-1	(Lot 68996CMV)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
12	Benzyl alcohol		1,000.4	µg/mL	+/-	5.8164	µg/mL	Gravimetric
	CAS # 100-51-6	(Lot SHBC1850V)			+/-	10.9531	µg/mL	Unstressed
	Purity 99%				+/-	18.5782	µg/mL	Stressed
13	2,2'-oxybis(1-chloropropane)		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 108-60-1	(Lot 2-EAW-18-3)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
14	2-Methylphenol (o-cresol)		1,003.6	µg/mL	+/-	5.8350	µg/mL	Gravimetric
	CAS # 95-48-7	(Lot SHBC1479V)			+/-	10.9881	µg/mL	Unstressed
	Purity 99%				+/-	18.6376	µg/mL	Stressed
15	Hexachloroethane		1,005.9	µg/mL	+/-	5.8484	µg/mL	Gravimetric
	CAS # 67-72-1	(Lot 4H3SF)			+/-	11.0133	µg/mL	Unstressed
	Purity 99%				+/-	18.6804	µg/mL	Stressed
16	Acetophenone		1,002.7	µg/mL	+/-	5.8298	µg/mL	Gravimetric
	CAS # 98-86-2	(Lot MKBR7156V)			+/-	10.9783	µg/mL	Unstressed
	Purity 99%				+/-	18.6209	µg/mL	Stressed
17	N-Nitroso-di-n-propylamine		1,003.9	µg/mL	+/-	5.8368	µg/mL	Gravimetric
	CAS # 621-64-7	(Lot OPAGF)			+/-	10.9914	µg/mL	Unstressed
	Purity 99%				+/-	18.6432	µg/mL	Stressed
18	4-Methylphenol (p-cresol)		500.4	µg/mL	+/-	2.9161	µg/mL	Gravimetric
	CAS # 106-44-5	(Lot 49396APV)			+/-	5.4823	µg/mL	Unstressed
	Purity 99%				+/-	9.2949	µg/mL	Stressed
19	3-Methylphenol (m-cresol)		500.2	µg/mL	+/-	2.9149	µg/mL	Gravimetric
	CAS # 108-39-4	(Lot SHBD0627V)			+/-	5.4801	µg/mL	Unstressed
	Purity 99%				+/-	9.2912	µg/mL	Stressed
20	Nitrobenzene		1,001.1	µg/mL	+/-	5.8205	µg/mL	Gravimetric
	CAS # 98-95-3	(Lot 65096APV)			+/-	10.9607	µg/mL	Unstressed
	Purity 99%				+/-	18.5912	µg/mL	Stressed
21	Isophorone		999.3	µg/mL	+/-	5.8100	µg/mL	Gravimetric
	CAS # 78-59-1	(Lot 06705DE)			+/-	10.9410	µg/mL	Unstressed
	Purity 97%				+/-	18.5577	µg/mL	Stressed
22	2-Nitrophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 88-75-5	(Lot BCBH7602V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
23	2,4-Dimethylphenol		1,003.2	µg/mL	+/-	5.8327	µg/mL	Gravimetric
	CAS # 105-67-9	(Lot 10165155)			+/-	10.9837	µg/mL	Unstressed
	Purity 99%				+/-	18.6302	µg/mL	Stressed

24	Bis(2-chloroethoxy)methane		1,004.5	µg/mL	+/-	5.8402	µg/mL	Gravimetric
	CAS # 111-91-1	(Lot 317200)			+/-	10.9980	µg/mL	Unstressed
	Purity 99%				+/-	18.6544	µg/mL	Stressed
25	2,4-Dichlorophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-83-2	(Lot BCBH1617V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
26	1,2,4-Trichlorobenzene		1,000.6	µg/mL	+/-	5.8176	µg/mL	Gravimetric
	CAS # 120-82-1	(Lot 26896BM)			+/-	10.9553	µg/mL	Unstressed
	Purity 99%				+/-	18.5819	µg/mL	Stressed
27	Naphthalene		1,002.3	µg/mL	+/-	5.8275	µg/mL	Gravimetric
	CAS # 91-20-3	(Lot MKBH4351V)			+/-	10.9739	µg/mL	Unstressed
	Purity 99%				+/-	18.6135	µg/mL	Stressed
28	2,6-Dichlorophenol		1,000.1	µg/mL	+/-	5.8147	µg/mL	Gravimetric
	CAS # 87-65-0	(Lot MKBN2776V)			+/-	10.9498	µg/mL	Unstressed
	Purity 99%				+/-	18.5726	µg/mL	Stressed
29	4-Chloroaniline		1,000.3	µg/mL	+/-	5.8157	µg/mL	Gravimetric
	CAS # 106-47-8	(Lot 12528PH)			+/-	10.9518	µg/mL	Unstressed
	Purity 98%				+/-	18.5761	µg/mL	Stressed
30	Hexachlorobutadiene		999.9	µg/mL	+/-	5.8135	µg/mL	Gravimetric
	CAS # 87-68-3	(Lot K22W009)			+/-	10.9475	µg/mL	Unstressed
	Purity 98%				+/-	18.5688	µg/mL	Stressed
31	2-Methylnaphthalene		998.6	µg/mL	+/-	5.8059	µg/mL	Gravimetric
	CAS # 91-57-6	(Lot 19399MJV)			+/-	10.9333	µg/mL	Unstressed
	Purity 96%				+/-	18.5446	µg/mL	Stressed
32	4-Chloro-3-methylphenol		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 59-50-7	(Lot STBC0769V)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
33	1-Methylnaphthalene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 90-12-0	(Lot 525000-10)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
34	1,2,4,5-Tetrachlorobenzene		1,000.2	µg/mL	+/-	5.8152	µg/mL	Gravimetric
	CAS # 95-94-3	(Lot 06024AIV)			+/-	10.9509	µg/mL	Unstressed
	Purity 99%				+/-	18.5745	µg/mL	Stressed
35	Hexachlorocyclopentadiene		1,002.4	µg/mL	+/-	5.8280	µg/mL	Gravimetric
	CAS # 77-47-4	(Lot 3140300)			+/-	10.9750	µg/mL	Unstressed
	Purity 99%				+/-	18.6154	µg/mL	Stressed
36	2,4,6-Trichlorophenol		1,001.4	µg/mL	+/-	5.8222	µg/mL	Gravimetric
	CAS # 88-06-2	(Lot MKBH7393V)			+/-	10.9640	µg/mL	Unstressed
	Purity 99%				+/-	18.5968	µg/mL	Stressed
37	2,4,5-Trichlorophenol		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 95-95-4	(Lot FHM01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
38	2-Chloronaphthalene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 91-58-7	(Lot FIJ01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
39	Biphenyl		1,006.1	µg/mL	+/-	5.8496	µg/mL	Gravimetric
	CAS # 92-52-4	(Lot 1277976)			+/-	11.0155	µg/mL	Unstressed
	Purity 99%				+/-	18.6841	µg/mL	Stressed

40	2-Nitroaniline		1,004.6	µg/mL	+/-	5.8408	µg/mL	Gravimetric
	CAS # 88-74-4	(Lot MKBK7597V)			+/-	10.9991	µg/mL	Unstressed
	Purity 99%				+/-	18.6562	µg/mL	Stressed
41	Acenaphthylene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 208-96-8	(Lot ER030707-01)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
42	1,3-Dinitrobenzene		1,006.4	µg/mL	+/-	5.8513	µg/mL	Gravimetric
	CAS # 99-65-0	(Lot BCBB1436V)			+/-	11.0188	µg/mL	Unstressed
	Purity 99%				+/-	18.6896	µg/mL	Stressed
43	Dimethylphthalate		1,001.0	µg/mL	+/-	5.8199	µg/mL	Gravimetric
	CAS # 131-11-3	(Lot 10117699)			+/-	10.9596	µg/mL	Unstressed
	Purity 99%				+/-	18.5894	µg/mL	Stressed
44	2,6-Dinitrotoluene		1,001.6	µg/mL	+/-	5.8234	µg/mL	Gravimetric
	CAS # 606-20-2	(Lot 1437483V)			+/-	10.9662	µg/mL	Unstressed
	Purity 99%				+/-	18.6005	µg/mL	Stressed
45	Acenaphthene		1,001.8	µg/mL	+/-	5.8246	µg/mL	Gravimetric
	CAS # 83-32-9	(Lot MKBP0384V)			+/-	10.9684	µg/mL	Unstressed
	Purity 99%				+/-	18.6042	µg/mL	Stressed
46	2,4-Dinitrophenol		2,008.7	µg/mL	+/-	11.6788	µg/mL	Gravimetric
	CAS # 51-28-5	(Lot MKBP5833V)			+/-	21.9927	µg/mL	Unstressed
	Purity 99%				+/-	37.3031	µg/mL	Stressed
47	Dibenzofuran		1,001.8	µg/mL	+/-	5.8246	µg/mL	Gravimetric
	CAS # 132-64-9	(Lot MKBK2375V)			+/-	10.9684	µg/mL	Unstressed
	Purity 99%				+/-	18.6042	µg/mL	Stressed
48	3-Nitroaniline		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 99-09-2	(Lot MKBH5131V)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
49	2,4-Dinitrotoluene		1,001.7	µg/mL	+/-	5.8240	µg/mL	Gravimetric
	CAS # 121-14-2	(Lot MKAA0690V)			+/-	10.9673	µg/mL	Unstressed
	Purity 99%				+/-	18.6024	µg/mL	Stressed
50	4-Nitrophenol		2,001.0	µg/mL	+/-	11.6340	µg/mL	Gravimetric
	CAS # 100-02-7	(Lot MKBK1842V)			+/-	21.9083	µg/mL	Unstressed
	Purity 99%				+/-	37.1601	µg/mL	Stressed
51	2,3,4,6-Tetrachlorophenol		1,002.1	µg/mL	+/-	5.8263	µg/mL	Gravimetric
	CAS # 58-90-2	(Lot FN10221307)			+/-	10.9717	µg/mL	Unstressed
	Purity 99%				+/-	18.6098	µg/mL	Stressed
52	Fluorene		1,000.5	µg/mL	+/-	5.8169	µg/mL	Gravimetric
	CAS # 86-73-7	(Lot 10174662)			+/-	10.9540	µg/mL	Unstressed
	Purity 98%				+/-	18.5797	µg/mL	Stressed
53	4-Chlorophenyl phenyl ether		1,000.4	µg/mL	+/-	5.8164	µg/mL	Gravimetric
	CAS # 7005-72-3	(Lot MKBL1347V)			+/-	10.9531	µg/mL	Unstressed
	Purity 99%				+/-	18.5782	µg/mL	Stressed
54	n-Hexadecane (C16)		1,002.2	µg/mL	+/-	5.8269	µg/mL	Gravimetric
	CAS # 544-76-3	(Lot SHBD4570V)			+/-	10.9728	µg/mL	Unstressed
	Purity 99%				+/-	18.6116	µg/mL	Stressed
55	Diethylphthalate		1,001.1	µg/mL	+/-	5.8205	µg/mL	Gravimetric
	CAS # 84-66-2	(Lot MKBJ3578V)			+/-	10.9607	µg/mL	Unstressed
	Purity 99%				+/-	18.5912	µg/mL	Stressed

56	Azobenzene		1,002.3	µg/mL	+/-	5.8275	µg/mL	Gravimetric
	CAS # 103-33-3	(Lot MKBS2559V)			+/-	10.9739	µg/mL	Unstressed
	Purity 99%				+/-	18.6135	µg/mL	Stressed
57	Diphenylamine		1,713.4	µg/mL	+/-	9.9619	µg/mL	Gravimetric
	CAS # 122-39-4	(Lot 07525MF)			+/-	18.7595	µg/mL	Unstressed
	Purity 99%				+/-	31.8192	µg/mL	Stressed
58	4-Nitroaniline		1,002.8	µg/mL	+/-	5.8304	µg/mL	Gravimetric
	CAS # 100-01-6	(Lot BCBG4702V)			+/-	10.9794	µg/mL	Unstressed
	Purity 99%				+/-	18.6228	µg/mL	Stressed
59	4,6-Dinitro-2-methylphenol (Dinitro-o-cresol)		2,002.0	µg/mL	+/-	11.6398	µg/mL	Gravimetric
	CAS # 534-52-1	(Lot LC06195V)			+/-	21.9193	µg/mL	Unstressed
	Purity 99%				+/-	37.1787	µg/mL	Stressed
60	4-Bromophenyl phenyl ether		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 101-55-3	(Lot STBB9729V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
61	Hexachlorobenzene		1,002.1	µg/mL	+/-	5.8260	µg/mL	Gravimetric
	CAS # 118-74-1	(Lot LC04221V)			+/-	10.9711	µg/mL	Unstressed
	Purity 98%				+/-	18.6089	µg/mL	Stressed
62	Pentachlorophenol		2,000.3	µg/mL	+/-	11.6299	µg/mL	Gravimetric
	CAS # 87-86-5	(Lot 140626JLM)			+/-	21.9007	µg/mL	Unstressed
	Purity 99%				+/-	37.1471	µg/mL	Stressed
63	Phenanthrene		999.0	µg/mL	+/-	5.8083	µg/mL	Gravimetric
	CAS # 85-01-8	(Lot MKBL6906V)			+/-	10.9379	µg/mL	Unstressed
	Purity 98%				+/-	18.5524	µg/mL	Stressed
64	n-Octadecane (C18)		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 593-45-3	(Lot OGCDK)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
65	Anthracene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 120-12-7	(Lot MKBK5208V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
66	Carbazole		1,000.1	µg/mL	+/-	5.8146	µg/mL	Gravimetric
	CAS # 86-74-8	(Lot S42950-417)			+/-	10.9497	µg/mL	Unstressed
	Purity 98%				+/-	18.5725	µg/mL	Stressed
67	Di-n-butylphthalate		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 84-74-2	(Lot MKBL8501V)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed
68	Fluoranthene		999.7	µg/mL	+/-	5.8123	µg/mL	Gravimetric
	CAS # 206-44-0	(Lot MKBQ6360V)			+/-	10.9454	µg/mL	Unstressed
	Purity 98%				+/-	18.5652	µg/mL	Stressed
69	Pyrene		999.1	µg/mL	+/-	5.8089	µg/mL	Gravimetric
	CAS # 129-00-0	(Lot BCBJ0984V)			+/-	10.9390	µg/mL	Unstressed
	Purity 98%				+/-	18.5543	µg/mL	Stressed
70	Benzyl butyl phthalate		1,001.2	µg/mL	+/-	5.8211	µg/mL	Gravimetric
	CAS # 85-68-7	(Lot 03027HV)			+/-	10.9618	µg/mL	Unstressed
	Purity 99%				+/-	18.5931	µg/mL	Stressed
71	Benz(a)anthracene		1,001.4	µg/mL	+/-	5.8222	µg/mL	Gravimetric
	CAS # 56-55-3	(Lot ER031412-01)			+/-	10.9640	µg/mL	Unstressed
	Purity 99%				+/-	18.5968	µg/mL	Stressed

72	Chrysene		1,006.5	µg/mL	+/-	5.8519	µg/mL	Gravimetric
	CAS # 218-01-9	(Lot ER120810-02)			+/-	11.0199	µg/mL	Unstressed
	Purity 99%				+/-	18.6915	µg/mL	Stressed
73	Bis(2-ethylhexyl)phthalate		1,000.5	µg/mL	+/-	5.8170	µg/mL	Gravimetric
	CAS # 117-81-7	(Lot MKBK2695V)			+/-	10.9542	µg/mL	Unstressed
	Purity 99%				+/-	18.5801	µg/mL	Stressed
74	Di-n-octyl phthalate		1,002.1	µg/mL	+/-	5.8263	µg/mL	Gravimetric
	CAS # 117-84-0	(Lot 3128600)			+/-	10.9717	µg/mL	Unstressed
	Purity 99%				+/-	18.6098	µg/mL	Stressed
75	Benzo(b)fluoranthene		1,000.1	µg/mL	+/-	5.8147	µg/mL	Gravimetric
	CAS # 205-99-2	(Lot ER03101401)			+/-	10.9498	µg/mL	Unstressed
	Purity 99%				+/-	18.5726	µg/mL	Stressed
76	Benzo(k)fluoranthene		1,003.3	µg/mL	+/-	5.8333	µg/mL	Gravimetric
	CAS # 207-08-9	(Lot ER041513-01)			+/-	10.9848	µg/mL	Unstressed
	Purity 99%				+/-	18.6321	µg/mL	Stressed
77	Benzo(a)pyrene		1,001.3	µg/mL	+/-	5.8216	µg/mL	Gravimetric
	CAS # 50-32-8	(Lot ER071309-02)			+/-	10.9629	µg/mL	Unstressed
	Purity 99%				+/-	18.5949	µg/mL	Stressed
78	Indeno(1,2,3-cd)pyrene		1,000.6	µg/mL	+/-	5.8176	µg/mL	Gravimetric
	CAS # 193-39-5	(Lot ER082107-02)			+/-	10.9553	µg/mL	Unstressed
	Purity 99%				+/-	18.5819	µg/mL	Stressed
79	Dibenz(a,h)anthracene		1,004.4	µg/mL	+/-	5.8397	µg/mL	Gravimetric
	CAS # 53-70-3	(Lot ER032211-01)			+/-	10.9969	µg/mL	Unstressed
	Purity 99%				+/-	18.6525	µg/mL	Stressed
80	Benzo(g,h,i)perylene		1,000.0	µg/mL	+/-	5.8141	µg/mL	Gravimetric
	CAS # 191-24-2	(Lot ER020708-08)			+/-	10.9487	µg/mL	Unstressed
	Purity 99%				+/-	18.5708	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Specific Reference Material Notes:

N-nitrosodiphenylamine 2000 ug/mL equivalent when used for GC analysis. Actual formulation is diphenylamine 1710 ug/mL. N-Nitrosodiphenylamine is prone to breakdown in the injection port and will be converted to diphenylamine. N-Nitrosodiphenylamine is also a reactive species that can initiate premature decomposition of other compounds in the mix. For these reasons diphenylamine is used in the preparation of this mixture. When comparing the response of this compound to mixtures manufactured using N-nitrosodiphenylamine, a difference in response will be observed.

Column:
30m x 0.25mm x 0.25µm
Rtx-5 (cat.#10223)

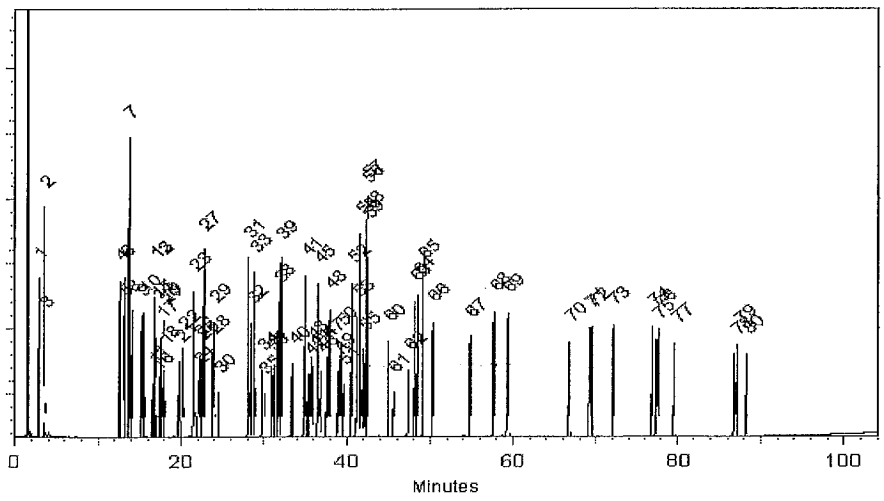
Carrier Gas:
hydrogen-constant pressure 10 psi

Temp. Program:
35°C (hold 3 min.) to 330°C
@ 3°C/min. (hold 3 min.)

Inj. Temp:
250°C

Det. Temp:
300°C

Det. Type:
FID



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

F. Joseph Tallon - Mix Technician **Date Mixed:** 24-Nov-2014 **Balance:** 1128360905

Jodi E. Breon - QA Analyst **Date Passed:** 05-Dec-2014

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

k is a coverage factor of 2, which gives a level of confidence of approximately 95%.

- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

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Reference Material Producer
Certificate #3222.01

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Testing Laboratory
Certificate #3222.02

FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 569730 **Lot No.:** A0108709

Description : 8270 List 1 / Std #9

8270 List 1 / Std #9 2,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 10 mL **Pkg Amt:** > 5 mL

Expiration Date : July 31, 2016 **Storage:** 10°C or colder

Handling: Contains carcinogen/reproductive toxin.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Benzidine **CAS #** 92-87-5 **Purity** 99% (Lot 141208JLM)	2,006.6 µg/mL	+/- 11.6665 µg/mL +/- 21.9697 µg/mL +/- 37.2641 µg/mL		Gravimetric Unstressed Stressed
2	3,3'-Dichlorobenzidine **CAS #** 91-94-1 **Purity** 99% (Lot 141205JLM)	2,001.0 µg/mL	+/- 11.6340 µg/mL +/- 21.9083 µg/mL +/- 37.1601 µg/mL		Gravimetric Unstressed Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

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FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 569731 **Lot No.:** A0107943

Description : 8270 List 1 / Std #10

8270 List 1 / Std #10 2,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : June 30, 2016 **Storage:** 10°C or colder

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)	
1	Indene	2,001.4 µg/mL	+/- 11.6363 µg/mL	Gravimetric
	CAS # 95-13-6 (Lot MKBP3098V)		+/- 22.5687 µg/mL	Unstressed
	Purity 99%		+/- 25.9700 µg/mL	Stressed
2	Benzoic acid	2,005.8 µg/mL	+/- 11.6619 µg/mL	Gravimetric
	CAS # 65-85-0 (Lot MKBL6689V)		+/- 22.6183 µg/mL	Unstressed
	Purity 99%		+/- 26.0271 µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Column:
30m x 0.25mm x 0.25µm
Rtx-5 (cat.#10223)

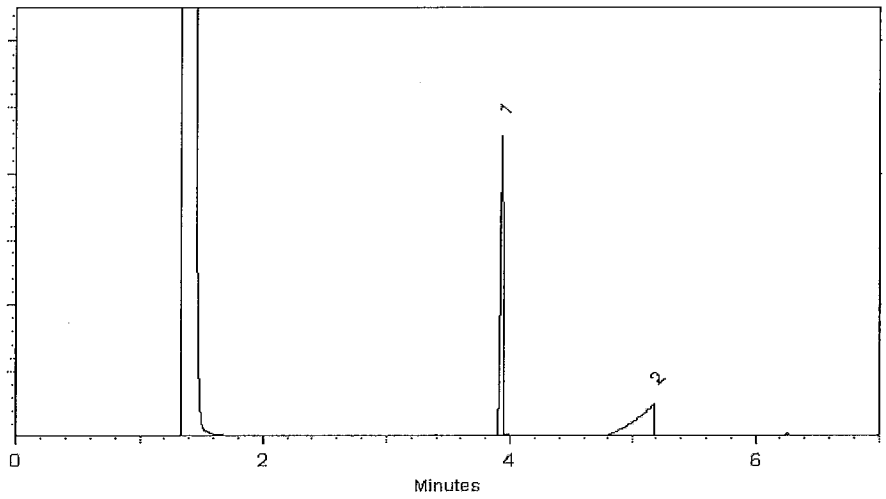
Carrier Gas:
hydrogen-constant pressure 10 psi.

Temp. Program:
75°C (hold 1 min.) to 330°C
@ 20°C/min. (hold 10 min.)

Inj. Temp:
250°C

Det. Temp:
330°C

Det. Type:
FID



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

F. Joseph Fallon - Mix Technician

Date Mixed: 22-Dec-2014 **Balance:** 1128360905

Tyler Brown - QA Analyst

Date Passed: 29-Dec-2014

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

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Testing Laboratory
Certificate #3222.02

FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 569731 **Lot No.:** A0107943

Description : 8270 List 1 / Std #10

8270 List 1 / Std #10 2,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : June 30, 2016 **Storage:** 10°C or colder

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)	
1	Indene	2,001.4 µg/mL	+/- 11.6363 µg/mL	Gravimetric
	CAS # 95-13-6 (Lot MKBP3098V)		+/- 22.5687 µg/mL	Unstressed
	Purity 99%		+/- 25.9700 µg/mL	Stressed
2	Benzoic acid	2,005.8 µg/mL	+/- 11.6619 µg/mL	Gravimetric
	CAS # 65-85-0 (Lot MKBL6689V)		+/- 22.6183 µg/mL	Unstressed
	Purity 99%		+/- 26.0271 µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Column:
30m x 0.25mm x 0.25µm
Rtx-5 (cat.#10223)

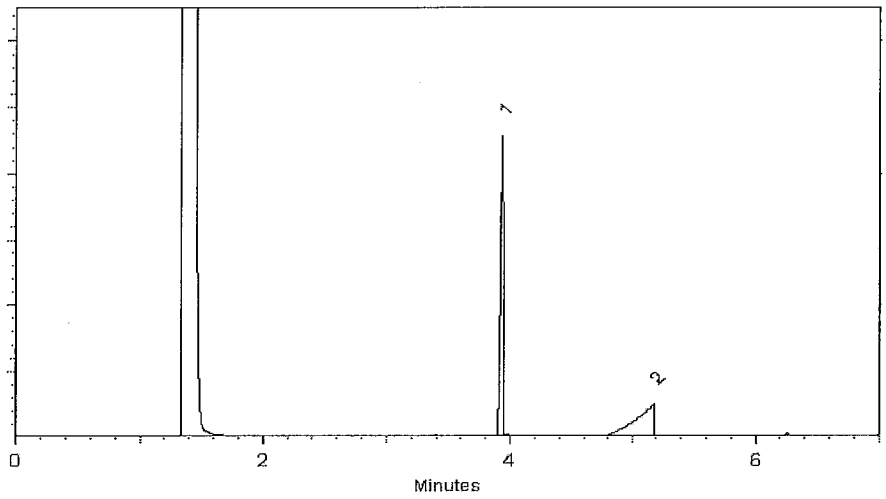
Carrier Gas:
hydrogen-constant pressure 10 psi.

Temp. Program:
75°C (hold 1 min.) to 330°C
@ 20°C/min. (hold 10 min.)

Inj. Temp:
250°C

Det. Temp:
330°C

Det. Type:
FID



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

F. Joseph Fallon - Mix Technician

Date Mixed: 22-Dec-2014 **Balance:** 1128360905

Tyler Brown - QA Analyst

Date Passed: 29-Dec-2014

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

RESTEK®
110 Benner Circle
Bellefonte, PA 16823-8812
Tel: (800)356-1688
Fax: (814)353-1309

www.restek.com

CERTIFIED REFERENCE MATERIAL

ACCREDITED
ISO Guide 34 Accredited
Reference Material Producer
Certificate #3222.01

Certificate of Analysis

ilac-MRA

ACCREDITED
ISO/IEC 17025 Accredited
Testing Laboratory
Certificate #3222.02

FOR LABORATORY USE ONLY-READ SDS PRIOR TO USE.

This Reference Material is intended for Laboratory Use Only as a standard for the qualitative and/or quantitative determination of the analyte(s) listed.

Catalog No. : 569732 **Lot No.:** A0108035

Description : 8270 List 1 / Std #11

8270 List 1 / Std #11 2,000 ug/ml, Methylene Chloride, 5 ml/ampul

Container Size : 5 mL **Pkg Amt:** > 5 mL

Expiration Date : June 30, 2016 **Storage:** 10°C or colder

Handling: This product is photosensitive.

CERTIFIED VALUES

Elution Order	Compound	Grav. Conc. (weight/volume)	Expanded Uncertainty (95% C.L.; K=2)		
1	Benzaldehyde	2,000.6 µg/mL	+/- 11.6317	µg/mL	Gravimetric
	CAS # 100-52-7 (Lot SHBD3510V)		+/- 64.1305	µg/mL	Unstressed
	Purity 99%		+/- 74.5493	µg/mL	Stressed
2	epsilon-Caprolactam	2,001.2 µg/mL	+/- 11.6351	µg/mL	Gravimetric
	CAS # 105-60-2 (Lot I16X016)		+/- 64.1498	µg/mL	Unstressed
	Purity 99%		+/- 74.5716	µg/mL	Stressed
3	Atrazine	2,004.3 µg/mL	+/- 11.6532	µg/mL	Gravimetric
	CAS # 1912-24-9 (Lot TZ8ED)		+/- 64.2490	µg/mL	Unstressed
	Purity 98%		+/- 74.6870	µg/mL	Stressed

Solvent: Methylene Chloride
CAS # 75-09-2
Purity 99%

Column:
30m x 0.25mm x 0.25µm
Rtx-5 (cat.#10223)

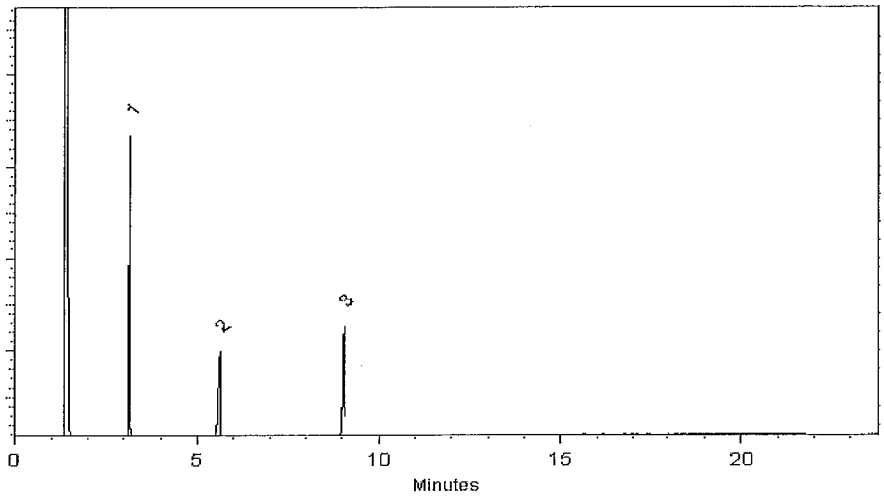
Carrier Gas:
hydrogen-constant pressure 10 psi.

Temp. Program:
75°C (hold 1 min.) to 330°C
@ 20°C/min. (hold 10 min.)

Inj. Temp:
250°C

Det. Temp:
330°C

Det. Type:
FID



This chromatogram represents a general set of testing conditions chosen for product acceptance. For optimal results in your lab, conditions should be adjusted for your specific instrument, method, and application.

F. Joseph Tallon - Mix Technician

Date Mixed: 30-Dec-2014 **Balance:** 1128360905

Jennifer L. Pollino - QC Analyst

Date Passed: 02-Jan-2015

Manufactured under Restek's ISO 9001:2008 Registered Quality System Certificate #FM 80397

General Certified Reference Material Notes

Expiration Notes:

- Expiration date valid for unopened ampul stored in compliance with the recommended conditions.
- Uncertainty, concentration, and expiration of the CRM are based on the unopened product being stored according to the recommended condition found in the storage field.

Purity Notes:

- Purity and/or chemical identity are determined by one or more of the following techniques: GC/FID, HPLC, GC/μECD, GC/MS, LC/MS, RI, and/or melting point.
- Compounds with a listed purity of less than 99% have been weight corrected to compensate for impurities and/or salts. A correction factor is used to calculate the amount of compound necessary to achieve the desired concentration of the parent compound in solution.
- Purity of isomeric compounds is reported as the sum of the isomers.
- Purity values are rounded to the nearest whole number.

Certified Uncertainty Value Notes:

- The uncertainties are determined in accordance with ISO Guides 34 and 35. The certified combined stressed uncertainty value (includes gravimetric uncertainty, homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty and were combined using the following formula:

$$U_{combined\ stressed} = k\sqrt{U^2_{gravimetric} + U^2_{homogeneity} + U^2_{storage\ stability} + U^2_{shipping\ stability}}$$

  k is a coverage factor of 2, which gives a level of confidence of approximately 95%.
- It is important to note that the shipping stability uncertainty was obtained under temperature extremes for specific time intervals; therefore, the certified combined stressed uncertainty value should only be applied to the product if it was stored at non-standard temperature conditions up to and including 7 days. Contact Restek Technical Service at www.restek.com/Contact-Us for use recommendations if your shipment was in-transit for more than 7 days at non-standard temperature conditions.
- Apply the certified combined unstressed uncertainty value if the product was received under standard shipping conditions. Apply the certified combined stressed uncertainty value if the product was received under non-standard conditions as specified below.

Label Conditions	Standard Conditions	Non-Standard Conditions
25°C Nominal (Room Temperature)	< 60°C	≥ 60°C up to 7 days
10°C or colder (Refrigerate)	< 40°C	≥ 40°C up to 7 days
0°C or colder (Freezer)	< 25°C	≥ 25°C up to 7 days

- Separate (not combined) uncertainty values for gravimetric uncertainty are also displayed on the certificate, if needed, separate homogeneity between-ampul uncertainty, storage stability uncertainty and shipping stability uncertainty values are available by contacting Restek Technical Service at www.restek.com/Contact-Us.
- The packaged amount is the minimum sample size for which uncertainty is valid. The ampules are over-filled to ensure that the minimum packaged amount can be sufficiently transferred.

Manufacturing Notes:

- Concentration is based upon gravimetric preparation using either a balance whose calibration has been verified daily using NIST traceable weights, and/or dilutions with Class A glassware.

Handling Notes:

- Samples should be transferred into deactivated vials for handling and storage. Restek supplies deactivated vials along with most standards packed in 2 mL ampules. Due to space constraints, Restek does not supply vials for larger volume ampules. Restek sells DMDCS for the purpose of glassware deactivation as catalog number 31861, which includes complete instructions. Restek will also deactivate larger volume vials from our inventory as a custom ordered item. Contact your Restek sales or customer service representative for details.
- If any undissolved material is visible inside the ampul, sonicate the unopened ampul until the material is completely dissolved.

Certification Summary

Client: Sundance Consulting, Inc
Project/Site: Fort Wingate, New Mexico

TestAmerica Job ID: 280-67711-2

Laboratory	Authority	Program	EPA Region	Certification ID
TestAmerica Denver	A2LA	DoD ELAP		2907.01
TestAmerica Denver	A2LA	ISO/IEC 17025		2907.01
TestAmerica Denver	Alaska (UST)	State Program	10	UST-30
TestAmerica Denver	Arizona	State Program	9	AZ0713
TestAmerica Denver	Arkansas DEQ	State Program	6	88-0687
TestAmerica Denver	California	State Program	9	2513
TestAmerica Denver	Connecticut	State Program	1	PH-0686
TestAmerica Denver	Florida	NELAP	4	E87667
TestAmerica Denver	Georgia	State Program	4	N/A
TestAmerica Denver	Illinois	NELAP	5	200017
TestAmerica Denver	Iowa	State Program	7	370
TestAmerica Denver	Kansas	NELAP	7	E-10166
TestAmerica Denver	Louisiana	NELAP	6	02096
TestAmerica Denver	Maine	State Program	1	CO0002
TestAmerica Denver	Minnesota	NELAP	5	8-999-405
TestAmerica Denver	Nevada	State Program	9	CO0026
TestAmerica Denver	New Jersey	NELAP	2	CO004
TestAmerica Denver	New York	NELAP	2	11964
TestAmerica Denver	North Carolina (WW/SW)	State Program	4	358
TestAmerica Denver	North Dakota	State Program	8	R-034
TestAmerica Denver	Oklahoma	State Program	6	8614
TestAmerica Denver	Oregon	NELAP	10	4025
TestAmerica Denver	Pennsylvania	NELAP	3	68-00664
TestAmerica Denver	South Carolina	State Program	4	72002001
TestAmerica Denver	Texas	NELAP	6	T104704183-13-8
TestAmerica Denver	USDA	Federal		P330-13-00202
TestAmerica Denver	Utah	NELAP	8	CO00026
TestAmerica Denver	Virginia	NELAP	3	460232
TestAmerica Denver	Washington	State Program	10	C583
TestAmerica Denver	West Virginia DEP	State Program	3	354
TestAmerica Denver	Wisconsin	State Program	5	999615430
TestAmerica Denver	Wyoming (UST)	A2LA	8	2907.01

Accreditation may not be offered or required for all methods and analytes reported in this package. Please contact your project manager for the laboratory's current list of certified methods and analytes.

Method 8270D

Semivolatile Organic Compounds (GC/MS) by Method 8270D

FORM II
GC/MS SEMI VOA SURROGATE RECOVERY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Matrix: Water Level: Low

GC Column (1): Vf-5MS (30.2 ID: 0.25 (mm)

Client Sample ID	Lab Sample ID	2FP #	PHL #	NBZ #	FBP #	TBP #	TPH #
TMW43042015	280-67711-1	65	66	66	64	67	65
DTW43042015	280-67711-2	65	66	66	66	67	64
TMW47042015	280-67711-4	51	53	53	53	60	63
	MB 280-272870/1-A	64	64	65	62	65	64
	LCS 280-272870/2-A	60	60	61	61	68	62
TMW43042015 MS	280-67711-1 MS	58	60	58	61	72	65
TMW43042015 MSD	280-67711-1 MSD	62	62	63	61	66	62

	QC LIMITS
2FP = 2-Fluorophenol (Surr)	41-135
PHL = Phenol-d5 (Surr)	46-135
NBZ = Nitrobenzene-d5 (Surr)	42-135
FBP = 2-Fluorobiphenyl	48-135
TBP = 2,4,6-Tribromophenol (Surr)	48-135
TPH = Terphenyl-d14 (Surr)	20-135

Column to be used to flag recovery values

FORM II 8270D

FORM III
GC/MS SEMI VOA LAB CONTROL SAMPLE RECOVERY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Matrix: Water Level: Low Lab File ID: G6_17421.D

Lab ID: LCS 280-272870/2-A Client ID:

COMPOUND	SPIKE ADDED (ug/L)	LCS CONCENTRATION (ug/L)	LCS % REC	QC LIMITS REC	#
Caprolactam	80.0	52.3	65	64-120	

Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE RECOVERY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Matrix: Water Level: Low Lab File ID: G6_17429.D

Lab ID: 280-67711-1 MS Client ID: TMW43042015 MS

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC	QC LIMITS REC	#
Caprolactam	90.4	2.7 U	62.1	69	64-120	

Column to be used to flag recovery and RPD values

FORM III 8270D

FORM III
GC/MS SEMI VOA MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Matrix: Water Level: Low Lab File ID: G6_17430.D

Lab ID: 280-67711-1 MSD Client ID: TMW43042015 MSD

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD % REC	% RPD	QC LIMITS RPD	QC LIMITS REC	#
Caprolactam	88.6	58.4	66	6	30	64-120	

Column to be used to flag recovery and RPD values

FORM III 8270D

FORM IV
GC/MS SEMI VOA METHOD BLANK SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab File ID: G6_17420.D Lab Sample ID: MB 280-272870/1-A

Matrix: Water Date Extracted: 04/15/2015 18:50

Instrument ID: SMS_G6 Date Analyzed: 04/18/2015 17:49

Level:(Low/Med) Low

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
	LCS 280-272870/2-A	G6_17421.D	04/18/2015 18:16
TMW43042015	280-67711-1	G6_17428.D	04/18/2015 21:13
TMW43042015 MS	280-67711-1 MS	G6_17429.D	04/18/2015 21:39
TMW43042015 MSD	280-67711-1 MSD	G6_17430.D	04/18/2015 22:06
DTW43042015	280-67711-2	G6_17431.D	04/18/2015 22:32
TMW47042015	280-67711-4	G6_17432.D	04/18/2015 22:59

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab File ID: G6_16636.D DFTPP Injection Date: 02/25/2015

Instrument ID: SMS_G6 DFTPP Injection Time: 11:44

Analysis Batch No.: 272059

M/E	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE	
51	10-80% of Base Peak	49.7	
68	Less than 2% of mass 69	0.7	(1.5)1
69	Mass 69 Relative abundance	47.6	
70	Less than 2% of mass 69	0.1	(0.2)1
127	10-80% of Base Peak	56.0	
197	Less than 2% of mass 198	0.0	
198	Base peak	100.0	
199	5-9% of mass 198	6.6	
275	10-60% of Base Peak	19.4	
365	Greater than 1% of mass 198	2.1	
441	present but less than 24% of mass 442	8.3	(14.8)2
442	Greater than 50% of mass 198	56.0	
443	15-24% of mass 442	10.7	(19.0)2

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	ICIS 280-272059/3	G6_16637.D	02/25/2015	11:53
	STD004 280-272059/4	G6_16638.D	02/25/2015	12:19
	STD010 280-272059/5	G6_16639.D	02/25/2015	12:46
	STD020 280-272059/6	G6_16640.D	02/25/2015	13:12
	STD050 280-272059/7	G6_16641.D	02/25/2015	13:39
	STD120 280-272059/8	G6_16642.D	02/25/2015	14:06
	STD160 280-272059/9	G6_16643.D	02/25/2015	14:32
	STD200 280-272059/10	G6_16644.D	02/25/2015	14:59
	ICV 280-272059/11	G6_16645.D	02/25/2015	15:26
	ICV 280-272059/12	G6_16646.D	02/25/2015	15:52
	ICV 280-272059/13	G6_16647.D	02/25/2015	16:19

FORM V
GC/MS SEMI VOA INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab File ID: G6_17413.D DFTPP Injection Date: 04/18/2015

Instrument ID: SMS_G6 DFTPP Injection Time: 15:09

Analysis Batch No.: 273380

M/E	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE	
51	10-80% of Base Peak	49.1	
68	Less than 2% of mass 69	0.6	(1.2)1
69	Mass 69 Relative abundance	48.3	
70	Less than 2% of mass 69	0.3	(0.6)1
127	10-80% of Base Peak	56.1	
197	Less than 2% of mass 198	0.1	
198	Base peak	100.0	
199	5-9% of mass 198	7.0	
275	10-60% of Base Peak	17.6	
365	Greater than 1% of mass 198	2.0	
441	present but less than 24% of mass 442	7.1	(14.8)2
442	Greater than 50% of mass 198	48.2	
443	15-24% of mass 442	9.4	(19.5)2

1-Value is % mass 69 2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
	CCV 280-273380/3	G6_17414.D	04/18/2015	15:20
	CCV 280-273380/4	G6_17416.D	04/18/2015	16:09
	MB 280-272870/1-A	G6_17420.D	04/18/2015	17:49
	LCS 280-272870/2-A	G6_17421.D	04/18/2015	18:16
TMW43042015	280-67711-1	G6_17428.D	04/18/2015	21:13
TMW43042015 MS	280-67711-1 MS	G6_17429.D	04/18/2015	21:39
TMW43042015 MSD	280-67711-1 MSD	G6_17430.D	04/18/2015	22:06
DTW43042015	280-67711-2	G6_17431.D	04/18/2015	22:32
TMW47042015	280-67711-4	G6_17432.D	04/18/2015	22:59

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Sample No.: ICIS 280-272059/3 Date Analyzed: 02/25/2015 11:53

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.25) ID: 0.25(mm)

Lab File ID (Standard): G6_16637.D Heated Purge: (Y/N) N

Calibration ID: 21947

		DCB		NPT		ANT	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
INITIAL CALIBRATION MID-POINT		187021	4.80	713420	6.02	385413	7.78
UPPER LIMIT		374042	5.30	1426840	6.52	770826	8.28
LOWER LIMIT		93511	4.30	356710	5.52	192707	7.28
LAB SAMPLE ID	CLIENT SAMPLE ID						
ICV 280-272059/11		188037	4.80	718467	6.02	393887	7.78
ICV 280-272059/12		185167	4.80	715813	6.02	400054	7.78
ICV 280-272059/13		186013	4.80	710514	6.02	389552	7.78
CCV 280-273380/3		216765	4.71	822108	5.95	444568	7.71
CCV 280-273380/4		196797	4.71	720823	5.95	395890	7.70
MB 280-272870/1-A		217388	4.71	813022	5.95	452207	7.70
LCS 280-272870/2-A		220573	4.71	836102	5.95	460509	7.70
280-67711-1	TMW43042015	206375	4.71	800081	5.95	441344	7.70
280-67711-1 MS	TMW43042015 MS	207648	4.71	815083	5.95	448804	7.70
280-67711-1 MSD	TMW43042015 MSD	222633	4.71	860621	5.95	488251	7.70
280-67711-2	DTW43042015	212896	4.71	809424	5.95	443429	7.70
280-67711-4	TMW47042015	201871	4.71	767447	5.95	436451	7.70

DCB = 1,4-Dichlorobenzene-d4 (IS)
NPT = Naphthalene-d8 (IS)
ANT = Acenaphthene-d10 (IS)

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

Column used to flag values outside QC limits

FORM VIII 8270D

FORM VIII
GC/MS SEMI VOA INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Sample No.: ICIS 280-272059/3 Date Analyzed: 02/25/2015 11:53

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.25) ID: 0.25(mm)

Lab File ID (Standard): G6_16637.D Heated Purge: (Y/N) N

Calibration ID: 21947

		PHN		CRY		PRY	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
INITIAL CALIBRATION MID-POINT		653325	9.28	616068	13.71	532020	17.90
UPPER LIMIT		1306650	9.78	1232136	14.21	1064040	18.40
LOWER LIMIT		326663	8.78	308034	13.21	266010	17.40
LAB SAMPLE ID	CLIENT SAMPLE ID						
ICV 280-272059/11		663955	9.29	625452	13.71	534138	17.91
ICV 280-272059/12		674469	9.28	649859	13.70	546537	17.90
ICV 280-272059/13		665613	9.28	633484	13.70	533431	17.90
CCV 280-273380/3		728379	9.21	645212	13.53	551738	17.66
CCV 280-273380/4		665198	9.21	600975	13.52	517824	17.65
MB 280-272870/1-A		769713	9.20	727965	13.51	630046	17.65
LCS 280-272870/2-A		780849	9.20	716077	13.52	613809	17.65
280-67711-1	TMW43042015	750086	9.20	707015	13.52	606974	17.65
280-67711-1 MS	TMW43042015 MS	777208	9.20	709191	13.52	608383	17.66
280-67711-1 MSD	TMW43042015 MSD	823179	9.20	756154	13.52	637043	17.65
280-67711-2	DTW43042015	757081	9.20	713388	13.51	605197	17.65
280-67711-4	TMW47042015	746157	9.20	684104	13.51	593758	17.65

PHN = Phenanthrene-d10 (IS)
CRY = Chrysene-d12 (IS)
PRY = Perylene-d12 (IS)

Area Limit = 50%-200% of internal standard area
RT Limit = ± 0.5 minutes of internal standard RT

Column used to flag values outside QC limits

FORM VIII 8270D

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Denver | Job No.: 280-67711-2

SDG No.:

Client Sample ID: TMW43042015 | Lab Sample ID: 280-67711-1

Matrix: Water | Lab File ID: G6_17428.D

Analysis Method: 8270D | Date Collected: 04/10/2015 08:25

Extract. Method: 3520C | Date Extracted: 04/15/2015 18:50

Sample wt/vol: 925(mL) | Date Analyzed: 04/18/2015 21:13

Con. Extract Vol.: 1(mL) | Dilution Factor: 1

Injection Volume: 0.5(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 273380 | Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
105-60-2	Caprolactam	2.7	U	5.4	2.7

CAS NO.	SURROGATE	%REC	Q	LIMITS
118-79-6	2,4,6-Tribromophenol (Surr)	67		48-135
321-60-8	2-Fluorobiphenyl	64		48-135
367-12-4	2-Fluorophenol (Surr)	65		41-135
4165-60-0	Nitrobenzene-d5 (Surr)	66		42-135
4165-62-2	Phenol-d5 (Surr)	66		46-135
1718-51-0	Terphenyl-d14 (Surr)	65		20-135

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17428.D
Lims ID: 280-67711-C-1-A Lab Sample ID: 280-67711-1
Client ID: TMW43042015
Sample Type: Client
Inject. Date: 18-Apr-2015 21:13:30 ALS Bottle#: 16 Worklist Smp#: 29
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: 280-67711-C-1-B
Operator ID: HOEFLERA Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:13 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: kiekeld Date: 20-Apr-2015 07:06:02

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	96	206375	40.0	
* 2 Naphthalene-d8	136	5.945	5.951	-0.006	100	800081	40.0	
* 3 Acenaphthene-d10	164	7.704	7.704	0.000	92	441344	40.0	
* 4 Phenanthrene-d10	188	9.204	9.210	-0.006	97	750086	40.0	
* 5 Chrysene-d12	240	13.516	13.516	0.000	96	707015	40.0	
* 6 Perylene-d12	264	17.651	17.651	0.000	95	606974	40.0	
$ 7 2-Fluorophenol	112	3.487	3.493	-0.006	92	488759	65.1	
$ 8 Phenol-d5	99	4.328	4.334	-0.006	99	621902	66.1	
$ 9 Nitrobenzene-d5	82	5.240	5.245	-0.005	90	541909	66.1	
$ 10 2-Fluorobiphenyl	172	6.998	7.004	-0.006	99	954256	64.0	
$ 11 2,4,6-Tribromophenol	330	8.498	8.504	-0.006	87	110712	66.7	
$ 12 Terphenyl-d14	244	11.233	11.239	-0.006	99	911263	64.9	
13 1,4-Dioxane	88		2.040				ND	
14 N-Nitrosodimethylamine	74		2.305				ND	
15 Pyridine	79		2.352				ND	
16 2-Picoline	93		2.987				ND	
17 N-Nitrosomethylethylamine	88		3.081				ND	
18 Methyl methanesulfonate	80		3.351				ND	
19 N-Nitrosodiethylamine	102		3.704				ND	
20 Pentachlorophenol_T	266		3.822				ND	
21 Ethyl methanesulfonate	79		3.969				ND	
22 Pentachloroethane	117		4.446				ND	
23 Phenol	94		4.346				ND	
24 Aniline	93		4.404				ND	
25 Bis(2-chloroethyl)ether	93		4.434				ND	
26 2-Chlorophenol	128		4.516				ND	
27 1,3-Dichlorobenzene	146		4.657				ND	
28 1,4-Dichlorobenzene	146		4.728				ND	
29 Benzyl alcohol	108		4.828				ND	
30 1,2-Dichlorobenzene	146		4.875				ND	
31 2-Methylphenol	108		4.916				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/ml	Flags
32 2,2'-oxybis[1-chloropropan	45		4.934				ND	
33 N-Nitrosopyrrolidine	100		5.075				ND	
34 N-Nitrosomorpholine	116		5.104				ND	
35 Benzidine_T	184		4.952				ND	
36 2-Toluidine	106		5.122				ND	
37 Benzaldehyde	106		5.015				ND	
38 3 & 4 Methylphenol	108		5.063				ND	
39 3-Methylphenol	108		5.063				ND	
40 4-Methylphenol	108		5.063				ND	
41 N-Nitrosodi-n-propylamine	70		5.069				ND	
42 Acetophenone	105		5.087				ND	
43 Hexachloroethane	117		5.198				ND	
44 Nitrobenzene	77		5.263				ND	
45 N-Nitrosopiperidine	114		5.410				ND	
46 Isophorone	82		5.475				ND	
47 o,o',o''-Triethylphosphoro	198		5.622				ND	
48 2-Nitrophenol	139		5.569				ND	
49 2,4-Dimethylphenol	107		5.581				ND	
50 Bis(2-chloroethoxy)methane	93		5.669				ND	
51 alpha,alpha-Dimethyl phene	58		5.792				ND	
52 Benzoic acid	105		5.687				ND	
53 2,4-Dichlorophenol	162		5.798				ND	
54 1,2,4-Trichlorobenzene	180		5.887				ND	
55 2,6-Dichlorophenol	162		6.022				ND	
56 Hexachloropropene	213		6.045				ND	
57 Naphthalene	128		5.975				ND	
58 4-Chloroaniline	127		6.016				ND	
59 Hexachlorobutadiene	225		6.069				ND	
60 N-Nitrosodi-n-butylamine	84		6.310				ND	
61 p-Phenylene diamine	108		6.363				ND	
62 Caprolactam	55		6.369				ND	
63 Safrole, Total	162		6.540				ND	
64 4-Chloro-3-methylphenol	107		6.481				ND	
65 2-Methylnaphthalene	142		6.651				ND	
66 Isosafrole Peak 1	162		6.834				ND	
67 1-Methylnaphthalene	142		6.757				ND	
68 Hexachlorocyclopentadiene	237		6.798				ND	
69 1,2,4,5-Tetrachlorobenzene	216		6.822				ND	
70 2,4,6-Trichlorophenol	196		6.934				ND	
71 Isosafrole Peak 2	104		7.063				ND	
72 2,4,5-Trichlorophenol	196		6.969				ND	
73 1-Chloronaphthalene	162		7.169				ND	
74 1,1'-Biphenyl	154		7.110				ND	
75 2-Chloronaphthalene	162		7.151				ND	
76 1,4-Naphthoquinone	158		7.328				ND	
77 2-Nitroaniline	65		7.251				ND	
78 1,4-Dinitrobenzene	168		7.381				ND	
79 Dimethyl phthalate	163		7.398				ND	
80 1,3-Dinitrobenzene	168		7.463				ND	
81 2,6-Dinitrotoluene	165		7.481				ND	
82 Acenaphthylene	152		7.569				ND	
83 3-Nitroaniline	138		7.663				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/ml	Flags
84 Acenaphthene	153		7.739				ND	
85 Pentachlorobenzene	250		7.863				ND	
86 2,4-Dinitrophenol	184		7.763				ND	
87 4-Nitrophenol	109		7.816				ND	
88 1-Naphthylamine	143		7.998				ND	
89 2,4-Dinitrotoluene	165		7.887				ND	
90 Dibenzofuran	168		7.910				ND	
91 2-Naphthylamine	143		8.075				ND	
92 2,3,4,6-Tetrachlorophenol	232		8.028				ND	
93 Thionazin	97		8.169				ND	
94 Diethyl phthalate	149		8.092				ND	
95 N-Nitro-o-toluidine	152		8.263				ND	
96 4-Chlorophenyl phenyl ethe	204		8.234				ND	
97 Diphenylamine	169		8.351				ND	
98 Fluorene	166		8.257				ND	
99 4-Nitroaniline	138		8.286				ND	
100 4,6-Dinitro-2-methylphenol	198		8.298				ND	
101 Sulfotepp	97		8.451				ND	
102 N-Nitrosodiphenylamine	169		8.357				ND	
103 Azobenzene	77		8.392				ND	
104 1,2-Diphenylhydrazine	77		8.392				ND	
105 Diallate Peak 1	86		8.610				ND	
106 1,3,5-Trinitrobenzene	213		8.622				ND	
107 Phorate	121		8.628				ND	
108 Phenacetin	108		8.645				ND	
109 Diallate Peak 2	86		8.704				ND	
110 Dimethoate	87		8.822				ND	
111 4-Bromophenyl phenyl ether	248		8.728				ND	
112 Hexachlorobenzene	284		8.810				ND	
113 4-Aminobiphenyl	169		8.998				ND	
114 Pentachloronitrobenzene	237		9.010				ND	
115 Pronamide	173		9.016				ND	
116 Pentachlorophenol	266		9.010				ND	
117 Disulfoton	88		9.145				ND	
118 Dinoseb	211		9.157				ND	
119 Phenanthrene	178		9.233				ND	
120 Anthracene	178		9.286				ND	
121 Methyl parathion	109		9.551				ND	
122 Carbazole	167		9.445				ND	
123 Di-n-butyl phthalate	149		9.739				ND	
124 Ethyl Parathion	109		9.980				ND	
125 4-Nitroquinoline-1-oxide	190		10.104				ND	
126 Methapyrilene	97		10.128				ND	
127 Isodrin	193		10.433				ND	
128 Fluoranthene	202		10.675				ND	
129 Benzidine	184		10.466				ND	
130 Aramite Peak 1	185		11.139				ND	
131 Pyrene	202		11.039				ND	
132 Aramite Peak 2	185		11.274				ND	
133 p-Dimethylamino azobenzene	120		11.492				ND	
134 Chlorobenzilate	251		11.545				ND	
135 3,3'-Dimethylbenzidine	212		12.180				ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/ml	Flags
136 Famphur	218		12.045				ND	
137 Butyl benzyl phthalate	149		12.151				ND	
138 2-Acetylaminofluorene	181		12.745				ND	
139 4,4'-Methylene bis(2-chlor	231		13.463				ND	
140 3,3'-Dichlorobenzidine	252		13.468				ND	
141 Benzo[a]anthracene	228		13.504				ND	
142 Bis(2-ethylhexyl) phthalat	149		13.492				ND	
143 Chrysene	228		13.598				ND	
144 Di-n-octyl phthalate	149		15.362				ND	
145 7,12-Dimethylbenz(a)anthra	256		16.480				ND	
146 Benzo[b]fluoranthene	252		16.509				ND	
147 Benzo[k]fluoranthene	252		16.592				ND	
148 Benzo[a]pyrene	252		17.492				ND	
149 3-Methylcholanthrene	268		18.550				ND	
150 Dibenz[a,j]acridine	279		20.350				ND	
151 Indeno[1,2,3-cd]pyrene	276		20.862				ND	
152 Dibenz(a,h)anthracene	278		20.939				ND	
153 Benzo[g,h,i]perylene	276		21.633				ND	
S 160 Aramite, Total	185		15.047				ND	
S 161 Isosafrole	162		15.047				ND	
S 162 Diallate	86		15.047				ND	
154 Total Cresols	1		0.000				ND	
155 Tetraethyl Pyrophosphate (	1		0.000				ND	
157 4,4'-DDE	246		5.063				ND	
158 4,4'-DDD	235		5.375				ND	
159 4,4'-DDT	235		5.604				ND	
S 163 Methyl Phenols,Total	1		0.000				ND	

Reagents:

MS-IS_00007 Amount Added: 20.00 Units: uL Run Reagent

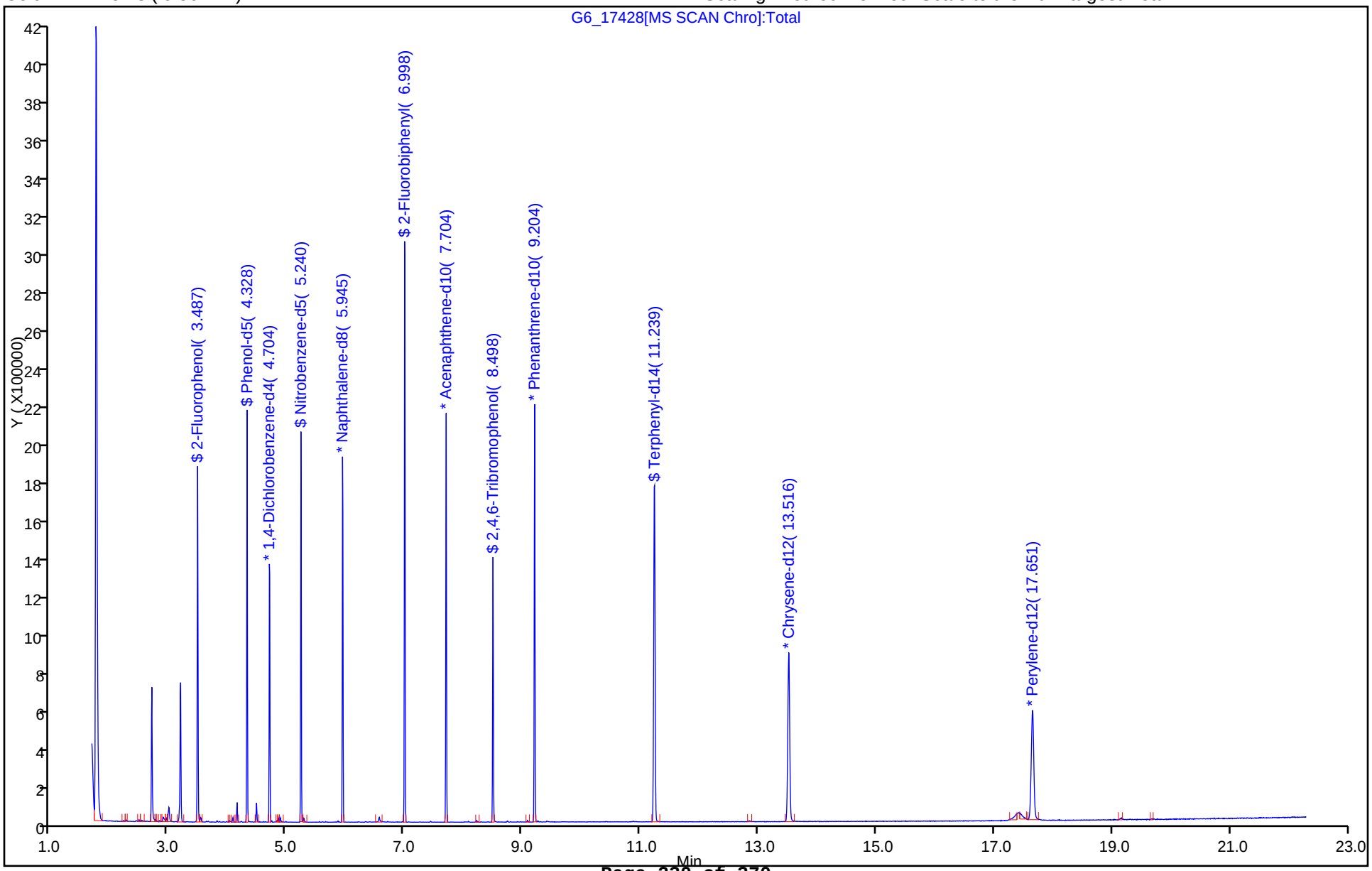
Report Date: 20-Apr-2015 07:09:15 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17428.D

Injection Date:	18-Apr-2015 21:13:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	280-67711-C-1-A	Lab Sample ID:	280-67711-1	Worklist Smp#:	29
Client ID:	TMW43042015				
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	16
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Denver | Job No.: 280-67711-2

SDG No.:

Client Sample ID: DTW43042015 | Lab Sample ID: 280-67711-2

Matrix: Water | Lab File ID: G6_17431.D

Analysis Method: 8270D | Date Collected: 04/10/2015 08:25

Extract. Method: 3520C | Date Extracted: 04/15/2015 18:50

Sample wt/vol: 1029(mL) | Date Analyzed: 04/18/2015 22:32

Con. Extract Vol.: 1(mL) | Dilution Factor: 1

Injection Volume: 0.5(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 273380 | Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
105-60-2	Caprolactam	2.4	U	4.9	2.4

CAS NO.	SURROGATE	%REC	Q	LIMITS
118-79-6	2,4,6-Tribromophenol (Surr)	67		48-135
321-60-8	2-Fluorobiphenyl	66		48-135
367-12-4	2-Fluorophenol (Surr)	65		41-135
4165-60-0	Nitrobenzene-d5 (Surr)	66		42-135
4165-62-2	Phenol-d5 (Surr)	66		46-135
1718-51-0	Terphenyl-d14 (Surr)	64		20-135

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17431.D
Lims ID: 280-67711-B-2-A Lab Sample ID: 280-67711-2
Client ID: DTW43042015
Sample Type: Client
Inject. Date: 18-Apr-2015 22:32:30 ALS Bottle#: 19 Worklist Smp#: 32
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: 280-67711-B-2-B
Operator ID: HOEFLERA Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:13 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: kiekeld Date: 20-Apr-2015 07:08:34

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	96	212896	40.0	
* 2 Naphthalene-d8	136	5.945	5.951	-0.006	100	809424	40.0	
* 3 Acenaphthene-d10	164	7.704	7.704	0.000	92	443429	40.0	
* 4 Phenanthrene-d10	188	9.204	9.210	-0.006	97	757081	40.0	
* 5 Chrysene-d12	240	13.510	13.516	-0.006	97	713388	40.0	
* 6 Perylene-d12	264	17.650	17.651	-0.001	96	605197	40.0	
$ 7 2-Fluorophenol	112	3.487	3.493	-0.006	92	503571	65.1	
$ 8 Phenol-d5	99	4.328	4.334	-0.006	99	640956	66.1	
$ 9 Nitrobenzene-d5	82	5.239	5.245	-0.006	90	543735	65.5	
$ 10 2-Fluorobiphenyl	172	6.998	7.004	-0.006	100	986894	65.8	
$ 11 2,4,6-Tribromophenol	330	8.498	8.504	-0.006	87	112309	67.3	
$ 12 Terphenyl-d14	244	11.233	11.239	-0.006	99	907066	64.0	
26 2-Chlorophenol	128		4.516				ND	
28 1,4-Dichlorobenzene	146		4.728				ND	
31 2-Methylphenol	108		4.916				ND	
41 N-Nitrosodi-n-propylamine	70		5.069				ND	
54 1,2,4-Trichlorobenzene	180		5.887				ND	
62 Caprolactam	55		6.369				ND	
64 4-Chloro-3-methylphenol	107		6.481				ND	
65 2-Methylnaphthalene	142		6.651				ND	
70 2,4,6-Trichlorophenol	196		6.934				ND	
84 Acenaphthene	153		7.739				ND	
87 4-Nitrophenol	109		7.816				ND	
89 2,4-Dinitrotoluene	165		7.887				ND	
116 Pentachlorophenol	266		9.010				ND	
120 Anthracene	178		9.286				ND	
131 Pyrene	202		11.039				ND	

Reagents:

MS-IS_00007 Amount Added: 20.00 Units: uL Run Reagent

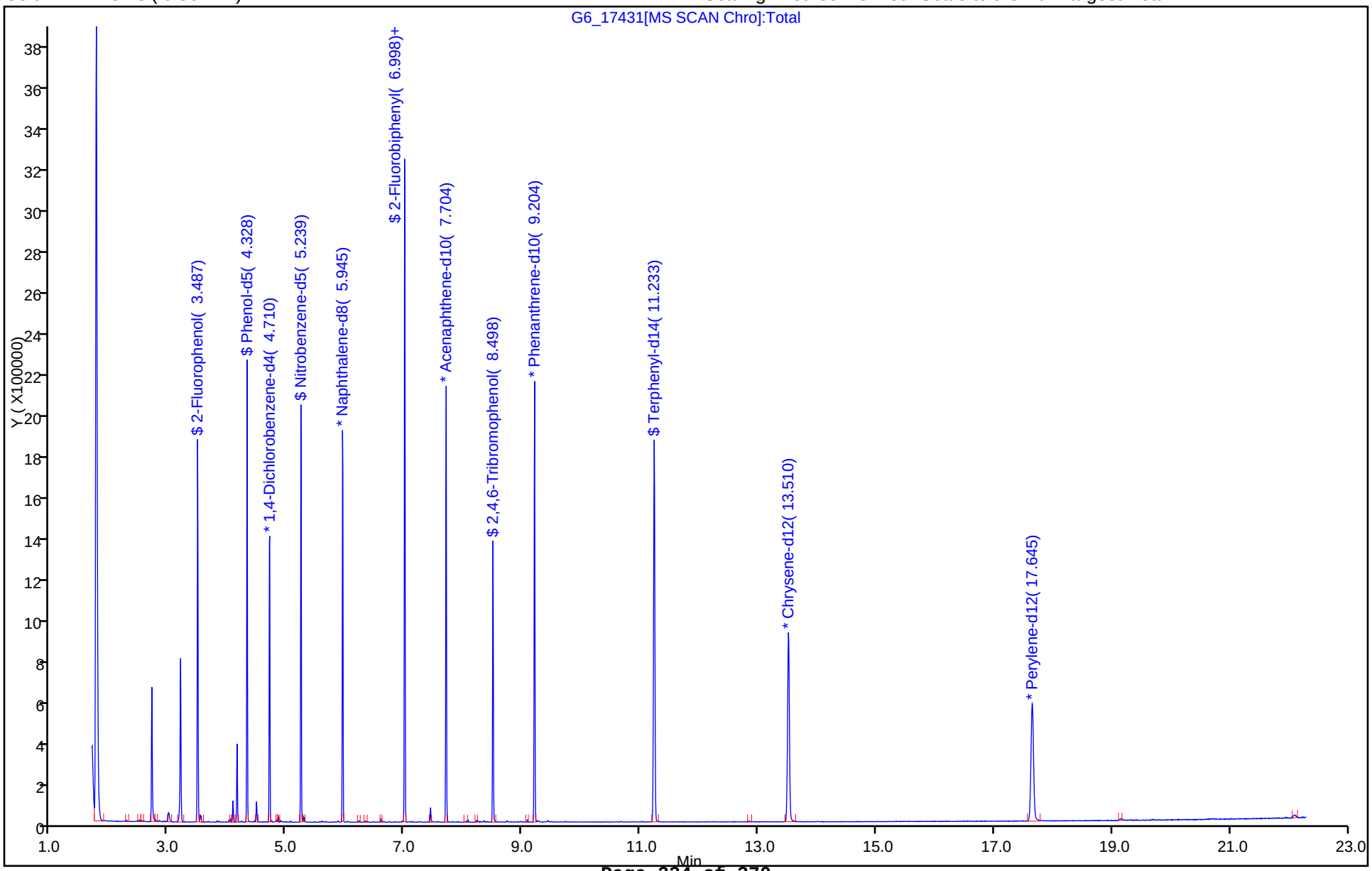
Report Date: 20-Apr-2015 07:09:17 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17431.D

Injection Date:	18-Apr-2015 22:32:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	280-67711-B-2-A	Lab Sample ID:	280-67711-2	Worklist Smp#:	32
Client ID:	DTW43042015				
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	19
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Denver | Job No.: 280-67711-2

SDG No.:

Client Sample ID: TMW47042015 | Lab Sample ID: 280-67711-4

Matrix: Water | Lab File ID: G6_17432.D

Analysis Method: 8270D | Date Collected: 04/10/2015 09:30

Extract. Method: 3520C | Date Extracted: 04/15/2015 18:50

Sample wt/vol: 963.2(mL) | Date Analyzed: 04/18/2015 22:59

Con. Extract Vol.: 1(mL) | Dilution Factor: 1

Injection Volume: 0.5(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 273380 | Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
105-60-2	Caprolactam	2.6	U	5.2	2.6

CAS NO.	SURROGATE	%REC	Q	LIMITS
118-79-6	2,4,6-Tribromophenol (Surr)	60		48-135
321-60-8	2-Fluorobiphenyl	53		48-135
367-12-4	2-Fluorophenol (Surr)	51		41-135
4165-60-0	Nitrobenzene-d5 (Surr)	53		42-135
4165-62-2	Phenol-d5 (Surr)	53		46-135
1718-51-0	Terphenyl-d14 (Surr)	63		20-135

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17432.D
Lims ID: 280-67711-C-4-A Lab Sample ID: 280-67711-4
Client ID: TMW47042015
Sample Type: Client
Inject. Date: 18-Apr-2015 22:59:30 ALS Bottle#: 20 Worklist Smp#: 33
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: 280-67711-C-4-B
Operator ID: HOEFLERA Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:13 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: kiekeld Date: 20-Apr-2015 07:08:45

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	96	201871	40.0	
* 2 Naphthalene-d8	136	5.945	5.951	-0.006	100	767447	40.0	
* 3 Acenaphthene-d10	164	7.704	7.704	0.000	92	436451	40.0	
* 4 Phenanthrene-d10	188	9.204	9.210	-0.006	97	746157	40.0	
* 5 Chrysene-d12	240	13.510	13.516	-0.006	96	684104	40.0	
* 6 Perylene-d12	264	17.645	17.651	-0.006	95	593758	40.0	
$ 7 2-Fluorophenol	112	3.487	3.493	-0.006	92	373127	50.8	
$ 8 Phenol-d5	99	4.328	4.334	-0.006	99	487221	53.0	
$ 9 Nitrobenzene-d5	82	5.240	5.245	-0.005	90	419917	53.4	
$ 10 2-Fluorobiphenyl	172	6.998	7.004	-0.006	100	778358	52.8	
$ 11 2,4,6-Tribromophenol	330	8.498	8.504	-0.006	88	98905	60.3	
$ 12 Terphenyl-d14	244	11.233	11.239	-0.006	99	859315	63.2	
26 2-Chlorophenol	128		4.516				ND	
28 1,4-Dichlorobenzene	146		4.728				ND	
31 2-Methylphenol	108		4.916				ND	
41 N-Nitrosodi-n-propylamine	70		5.069				ND	
54 1,2,4-Trichlorobenzene	180		5.887				ND	
62 Caprolactam	55		6.369				ND	
64 4-Chloro-3-methylphenol	107		6.481				ND	
65 2-Methylnaphthalene	142		6.651				ND	
70 2,4,6-Trichlorophenol	196		6.934				ND	
84 Acenaphthene	153		7.739				ND	
87 4-Nitrophenol	109		7.816				ND	
89 2,4-Dinitrotoluene	165		7.887				ND	
116 Pentachlorophenol	266		9.010				ND	
120 Anthracene	178		9.286				ND	
131 Pyrene	202		11.039				ND	

Reagents:

MS-IS_00007 Amount Added: 20.00 Units: uL Run Reagent

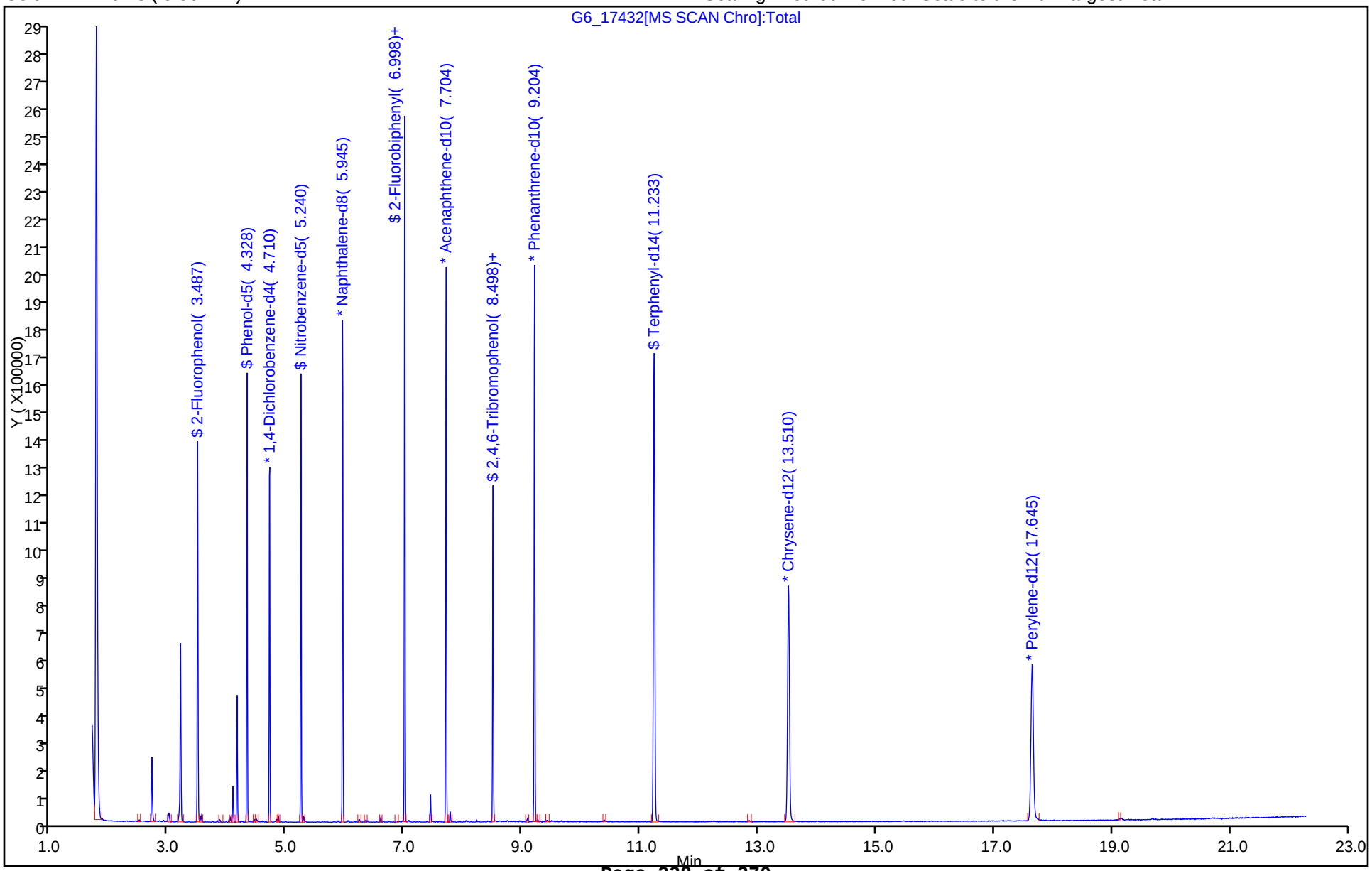
Report Date: 20-Apr-2015 07:09:18 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17432.D

Injection Date:	18-Apr-2015 22:59:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	280-67711-C-4-A	Lab Sample ID:	280-67711-4	Worklist Smp#:	33
Client ID:	TMW47042015				
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	20
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD004 280-272059/4	G6_16638.D
Level 2	STD010 280-272059/5	G6_16639.D
Level 3	STD020 280-272059/6	G6_16640.D
Level 4	STD050 280-272059/7	G6_16641.D
Level 5	ICIS 280-272059/3	G6_16637.D
Level 6	STD120 280-272059/8	G6_16642.D
Level 7	STD160 280-272059/9	G6_16643.D
Level 8	STD200 280-272059/10	G6_16644.D

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
1,4-Dioxane	0.6238 0.6039	0.6479 0.5901	0.6303 0.5885	0.6040	0.6208	Ave		0.6137				3.4		20.0			
N-Nitrosodimethylamine	0.9428 0.9384	0.9324 0.9251	0.9075 0.9268	0.9449	0.9390	Ave		0.9321				1.3		20.0			
Pyridine	1.6094 1.6115	1.5930 1.5754	1.5655 1.5629	1.5930	1.6057	Ave		1.5895				1.2		20.0			
Phenol	1.8894 1.8069	1.8459 1.7739	1.8225 1.7674	1.8507	1.8678	Ave		1.8281			0.8000	2.4		20.0			
Aniline	2.3768 2.2326	2.1918 2.2254	2.1905 2.1802	2.2465	2.2323	Ave		2.2345				2.8		20.0			
Bis(2-chloroethyl)ether	1.5059 1.4502	1.4945 1.3267	1.4527 1.3207	1.4517	1.4789	Ave		1.4352			0.7000	5.0		20.0			
2-Chlorophenol	1.5378 1.4906	1.5218 1.4477	1.4794 1.4288	1.5127	1.5199	Ave		1.4924			0.8000	2.6		20.0			
1,3-Dichlorobenzene	1.5887 1.5052	1.5554 1.4508	1.5140 1.4356	1.5306	1.5345	Ave		1.5143				3.4		20.0			
1,4-Dichlorobenzene	1.5919 1.5145	1.5614 1.4693	1.5561 1.4271	1.5552	1.5637	Ave		1.5299				3.6		20.0			
Benzyl alcohol	0.9436 0.9634	0.9013 0.9386	0.9191 0.9314	0.9512	0.9727	Ave		0.9402				2.5		20.0			
1,2-Dichlorobenzene	1.5762 1.4506	1.5061 1.4020	1.4846 1.3840	1.4764	1.4904	Ave		1.4713				4.1		20.0			
2-Methylphenol	1.3905 1.3558	1.3706 1.3166	1.3493 1.3142	1.3800	1.3655	Ave		1.3553			0.7000	2.1		20.0			
bis (2-chloroisopropyl) ether	2.2287 2.1058	2.1607 2.0300	2.1132 2.0111	2.1516	2.0987	Ave		2.1125			0.0100	3.3		20.0			
3 & 4 Methylphenol	1.4380 1.4065	1.3807 1.3842	1.4127 1.3677	1.4193	1.4295	Ave		1.4048				1.8		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
3-Methylphenol	1.4380 1.4065	1.3807 1.3842	1.4127 1.3677	1.4193	1.4295	Ave		1.4048				1.8		20.0			
4-Methylphenol	1.4380 1.4065	1.3807 1.3842	1.4127 1.3677	1.4193	1.4295	Ave		1.4048			0.6000	1.8		20.0			
N-Nitrosodi-n-propylamine	1.0616 1.0289	0.9941 0.9958	1.0192 0.9800	1.0472	1.0302	Ave		1.0196			0.5000	2.8		20.0			
Acetophenone	2.1078 1.9618	2.0248 1.8890	1.9599 1.8742	2.0034	1.9742	Ave		1.9744			0.0100	3.8		20.0			
Hexachloroethane	0.6768 0.6538	0.6642 0.6300	0.6557 0.6255	0.6686	0.6777	Ave		0.6565			0.3000	3.0		20.0			
Nitrobenzene	0.3952 0.4001	0.4067 0.3872	0.3929 0.3764	0.3996	0.4027	Ave		0.3951				2.4		20.0			
Isophorone	0.6745 0.6778	0.6572 0.6688	0.6649 0.6533	0.6861	0.6860	Ave		0.6711			0.4000	1.8		20.0			
2,4-Dimethylphenol	0.3673 0.3568	0.3792 0.3487	0.3639 0.3319	0.3677	0.3710	Ave		0.3608			0.2000	4.1		20.0			
2-Nitrophenol	0.2010 0.1974	0.2008 0.1901	0.1990 0.1838	0.2024	0.2008	Ave		0.1969			0.1000	3.3		20.0			
Benzoic acid	0.2283 0.3050	0.2539 0.3193	0.2690 0.3060	0.2845	0.2990	Ave		0.2831				10.9		20.0			
Bis(2-chloroethoxy)methane	0.4554 0.4368	0.4570 0.4303	0.4453 0.4126	0.4466	0.4448	Ave		0.4411			0.3000	3.3		20.0			
2,4-Dichlorophenol	0.3071 0.2979	0.2949 0.2915	0.2959 0.2808	0.2954	0.3007	Ave		0.2955			0.2000	2.6		20.0			
1,2,4-Trichlorobenzene	0.3128 0.3001	0.3066 0.2894	0.3030 0.2824	0.3061	0.3017	Ave		0.3003				3.3		20.0			
Naphthalene	1.1213 1.0599	1.0974 1.0207	1.0807 0.9865	1.0695	1.0760	Ave		1.0640			0.7000	4.0		20.0			
4-Chloroaniline	0.4969 0.4804	0.4885 0.4712	0.4771 0.4573	0.4854	0.4931	Ave		0.4813			0.0100	2.7		20.0			
Hexachlorobutadiene	0.1780 0.1621	0.1655 0.1604	0.1633 0.1522	0.1618	0.1658	Ave		0.1636			0.0100	4.4		20.0			
Caprolactam	0.1575 0.2059	0.1794 0.2017	0.1815 0.2002	0.1874	0.2000	Ave		0.1892				8.6		20.0			
4-Chloro-3-methylphenol	0.3232 0.3198	0.3181 0.3124	0.3112 0.3052	0.3170	0.3237	Ave		0.3163			0.2000	2.0		20.0			
2-Methylnaphthalene	0.7077 0.6803	0.7083 0.6577	0.6911 0.6323	0.6962	0.6952	Ave		0.6836			0.4000	3.9		20.0			
1-Methylnaphthalene	0.6652 0.6382	0.6677 0.6166	0.6544 0.5913	0.6561	0.6543	Ave		0.6430				4.1		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
Hexachlorocyclopentadiene	0.2947 0.3198	0.2965 0.3184	0.3107 0.3146	0.3234	0.3361	Ave		0.3143			0.0500	4.4		20.0			
1,2,4,5-Tetrachlorobenzene	0.2970 0.2796	0.2871 0.2748	0.2894 0.2692	0.2921	0.2876	Ave		0.2846			0.0100	3.3		20.0			
2,4,6-Trichlorophenol	0.3657 0.3545	0.3558 0.3520	0.3604 0.3446	0.3676	0.3727	Ave		0.3592			0.2000	2.6		20.0			
2,4,5-Trichlorophenol	0.3946 0.3800	0.3801 0.3716	0.3848 0.3684	0.3863	0.3949	Ave		0.3826			0.2000	2.5		20.0			
1,1'-Biphenyl	1.5894 1.4447	1.5563 1.4388	1.5496 1.3861	1.5342	1.5450	Ave		1.5055				4.8		20.0			
2-Chloronaphthalene	1.1700 1.1082	1.1772 1.0827	1.1682 1.0515	1.1586	1.1648	Ave		1.1352			0.8000	4.2		20.0			
2-Nitroaniline	0.3890 0.4212	0.4091 0.4207	0.4239 0.4111	0.4270	0.4327	Ave		0.4168			0.0100	3.3		20.0			
Dimethyl phthalate	1.5586 1.2704	1.3548 1.2539	1.3221 1.2346	1.3203	1.3386	Ave		1.3317			0.0100	7.6		20.0			
1,3-Dinitrobenzene	0.1942 0.2516	0.2110 0.2498	0.2336 0.2509	0.2438	0.2549	Ave		0.2362				9.4		20.0			
2,6-Dinitrotoluene	0.2960 0.3045	0.2992 0.3071	0.3210 0.3014	0.3137	0.3193	Ave		0.3078			0.2000	3.0		20.0			
Acenaphthylene	1.8624 1.8257	1.8627 1.7954	1.8590 1.7480	1.8945	1.8979	Ave		1.8432			0.9000	2.8		20.0			
3-Nitroaniline	0.3671 0.3971	0.3907 0.4035	0.3969 0.4015	0.4156	0.4117	Ave		0.3980			0.0100	3.7		20.0			
Acenaphthene	1.2405 1.1570	1.2286 1.1392	1.2217 1.0937	1.2241	1.2147	Ave		1.1899			0.9000	4.5		20.0			
2,4-Dinitrophenol	0.1476 0.2290	0.1713 0.2290	0.2002 0.2287	0.2216	0.2288	Lin2	-0.709	0.2271			0.0100				0.9980		0.9900
4-Nitrophenol	0.2212 0.2119	0.2128 0.2099	0.2124 0.2061	0.2174	0.2247	Ave		0.2146			0.0100	2.8		20.0			
2,4-Dinitrotoluene	0.3711 0.4093	0.3874 0.4073	0.4097 0.4040	0.4194	0.4204	Ave		0.4036			0.2000	4.1		20.0			
Dibenzofuran	1.7496 1.6359	1.7143 1.6052	1.7150 1.5494	1.7178	1.7204	Ave		1.6760			0.8000	4.2		20.0			
2,3,4,6-Tetrachlorophenol	0.2847 0.3155	0.3085 0.3095	0.3060 0.3115	0.3205	0.3251	Ave		0.3102			0.0100	3.9		20.0			
Diethyl phthalate	1.4120 1.3118	1.3980 1.2889	1.4006 1.2647	1.3831	1.3855	Ave		1.3556			0.0100	4.3		20.0			
4-Chlorophenyl phenyl ether	0.6326 0.5864	0.6274 0.5832	0.6207 0.5690	0.6193	0.6132	Ave		0.6065			0.4000	3.9		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
Fluorene	1.4028 1.3292	1.4196 1.2961	1.4070 1.2662	1.4084	1.3789	Ave		1.3635			0.9000	4.3		20.0			
4-Nitroaniline	0.3747 0.3850	0.3884 0.3879	0.3917 0.3738	0.4022	0.4094	Ave		0.3891			0.0100	3.2		20.0			
4,6-Dinitro-2-methylphenol	0.1138 0.1466	0.1344 0.1468	0.1430 0.1444	0.1507	0.1524	Ave		0.1415			0.0100	8.8		20.0			
N-Nitrosodiphenylamine	0.6054 0.5842	0.6159 0.5651	0.6141 0.5562	0.6024	0.5971	Ave		0.5925			0.0100	3.7		20.0			
1,2-Diphenylhydrazine(as Azobenzene)	1.5351 1.4624	1.4890 1.4353	1.4994 1.4075	1.5191	1.5093	Ave		1.4821				3.0		20.0			
Azobenzene	1.5519 1.4785	1.5054 1.4510	1.5159 1.4229	1.5357	1.5259	Ave		1.4984				3.0		20.0			
4-Bromophenyl phenyl ether	0.1791 0.1876	0.1939 0.1840	0.1898 0.1803	0.1924	0.1923	Ave		0.1874			0.1000	3.0		20.0			
Hexachlorobenzene	0.1990 0.1896	0.1977 0.1875	0.2002 0.1880	0.1991	0.1965	Ave		0.1947			0.1000	2.8		20.0			
Pentachlorophenol	0.1133 0.1315	0.1202 0.1292	0.1285 0.1296	0.1356	0.1360	Ave		0.1280			0.0500	6.0		20.0			
Phenanthrene	1.1647 1.0730	1.1605 1.0522	1.1306 1.0273	1.1385	1.1283	Ave		1.1094			0.7000	4.7		20.0			
Anthracene	1.1822 1.1094	1.1766 1.0889	1.1645 1.0583	1.1597	1.1547	Ave		1.1368			0.7000	4.0		20.0			
Carbazole	1.1800 1.1376	1.1904 1.1151	1.1973 1.0890	1.1840	1.1774	Ave		1.1589			0.0100	3.4		20.0			
Di-n-butyl phthalate	1.3294 1.3535	1.3595 1.3335	1.3832 1.3196	1.4207	1.4049	Ave		1.3630			0.0100	2.7		20.0			
Fluoranthene	1.2052 1.1977	1.2468 1.1836	1.2384 1.1659	1.2350	1.2483	Ave		1.2151			0.6000	2.6		20.0			
Pyrene	1.3471 1.3422	1.3443 1.3128	1.3236 1.2915	1.3549	1.3687	Ave		1.3356			0.6000	1.9		20.0			
Famphur	0.4275 0.4189	0.4372 0.3872	0.4432 0.3507	0.4559	0.4518	Ave		0.4215				8.5		20.0			
Butyl benzyl phthalate	0.5480 0.6654	0.5938 0.6573	0.6115 0.6497	0.6472	0.6721	Ave		0.6306			0.0100	6.8		20.0			
3,3'-Dichlorobenzidine	0.3550 0.4127	0.3739 0.4060	0.3796 0.3966	0.4110	0.4275	Ave		0.3953			0.0100	6.1		20.0			
Bis(2-ethylhexyl) phthalate	0.7171 0.8671	0.7740 0.8474	0.7980 0.8287	0.8610	0.8965	Ave		0.8237			0.0100	7.1		20.0			
Benzo[a]anthracene	1.1704 1.1473	1.1594 1.1299	1.1437 1.1089	1.1663	1.1712	Ave		1.1496			0.8000	1.9		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD CURVE EVALUATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

ANALYTE	RRF					CURVE TYPE	COEFFICIENT			#	MIN RRF	%RSD	#	MAX %RSD	R^2 OR COD	#	MIN R^2 OR COD
	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5		B	M1	M2								
Chrysene	1.0715 1.0812	1.0706 1.0677	1.0653 1.0532	1.0881	1.0910	Ave		1.0736			0.7000	1.2		20.0			
Di-n-octyl phthalate	0.9715 1.5292	1.1522 1.5274	1.2430 1.5267	1.4190	1.5392	Lin2	-2.366	1.4972			0.0100				0.9960		0.9900
Benzo[b]fluoranthene	1.0676 1.1983	1.1421 1.2214	1.1629 1.1805	1.1650	1.2102	Ave		1.1685			0.7000	4.2		20.0			
Benzo[k]fluoranthene	1.1189 1.2408	1.0998 1.2306	1.1683 1.2469	1.2353	1.2599	Ave		1.2001			0.7000	5.2		20.0			
Benzo[a]pyrene	0.9584 1.1859	1.0416 1.1888	1.0918 1.1801	1.1657	1.1745	Ave		1.1233			0.7000	7.6		20.0			
Indeno[1,2,3-cd]pyrene	0.7186 0.8643	0.7610 0.8576	0.7749 0.8581	0.8305	0.8680	Ave		0.8166			0.5000	7.0		20.0			
Dibenz(a,h)anthracene	0.7853 0.9916	0.8077 0.9941	0.9011 0.9801	0.9711	0.9998	Ave		0.9288			0.4000	9.4		20.0			
Benzo[g,h,i]perylene	0.9134 1.0580	0.9613 1.0482	0.9725 1.0450	1.0270	1.0597	Ave		1.0106			0.5000	5.4		20.0			
2-Fluorophenol (Surr)	1.4984 1.4619	1.4670 1.4249	1.4461 1.4139	1.4548	1.4680	Ave		1.4544				1.8		20.0			
Phenol-d5 (Surr)	1.8890 1.8249	1.7818 1.7845	1.8331 1.7781	1.8364	1.8548	Ave		1.8228				2.2		20.0			
Nitrobenzene-d5 (Surr)	0.4132 0.4157	0.4145 0.4054	0.4052 0.3934	0.4128	0.4209	Ave		0.4101				2.1		20.0			
2-Fluorobiphenyl	1.4249 1.3139	1.3935 1.2755	1.3897 1.2507	1.3777	1.3915	Ave		1.3522				4.7		20.0			
2,4,6-Tribromophenol (Surr)	0.1405 0.1566	0.1391 0.1540	0.1478 0.1563	0.1539	0.1551	Ave		0.1504				4.7		20.0			
Terphenyl-d14 (Surr)	0.7894 0.7962	0.8165 0.7812	0.7915 0.7699	0.8063	0.8080	Ave		0.7949				1.9		20.0			

Note: The m1 coefficient is the same as Ave RRF for an Ave curve type.

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

Calibration Files:

LEVEL:	LAB SAMPLE ID:	LAB FILE ID:
Level 1	STD004 280-272059/4	G6_16638.D
Level 2	STD010 280-272059/5	G6_16639.D
Level 3	STD020 280-272059/6	G6_16640.D
Level 4	STD050 280-272059/7	G6_16641.D
Level 5	ICIS 280-272059/3	G6_16637.D
Level 6	STD120 280-272059/8	G6_16642.D
Level 7	STD160 280-272059/9	G6_16643.D
Level 8	STD200 280-272059/10	G6_16644.D

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
1,4-Dioxane	DCB	Ave	11382 333966	29946 439534	59534 558095	142371	232210	4.00 120	10.0 160	20.0 200	50.0	80.0
N-Nitrosodimethylamine	DCB	Ave	17202 518985	43094 689099	85710 878936	222746	351228	4.00 120	10.0 160	20.0 200	50.0	80.0
Pyridine	DCB	Ave	29364 891222	73629 1173437	147856 1482135	375517	600581	4.00 120	10.0 160	20.0 200	50.0	80.0
Phenol	DCB	Ave	34473 999282	85314 1321311	172136 1676043	436251	698622	4.00 120	10.0 160	20.0 200	50.0	80.0
Aniline	DCB	Ave	43366 1234749	101304 1657573	206891 2067510	529556	834965	4.00 120	10.0 160	20.0 200	50.0	80.0
Bis(2-chloroethyl)ether	DCB	Ave	27476 802000	69073 988186	137208 1252467	342205	553186	4.00 120	10.0 160	20.0 200	50.0	80.0
2-Chlorophenol	DCB	Ave	28058 824383	70338 1078332	139729 1354936	356590	568507	4.00 120	10.0 160	20.0 200	50.0	80.0
1,3-Dichlorobenzene	DCB	Ave	28987 832464	71887 1080606	142995 1361408	360797	573965	4.00 120	10.0 160	20.0 200	50.0	80.0
1,4-Dichlorobenzene	DCB	Ave	29045 837561	72168 1094435	146967 1353318	366601	584889	4.00 120	10.0 160	20.0 200	50.0	80.0
Benzyl alcohol	DCB	Ave	17216 532820	41657 699135	86805 883290	224224	363847	4.00 120	10.0 160	20.0 200	50.0	80.0
1,2-Dichlorobenzene	DCB	Ave	28758 802221	69610 1044311	140215 1312433	348029	557476	4.00 120	10.0 160	20.0 200	50.0	80.0
2-Methylphenol	DCB	Ave	25371 749809	63348 980672	127434 1246269	325292	510773	4.00 120	10.0 160	20.0 200	50.0	80.0
bis (2-chloroisopropyl) ether	DCB	Ave	40664 1164573	99866 1512052	199591 1907157	507190	785020	4.00 120	10.0 160	20.0 200	50.0	80.0
3 & 4 Methylphenol	DCB	Ave	26237 777835	63816 1031059	133429 1296981	334564	534702	4.00 120	10.0 160	20.0 200	50.0	80.0
3-Methylphenol	DCB	Ave	26237 777835	63816 1031059	133429 1296981	334564	534702	4.00 120	10.0 160	20.0 200	50.0	80.0

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
4-Methylphenol	DCB	Ave	26237 777835	63816 1031059	133429 1296981	334564	534702	4.00 120	10.0 160	20.0 200	50.0	80.0
N-Nitrosodi-n-propylamine	DCB	Ave	19369 569051	45947 741748	96259 929339	246852	385350	4.00 120	10.0 160	20.0 200	50.0	80.0
Acetophenone	DCB	Ave	38458 1084969	93583 1407057	185111 1777329	472247	738421	4.00 120	10.0 160	20.0 200	50.0	80.0
Hexachloroethane	DCB	Ave	12349 361563	30699 469286	61931 593205	157602	253482	4.00 120	10.0 160	20.0 200	50.0	80.0
Nitrobenzene	NPT	Ave	27855 842823	71302 1089690	141833 1376078	365089	574579	4.00 120	10.0 160	20.0 200	50.0	80.0
Isophorone	NPT	Ave	47544 1427939	115206 1881926	240037 2388530	626824	978816	4.00 120	10.0 160	20.0 200	50.0	80.0
2,4-Dimethylphenol	NPT	Ave	25893 751706	66475 981303	131383 1213524	335864	529293	4.00 120	10.0 160	20.0 200	50.0	80.0
2-Nitrophenol	NPT	Ave	14166 415910	35195 534875	71833 672106	184945	286572	4.00 120	10.0 160	20.0 200	50.0	80.0
Benzoic acid	NPT	Ave	32187 1285043	89023 1796945	194250 2237473	519854	853321	8.00 240	20.0 320	40.0 400	100	160
Bis(2-chloroethoxy)methane	NPT	Ave	32100 920059	80123 1210771	160771 1508491	407948	634640	4.00 120	10.0 160	20.0 200	50.0	80.0
2,4-Dichlorophenol	NPT	Ave	21647 627452	51694 820144	106817 1026716	269879	429047	4.00 120	10.0 160	20.0 200	50.0	80.0
1,2,4-Trichlorobenzene	NPT	Ave	22051 632193	53749 814451	109397 1032581	279647	430534	4.00 120	10.0 160	20.0 200	50.0	80.0
Naphthalene	NPT	Ave	79041 2232749	192392 2872188	390163 3606766	977042	1535294	4.00 120	10.0 160	20.0 200	50.0	80.0
4-Chloroaniline	NPT	Ave	35029 1011967	85639 1326050	172259 1671870	443432	703626	4.00 120	10.0 160	20.0 200	50.0	80.0
Hexachlorobutadiene	NPT	Ave	12544 341381	29013 451504	58938 556376	147770	236535	4.00 120	10.0 160	20.0 200	50.0	80.0
Caprolactam	NPT	Ave	11099 433839	31448 567672	65538 732007	171197	285328	4.00 120	10.0 160	20.0 200	50.0	80.0
4-Chloro-3-methylphenol	NPT	Ave	22782 673785	55761 879210	112366 1115868	289585	461868	4.00 120	10.0 160	20.0 200	50.0	80.0
2-Methylnaphthalene	NPT	Ave	49887 1433043	124165 1850886	249498 2311519	636017	991995	4.00 120	10.0 160	20.0 200	50.0	80.0
1-Methylnaphthalene	NPT	Ave	46889 1344403	117053 1735161	236260 2161606	599383	933568	4.00 120	10.0 160	20.0 200	50.0	80.0
Hexachlorocyclopentadiene	ANT	Ave	11449 376192	28473 489123	60610 621756	159056	259090	4.00 120	10.0 160	20.0 200	50.0	80.0
1,2,4,5-Tetrachlorobenzene	NPT	Ave	20934 588950	50324 773416	104495 984311	266816	410334	4.00 120	10.0 160	20.0 200	50.0	80.0

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
2,4,6-Trichlorophenol	ANT	Ave	14208 416933	34170 540716	70312 681085	180797	287284	4.00 120	10.0 160	20.0 200	50.0	80.0
2,4,5-Trichlorophenol	ANT	Ave	15333 446976	36506 570902	75074 728184	189993	304384	4.00 120	10.0 160	20.0 200	50.0	80.0
1,1'-Biphenyl	ANT	Ave	61753 1699215	149463 2210281	302324 2739578	754596	1190895	4.00 120	10.0 160	20.0 200	50.0	80.0
2-Chloronaphthalene	ANT	Ave	45458 1303430	113058 1663115	227907 2078262	569839	897878	4.00 120	10.0 160	20.0 200	50.0	80.0
2-Nitroaniline	ANT	Ave	15113 495382	39287 646189	82700 812592	210036	333520	4.00 120	10.0 160	20.0 200	50.0	80.0
Dimethyl phthalate	ANT	Ave	60554 1494143	130115 1926232	257927 2440056	649391	1031864	4.00 120	10.0 160	20.0 200	50.0	80.0
1,3-Dinitrobenzene	ANT	Ave	7547 295917	20263 383728	45571 495821	119893	196521	4.00 120	10.0 160	20.0 200	50.0	80.0
2,6-Dinitrotoluene	ANT	Ave	11500 358122	28732 471768	62625 595647	154313	246140	4.00 120	10.0 160	20.0 200	50.0	80.0
Acenaphthylene	ANT	Ave	72357 2147354	178896 2757941	362680 3454854	931810	1462928	4.00 120	10.0 160	20.0 200	50.0	80.0
3-Nitroaniline	ANT	Ave	14261 467084	37522 619865	77435 793558	204409	317373	4.00 120	10.0 160	20.0 200	50.0	80.0
Acenaphthene	ANT	Ave	48195 1360779	117997 1750014	238340 2161568	602051	936356	4.00 120	10.0 160	20.0 200	50.0	80.0
2,4-Dinitrophenol	ANT	Lin2	11473 538794	32905 703654	78117 903939	217993	352771	8.00 240	20.0 320	40.0 400	100	160
4-Nitrophenol	ANT	Ave	17189 498510	40880 644961	82891 814776	213891	346346	8.00 240	20.0 320	40.0 400	100	160
2,4-Dinitrotoluene	ANT	Ave	14418 481458	37208 625683	79927 798404	206301	324018	4.00 120	10.0 160	20.0 200	50.0	80.0
Dibenzofuran	ANT	Ave	67976 1924090	164638 2465817	334583 3062264	844896	1326145	4.00 120	10.0 160	20.0 200	50.0	80.0
2,3,4,6-Tetrachlorophenol	ANT	Ave	11063 371083	29631 475470	59694 615581	157611	250593	4.00 120	10.0 160	20.0 200	50.0	80.0
Diethyl phthalate	ANT	Ave	54858 1542842	134263 1980010	273258 2499587	680249	1067965	4.00 120	10.0 160	20.0 200	50.0	80.0
4-Chlorophenyl phenyl ether	ANT	Ave	24578 689660	60258 895954	121088 1124576	304597	472676	4.00 120	10.0 160	20.0 200	50.0	80.0
Fluorene	ANT	Ave	54503 1563412	136339 1991029	274499 2502594	692688	1062927	4.00 120	10.0 160	20.0 200	50.0	80.0
4-Nitroaniline	ANT	Ave	14556 452856	37301 595856	76419 738699	197830	315582	4.00 120	10.0 160	20.0 200	50.0	80.0
4,6-Dinitro-2-methylphenol	PHN	Ave	14971 578792	43031 761080	93754 958230	249423	398341	8.00 240	20.0 320	40.0 400	100	160

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
N-Nitrosodiphenylamine	PHN	Ave	39816 1152861	98599 1464837	201342 1845140	498390	780136	4.00 120	10.0 160	20.0 200	50.0	80.0
1,2-Diphenylhydrazine(as Azobenzene)	ANT	Ave	60296 1738928	144575 2228985	295740 2812289	755331	1176176	4.04 121	10.1 162	20.2 202	50.5	80.9
Azobenzene	ANT	Ave	60296 1738928	144575 2228985	295740 2812289	755331	1176176	4.00 120	10.0 160	20.0 200	50.0	80.0
4-Bromophenyl phenyl ether	PHN	Ave	11781 370159	31038 477097	62227 598149	159165	251254	4.00 120	10.0 160	20.0 200	50.0	80.0
Hexachlorobenzene	PHN	Ave	13090 374109	31646 486080	65622 623678	164764	256694	4.00 120	10.0 160	20.0 200	50.0	80.0
Pentachlorophenol	PHN	Ave	14903 519082	38486 669774	84234 859647	224380	355328	8.00 240	20.0 320	40.0 400	100	160
Phenanthrene	PHN	Ave	76605 2117512	185797 2727512	370680 3407975	941975	1474345	4.00 120	10.0 160	20.0 200	50.0	80.0
Anthracene	PHN	Ave	77754 2189501	188375 2822765	381794 3510820	959498	1508829	4.00 120	10.0 160	20.0 200	50.0	80.0
Carbazole	PHN	Ave	77606 2245142	190570 2890531	392545 3612710	979659	1538487	4.00 120	10.0 160	20.0 200	50.0	80.0
Di-n-butyl phthalate	PHN	Ave	87433 2671086	217656 3456767	453491 4377647	1175500	1835676	4.00 120	10.0 160	20.0 200	50.0	80.0
Fluoranthene	PHN	Ave	79265 2363681	199608 3068105	405997 3867766	1021800	1631054	4.00 120	10.0 160	20.0 200	50.0	80.0
Pyrene	CRY	Ave	85070 2459885	206008 3191260	418438 4045386	1068582	1686444	4.00 120	10.0 160	20.0 200	50.0	80.0
Famphur	CRY	Ave	26998 767713	66995 941175	140106 1098625	359540	556669	4.00 120	10.0 160	20.0 200	50.0	80.0
Butyl benzyl phthalate	CRY	Ave	34606 1219463	91001 1597755	193326 2035238	510447	828145	4.00 120	10.0 160	20.0 200	50.0	80.0
3,3'-Dichlorobenzidine	CRY	Ave	22416 756387	57301 987057	120010 1242385	324147	526780	4.00 120	10.0 160	20.0 200	50.0	80.0
Bis(2-ethylhexyl) phthalate	CRY	Ave	45285 1589107	118615 2060022	252271 2595812	679050	1104576	4.00 120	10.0 160	20.0 200	50.0	80.0
Benzo[a]anthracene	CRY	Ave	73908 2102748	177675 2746652	361571 3473559	919851	1443039	4.00 120	10.0 160	20.0 200	50.0	80.0
Chrysene	CRY	Ave	67664 1981581	164066 2595420	336779 3298874	858164	1344205	4.00 120	10.0 160	20.0 200	50.0	80.0
Di-n-octyl phthalate	CRY	Lin2	61347 2802646	176570 3712823	392967 4782051	1119188	1896507	4.00 120	10.0 160	20.0 200	50.0	80.0
Benzo[b]fluoranthene	PRY	Ave	57364 1901054	151541 2549534	312315 3151061	791973	1287729	4.00 120	10.0 160	20.0 200	50.0	80.0
Benzo[k]fluoranthene	PRY	Ave	60118 1968536	145931 2568840	313777 3328296	839735	1340597	4.00 120	10.0 160	20.0 200	50.0	80.0

FORM VI
GC/MS SEMI VOA INITIAL CALIBRATION DATA
INTERNAL STANDARD RESPONSE AND CONCENTRATION

Lab Name: TestAmerica Denver Job No.: 280-67711-2 Analy Batch No.: 272059

SDG No.:

Instrument ID: SMS_G6 GC Column: Vf-5MS (30.2 ID: 0.25(mm) Heated Purge: (Y/N) N

Calibration Start Date: 02/25/2015 11:53 Calibration End Date: 02/25/2015 14:59 Calibration ID: 21947

ANALYTE	IS REF	CURVE TYPE	RESPONSE					CONCENTRATION (UG/ML)				
			LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5	LVL 1 LVL 6	LVL 2 LVL 7	LVL 3 LVL 8	LVL 4	LVL 5
Benzo[a]pyrene	PRY	Ave	51493 1881464	138204 2481467	293226 3149998	792402	1249715	4.00 120	10.0 160	20.0 200	50.0	80.0
Indeno[1,2,3-cd]pyrene	CRY	Ave	45380 1584058	116621 2084773	244980 2687797	654992	1069526	4.00 120	10.0 160	20.0 200	50.0	80.0
Dibenz(a,h)anthracene	PRY	Ave	42195 1573114	107166 2075077	242009 2616064	660134	1063872	4.00 120	10.0 160	20.0 200	50.0	80.0
Benzo[g,h,i]perylene	PRY	Ave	49079 1678491	127555 2188108	261176 2789274	698128	1127580	4.00 120	10.0 160	20.0 200	50.0	80.0
2-Fluorophenol (Surr)	DCB	Ave	27340 808477	67803 1061373	136584 1340808	342921	549103	4.00 120	10.0 160	20.0 200	50.0	80.0
Phenol-d5 (Surr)	DCB	Ave	34467 1009270	82351 1329218	173137 1686187	432891	693790	4.00 120	10.0 160	20.0 200	50.0	80.0
Nitrobenzene-d5 (Surr)	NPT	Ave	29128 875604	72669 1140831	146288 1438315	377126	600607	4.00 120	10.0 160	20.0 200	50.0	80.0
2-Fluorobiphenyl	ANT	Ave	55360 1545378	133834 1959377	271127 2471944	677613	1072615	4.00 120	10.0 160	20.0 200	50.0	80.0
2,4,6-Tribromophenol (Surr)	ANT	Ave	5458 184203	13361 236631	28834 308954	75692	119592	4.00 120	10.0 160	20.0 200	50.0	80.0
Terphenyl-d14 (Surr)	CRY	Ave	49847 1459232	125136 1898972	250222 2411505	635940	995559	4.00 120	10.0 160	20.0 200	50.0	80.0

Curve Type Legend:

Ave = Average ISTD
Lin2 = Linear 1/conc^2 ISTD

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16637.D
Lims ID: ICIS HSL
Client ID:
Sample Type: ICIS Calib Level: 5
Inject. Date: 25-Feb-2015 11:53:30 ALS Bottle#: 2 Worklist Smp#: 3
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: ICIS HSL
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:37 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: kiekeld Date: 10-Apr-2015 08:10:13

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	96	187021	40.0	40.0	
* 2 Naphthalene-d8	136	6.022	6.022	0.000	100	713420	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	92	385413	40.0	40.0	
* 4 Phenanthrene-d10	188	9.280	9.280	0.000	97	653325	40.0	40.0	
* 5 Chrysene-d12	240	13.710	13.710	0.000	96	616068	40.0	40.0	
* 6 Perylene-d12	264	17.903	17.903	0.000	95	532020	40.0	40.0	
$ 7 2-Fluorophenol	112	3.622	3.622	0.000	94	549103	80.0	80.8	
$ 8 Phenol-d5	99	4.410	4.410	0.000	98	693790	80.0	81.4	
$ 9 Nitrobenzene-d5	82	5.322	5.322	0.000	91	600607	80.0	82.1	
$ 10 2-Fluorobiphenyl	172	7.069	7.069	0.000	100	1072615	80.0	82.3	
$ 11 2,4,6-Tribromophenol	330	8.575	8.575	0.000	88	119592	80.0	82.5	
$ 12 Terphenyl-d14	244	11.369	11.369	0.000	98	995559	80.0	81.3	
13 1,4-Dioxane	88	2.352	2.352	0.000	98	232210	80.0	80.9	
14 N-Nitrosodimethylamine	74	2.581	2.581	0.000	90	351228	80.0	80.6	
15 Pyridine	79	2.622	2.622	0.000	91	600581	80.0	80.8	
23 Phenol	94	4.422	4.422	0.000	99	698622	80.0	81.7	
24 Aniline	93	4.498	4.498	0.000	98	834965	80.0	79.9	
25 Bis(2-chloroethyl)ether	93	4.528	4.528	0.000	90	553186	80.0	82.4	
26 2-Chlorophenol	128	4.604	4.604	0.000	97	568507	80.0	81.5	
27 1,3-Dichlorobenzene	146	4.751	4.751	0.000	97	573965	80.0	81.1	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	93	584889	80.0	81.8	
29 Benzyl alcohol	108	4.904	4.904	0.000	93	363847	80.0	82.8	
30 1,2-Dichlorobenzene	146	4.963	4.963	0.000	96	557476	80.0	81.0	
31 2-Methylphenol	108	4.981	4.981	0.000	96	510773	80.0	80.6	
32 2,2'-oxybis[1-chloropropan	45	5.010	5.010	0.000	93	785020	80.0	79.5	
38 3 & 4 Methylphenol	108	5.128	5.128	0.000	98	534702	80.0	81.4	
39 3-Methylphenol	108	5.128	5.128	0.000	98	534702	80.0	81.4	
40 4-Methylphenol	108	5.128	5.128	0.000	93	534702	80.0	81.4	
41 N-Nitrosodi-n-propylamine	70	5.145	5.145	0.000	90	385350	80.0	80.8	
42 Acetophenone	105	5.163	5.163	0.000	96	738421	80.0	80.0	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.281	5.281	0.000	96	253482	80.0	82.6	
44 Nitrobenzene	77	5.340	5.340	0.000	91	574579	80.0	81.5	
46 Isophorone	82	5.545	5.545	0.000	99	978816	80.0	81.8	
48 2-Nitrophenol	139	5.645	5.645	0.000	93	286572	80.0	81.6	
49 2,4-Dimethylphenol	107	5.640	5.640	0.000	92	529293	80.0	82.2	
50 Bis(2-chloroethoxy)methane	93	5.734	5.734	0.000	98	634640	80.0	80.7	
52 Benzoic acid	105	5.728	5.728	0.000	90	853321	160.0	169.0	
53 2,4-Dichlorophenol	162	5.863	5.863	0.000	96	429047	80.0	81.4	
54 1,2,4-Trichlorobenzene	180	5.951	5.951	0.000	94	430534	80.0	80.4	
57 Naphthalene	128	6.040	6.040	0.000	98	1535294	80.0	80.9	
58 4-Chloroaniline	127	6.081	6.081	0.000	96	703626	80.0	82.0	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	97	236535	80.0	81.1	
62 Caprolactam	55	6.428	6.428	0.000	78	285328	80.0	84.6	M
64 4-Chloro-3-methylphenol	107	6.528	6.528	0.000	97	461868	80.0	81.9	
65 2-Methylnaphthalene	142	6.716	6.716	0.000	93	991995	80.0	81.4	
67 1-Methylnaphthalene	142	6.822	6.822	0.000	94	933568	80.0	81.4	
68 Hexachlorocyclopentadiene	237	6.863	6.863	0.000	97	259090	80.0	85.6	
69 1,2,4,5-Tetrachlorobenzene	216	6.887	6.887	0.000	98	410334	80.0	80.8	
70 2,4,6-Trichlorophenol	196	6.992	6.992	0.000	95	287284	80.0	83.0	
72 2,4,5-Trichlorophenol	196	7.028	7.028	0.000	93	304384	80.0	82.6	
74 1,1'-Biphenyl	154	7.181	7.181	0.000	95	1190895	80.0	82.1	
75 2-Chloronaphthalene	162	7.216	7.216	0.000	97	897878	80.0	82.1	
77 2-Nitroaniline	65	7.316	7.316	0.000	84	333520	80.0	83.0	
79 Dimethyl phthalate	163	7.463	7.463	0.000	98	1031864	80.0	80.4	
80 1,3-Dinitrobenzene	168	7.528	7.528	0.000	84	196521	80.0	86.3	
81 2,6-Dinitrotoluene	165	7.545	7.545	0.000	93	246140	80.0	83.0	
82 Acenaphthylene	152	7.639	7.639	0.000	99	1462928	80.0	82.4	
83 3-Nitroaniline	138	7.728	7.728	0.000	94	317373	80.0	82.8	
84 Acenaphthene	153	7.810	7.810	0.000	96	936356	80.0	81.7	
86 2,4-Dinitrophenol	184	7.828	7.828	0.000	84	352771	160.0	164.4	
87 4-Nitrophenol	109	7.869	7.869	0.000	95	346346	160.0	167.5	
89 2,4-Dinitrotoluene	165	7.951	7.951	0.000	91	324018	80.0	83.3	
90 Dibenzofuran	168	7.981	7.981	0.000	97	1326145	80.0	82.1	
92 2,3,4,6-Tetrachlorophenol	232	8.092	8.092	0.000	77	250593	80.0	83.9	
94 Diethyl phthalate	149	8.157	8.157	0.000	98	1067965	80.0	81.8	
96 4-Chlorophenyl phenyl ethe	204	8.298	8.298	0.000	97	472676	80.0	80.9	
98 Fluorene	166	8.328	8.328	0.000	95	1062927	80.0	80.9	
99 4-Nitroaniline	138	8.357	8.357	0.000	85	315582	80.0	84.2	
100 4,6-Dinitro-2-methylphenol	198	8.363	8.363	0.000	85	398341	160.0	172.3	
102 N-Nitrosodiphenylamine	169	8.422	8.422	0.000	63	780136	80.0	80.6	
103 Azobenzene	77	8.463	8.463	0.000	99	1176176	80.0	81.5	
104 1,2-Diphenylhydrazine	77	8.463	8.463	0.000	100	1176176	80.9	82.4	
111 4-Bromophenyl phenyl ether	248	8.798	8.798	0.000	72	251254	80.0	82.1	
112 Hexachlorobenzene	284	8.881	8.881	0.000	92	256694	80.0	80.7	
116 Pentachlorophenol	266	9.080	9.080	0.000	91	355328	160.0	170.0	
119 Phenanthrene	178	9.310	9.310	0.000	98	1474345	80.0	81.4	
120 Anthracene	178	9.357	9.357	0.000	98	1508829	80.0	81.3	
122 Carbazole	167	9.522	9.522	0.000	96	1538487	80.0	81.3	
123 Di-n-butyl phthalate	149	9.822	9.822	0.000	100	1835676	80.0	82.5	
128 Fluoranthene	202	10.786	10.786	0.000	98	1631054	80.0	82.2	
131 Pyrene	202	11.157	11.157	0.000	97	1686444	80.0	82.0	
136 Famphur	218	12.198	12.198	0.000	96	556669	80.0	85.7	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
137 Butyl benzyl phthalate	149	12.298	12.298	0.000	98	828145	80.0	85.3	
140 3,3'-Dichlorobenzidine	252	13.639	13.639	0.000	76	526780	80.0	86.5	
141 Benzo[a]anthracene	228	13.680	13.680	0.000	99	1443039	80.0	81.5	
142 Bis(2-ethylhexyl) phthalat	149	13.668	13.668	0.000	97	1104576	80.0	87.1	
143 Chrysene	228	13.774	13.774	0.000	98	1344205	80.0	81.3	
144 Di-n-octyl phthalate	149	15.574	15.574	0.000	99	1896507	80.0	83.8	
146 Benzo[b]fluoranthene	252	16.727	16.727	0.000	98	1287729	80.0	82.9	
147 Benzo[k]fluoranthene	252	16.815	16.815	0.000	98	1340597	80.0	84.0	
148 Benzo[a]pyrene	252	17.727	17.727	0.000	80	1249715	80.0	83.6	
151 Indeno[1,2,3-cd]pyrene	276	21.127	21.127	0.000	96	1069526	80.0	85.0	M
152 Dibenz(a,h)anthracene	278	21.203	21.203	0.000	96	1063872	80.0	86.1	
153 Benzo[g,h,i]perylene	276	21.897	21.897	0.000	95	1127580	80.0	83.9	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

MS-HSLA080_00015 Amount Added: 200.00 Units: uL

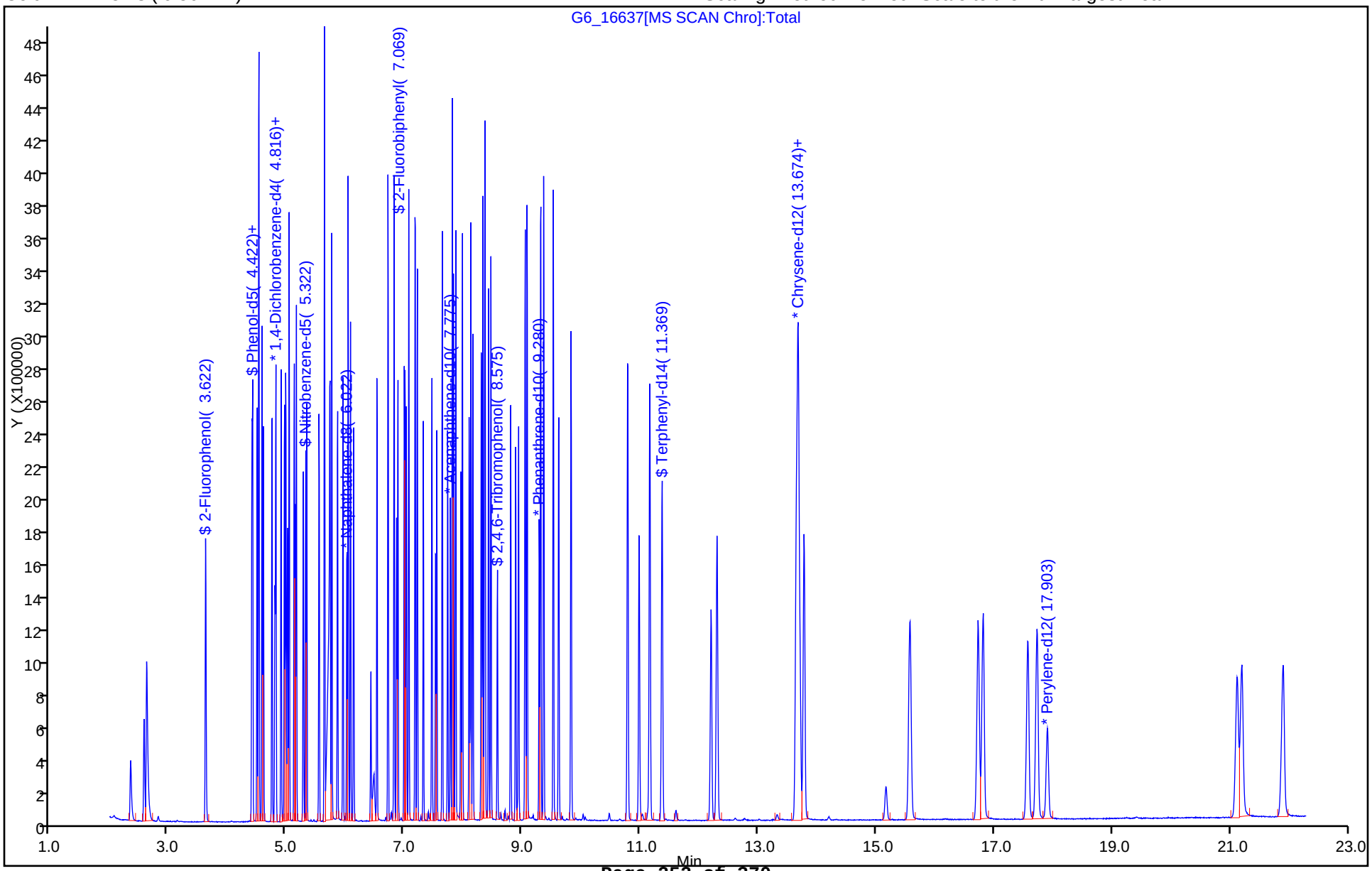
Report Date: 10-Apr-2015 12:09:38 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16637.D

Injection Date:	25-Feb-2015 11:53:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	ICIS HSL			Worklist Smp#:	3
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	2
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



Manual Integration/User Assign Peak Report

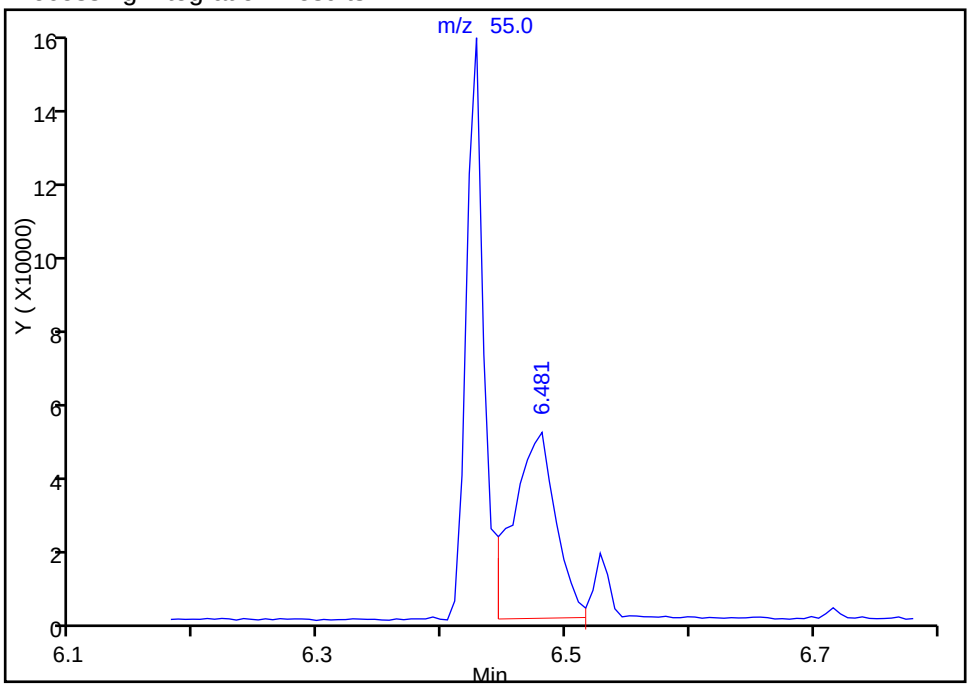
TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16637.D
Injection Date: 25-Feb-2015 11:53:30 Instrument ID: SMS_G6
Lims ID: ICIS HSL
Client ID:
Operator ID: KIEKELD ALS Bottle#: 2 Worklist Smp#: 3
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Method: SMS_G6_8270D Limit Group: MSSV - 8270D
Column: VF-5ms (0.50 mm) Detector MS SCAN

62 Caprolactam, CAS: 105-60-2

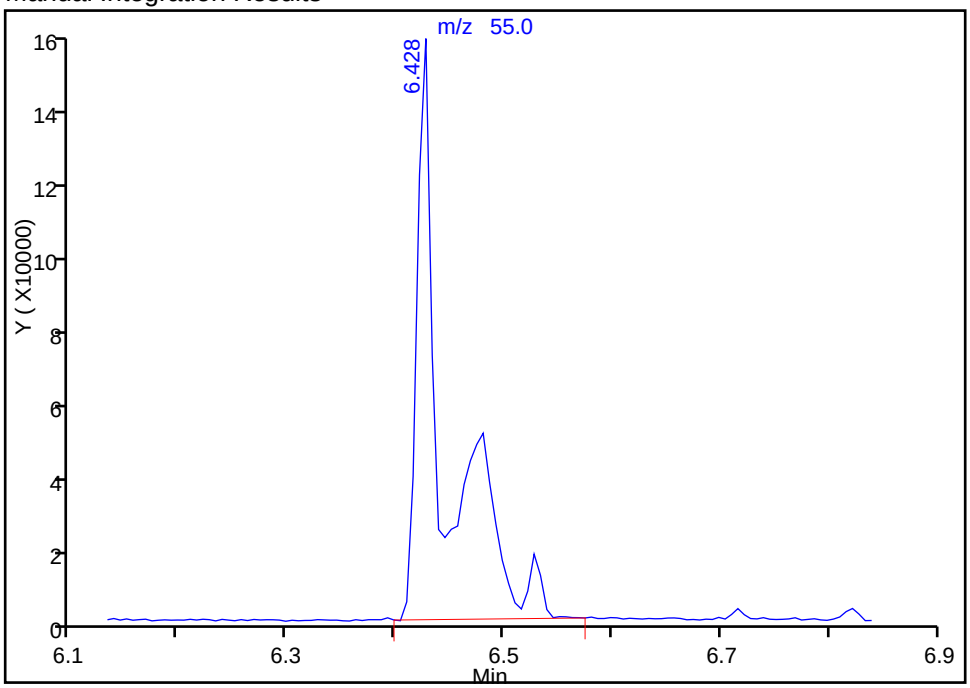
Processing Integration Results

RT: 6.48
Area: 122405
Amount: 80.000000
Amount Units: ug/ml



Manual Integration Results

RT: 6.43
Area: 285328
Amount: 84.551985
Amount Units: ug/ml



Reviewer: kiekeld, 10-Apr-2015 08:28:05
Audit Action: Manually Integrated
Audit Reason: Split Peak

Manual Integration/User Assign Peak Report

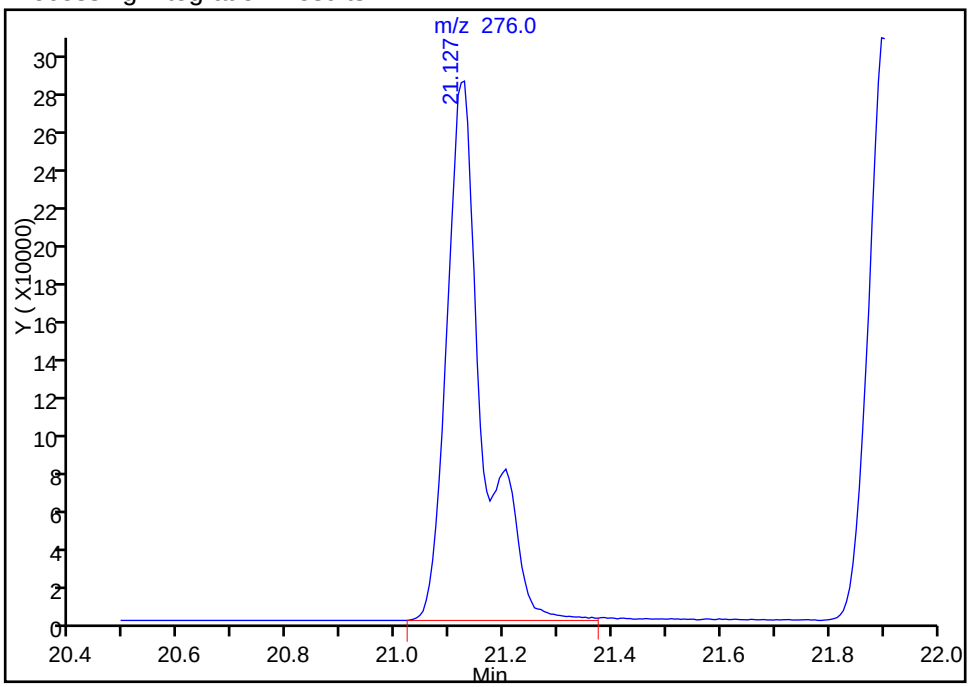
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16637.D				
Injection Date:	25-Feb-2015 11:53:30	Instrument ID:	SMS_G6		
Lims ID:	ICIS HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	2	Worklist Smp#:	3
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

151 Indeno[1,2,3-cd]pyrene, CAS: 193-39-5

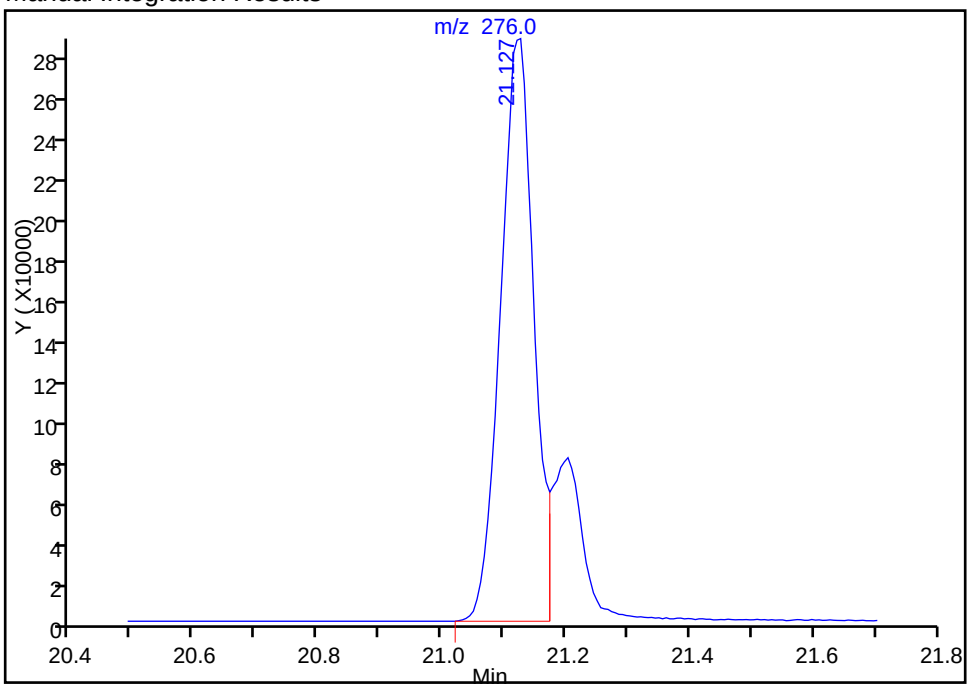
Processing Integration Results

RT: 21.13
Area: 1331166
Amount: 80.000000
Amount Units: ug/ml



Manual Integration Results

RT: 21.13
Area: 1069526
Amount: 85.035106
Amount Units: ug/ml



Reviewer: kiekeld, 10-Apr-2015 08:28:05
Audit Action: Split an Integrated Peak
Audit Reason: Shouldering

Report Date: 10-Apr-2015 12:09:39 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16638.D
Lims ID: STD004 HSL
Client ID:
Sample Type: IC Calib Level: 1
Inject. Date: 25-Feb-2015 12:19:30 ALS Bottle#: 3 Worklist Smp#: 4
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: STD004 HSL
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:38 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 08:51:31

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	97	182457	40.0	40.0	
* 2 Naphthalene-d8	136	6.016	6.022	-0.006	100	704885	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	92	388522	40.0	40.0	
* 4 Phenanthrene-d10	188	9.280	9.280	0.000	97	657696	40.0	40.0	
* 5 Chrysene-d12	240	13.698	13.710	-0.012	97	631485	40.0	40.0	
* 6 Perylene-d12	264	17.898	17.903	-0.005	95	537308	40.0	40.0	
$ 7 2-Fluorophenol	112	3.622	3.622	0.000	93	27340	4.00	4.12	
$ 8 Phenol-d5	99	4.404	4.410	-0.006	98	34467	4.00	4.15	
$ 9 Nitrobenzene-d5	82	5.322	5.322	0.000	91	29128	4.00	4.03	
$ 10 2-Fluorobiphenyl	172	7.069	7.069	0.000	99	55360	4.00	4.22	
$ 11 2,4,6-Tribromophenol	330	8.575	8.575	0.000	82	5458	4.00	3.74	
$ 12 Terphenyl-d14	244	11.363	11.369	-0.006	98	49847	4.00	3.97	
13 1,4-Dioxane	88	2.351	2.352	-0.001	99	11382	4.00	4.07	
14 N-Nitrosodimethylamine	74	2.581	2.581	0.000	93	17202	4.00	4.05	
15 Pyridine	79	2.628	2.622	0.006	87	29364	4.00	4.05	
23 Phenol	94	4.416	4.422	-0.006	98	34473	4.00	4.13	
24 Aniline	93	4.493	4.498	-0.006	98	43366	4.00	4.25	
25 Bis(2-chloroethyl)ether	93	4.528	4.528	0.000	91	27476	4.00	4.20	
26 2-Chlorophenol	128	4.604	4.604	0.000	96	28058	4.00	4.12	
27 1,3-Dichlorobenzene	146	4.751	4.751	0.000	97	28987	4.00	4.20	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	92	29045	4.00	4.16	
29 Benzyl alcohol	108	4.904	4.904	0.000	93	17216	4.00	4.01	
30 1,2-Dichlorobenzene	146	4.957	4.963	-0.006	95	28758	4.00	4.29	
31 2-Methylphenol	108	4.981	4.981	0.000	95	25371	4.00	4.10	
32 2,2'-oxybis[1-chloropropan	45	5.010	5.010	0.000	94	40664	4.00	4.22	
38 3 & 4 Methylphenol	108	5.122	5.128	-0.006	97	26237	4.00	4.09	
39 3-Methylphenol	108	5.122	5.128	-0.006	97	26237	4.00	4.09	
40 4-Methylphenol	108	5.122	5.128	-0.006	95	26237	4.00	4.09	
41 N-Nitrosodi-n-propylamine	70	5.140	5.145	-0.005	89	19369	4.00	4.16	
42 Acetophenone	105	5.163	5.163	0.000	96	38458	4.00	4.27	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.281	5.281	0.000	95	12349	4.00	4.12	
44 Nitrobenzene	77	5.340	5.340	0.000	92	27855	4.00	4.00	
46 Isophorone	82	5.545	5.545	0.000	99	47544	4.00	4.02	
48 2-Nitrophenol	139	5.639	5.645	-0.006	95	14166	4.00	4.08	
49 2,4-Dimethylphenol	107	5.634	5.640	-0.006	93	25893	4.00	4.07	
50 Bis(2-chloroethoxy)methane	93	5.734	5.734	0.000	98	32100	4.00	4.13	
52 Benzoic acid	105	5.657	5.728	-0.071	92	32187	8.00	6.45	
53 2,4-Dichlorophenol	162	5.857	5.863	-0.006	95	21647	4.00	4.16	
54 1,2,4-Trichlorobenzene	180	5.951	5.951	0.000	94	22051	4.00	4.17	
57 Naphthalene	128	6.039	6.040	-0.001	98	79041	4.00	4.22	
58 4-Chloroaniline	127	6.081	6.081	0.000	96	35029	4.00	4.13	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	96	12544	4.00	4.35	
62 Caprolactam	55	6.416	6.428	-0.012	76	11099	4.00	3.33	
64 4-Chloro-3-methylphenol	107	6.528	6.528	0.000	97	22782	4.00	4.09	
65 2-Methylnaphthalene	142	6.716	6.716	0.000	93	49887	4.00	4.14	
67 1-Methylnaphthalene	142	6.822	6.822	0.000	93	46889	4.00	4.14	
68 Hexachlorocyclopentadiene	237	6.863	6.863	0.000	97	11449	4.00	3.75	
69 1,2,4,5-Tetrachlorobenzene	216	6.881	6.887	-0.005	97	20934	4.00	4.17	
70 2,4,6-Trichlorophenol	196	6.992	6.992	0.000	93	14208	4.00	4.07	
72 2,4,5-Trichlorophenol	196	7.028	7.028	0.000	92	15333	4.00	4.13	
74 1,1'-Biphenyl	154	7.175	7.181	-0.006	95	61753	4.00	4.22	
75 2-Chloronaphthalene	162	7.210	7.216	-0.006	97	45458	4.00	4.12	
77 2-Nitroaniline	65	7.316	7.316	0.000	84	15113	4.00	3.73	
79 Dimethyl phthalate	163	7.457	7.463	-0.006	98	60554	4.00	4.68	
80 1,3-Dinitrobenzene	168	7.522	7.528	-0.006	81	7547	4.00	3.29	
81 2,6-Dinitrotoluene	165	7.539	7.545	-0.006	91	11500	4.00	3.85	
82 Acenaphthylene	152	7.639	7.639	0.000	98	72357	4.00	4.04	
83 3-Nitroaniline	138	7.728	7.728	0.000	91	14261	4.00	3.69	
84 Acenaphthene	153	7.810	7.810	0.000	95	48195	4.00	4.17	
86 2,4-Dinitrophenol	184	7.828	7.828	0.000	77	11473	8.00	8.32	
87 4-Nitrophenol	109	7.857	7.869	-0.012	95	17189	8.00	8.25	
89 2,4-Dinitrotoluene	165	7.951	7.951	0.000	91	14418	4.00	3.68	
90 Dibenzofuran	168	7.975	7.981	-0.006	96	67976	4.00	4.18	
92 2,3,4,6-Tetrachlorophenol	232	8.092	8.092	0.000	78	11063	4.00	3.67	
94 Diethyl phthalate	149	8.151	8.157	-0.006	97	54858	4.00	4.17	
96 4-Chlorophenyl phenyl ethe	204	8.298	8.298	0.000	96	24578	4.00	4.17	
98 Fluorene	166	8.322	8.328	-0.006	96	54503	4.00	4.12	
99 4-Nitroaniline	138	8.345	8.357	-0.012	83	14556	4.00	3.85	
100 4,6-Dinitro-2-methylphenol	198	8.357	8.363	-0.006	77	14971	8.00	6.43	
102 N-Nitrosodiphenylamine	169	8.422	8.422	0.000	63	39816	4.00	4.09	
103 Azobenzene	77	8.463	8.463	0.000	99	60296	4.00	4.14	
104 1,2-Diphenylhydrazine	77	8.463	8.463	0.000	99	60296	4.04	4.19	
111 4-Bromophenyl phenyl ether	248	8.798	8.798	0.000	75	11781	4.00	3.82	
112 Hexachlorobenzene	284	8.880	8.881	0.000	90	13090	4.00	4.09	
116 Pentachlorophenol	266	9.075	9.080	-0.005	87	14903	8.00	7.08	
119 Phenanthrene	178	9.304	9.310	-0.006	98	76605	4.00	4.20	
120 Anthracene	178	9.357	9.357	0.000	98	77754	4.00	4.16	
122 Carbazole	167	9.522	9.522	0.000	96	77606	4.00	4.07	
123 Di-n-butyl phthalate	149	9.822	9.822	0.000	99	87433	4.00	3.90	
128 Fluoranthene	202	10.780	10.786	-0.006	97	79265	4.00	3.97	
131 Pyrene	202	11.151	11.157	-0.006	97	85070	4.00	4.03	
136 Famphur	218	12.192	12.198	-0.006	94	26998	4.00	4.06	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
137 Butyl benzyl phthalate	149	12.292	12.298	-0.006	98	34606	4.00	3.48	
140 3,3'-Dichlorobenzidine	252	13.627	13.639	-0.012	75	22416	4.00	3.59	
141 Benzo[a]anthracene	228	13.674	13.680	-0.006	97	73908	4.00	4.07	
142 Bis(2-ethylhexyl) phthalat	149	13.668	13.668	0.000	77	45285	4.00	3.48	
143 Chrysene	228	13.762	13.774	-0.012	97	67664	4.00	3.99	
144 Di-n-octyl phthalate	149	15.562	15.574	-0.012	99	61347	4.00	4.18	
146 Benzo[b]fluoranthene	252	16.715	16.727	-0.012	98	57364	4.00	3.65	
147 Benzo[k]fluoranthene	252	16.798	16.815	-0.017	98	60118	4.00	3.73	
148 Benzo[a]pyrene	252	17.715	17.727	-0.012	80	51493	4.00	3.41	
151 Indeno[1,2,3-cd]pyrene	276	21.097	21.127	-0.030	95	45380	4.00	3.52	
152 Dibenz(a,h)anthracene	278	21.180	21.203	-0.023	94	42195	4.00	3.38	
153 Benzo[g,h,i]perylene	276	21.874	21.897	-0.023	94	49079	4.00	3.62	

Reagents:

MS-HSLA004_00015 Amount Added: 200.00 Units: uL

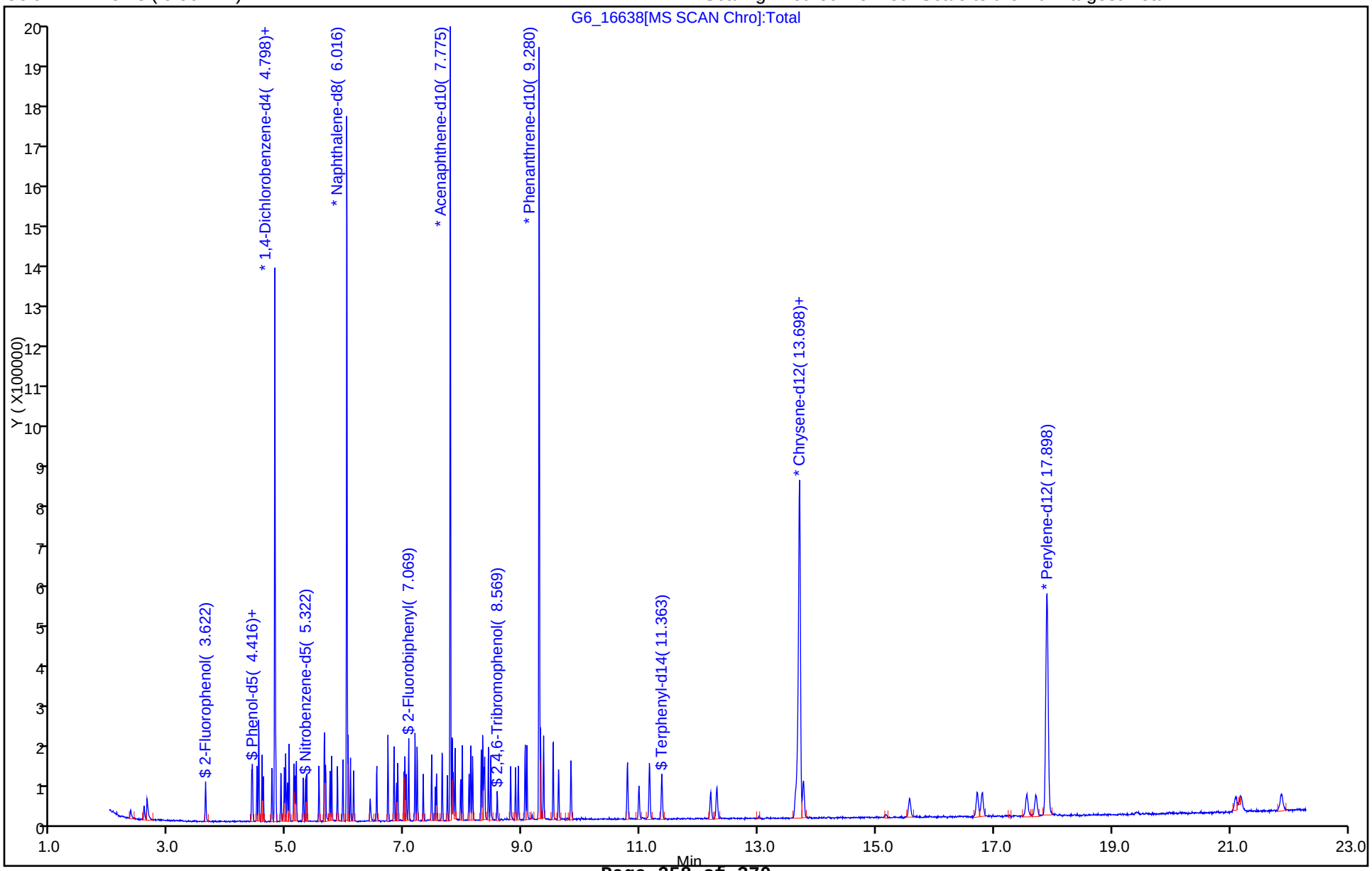
Report Date: 10-Apr-2015 12:09:39 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16638.D

Injection Date:	25-Feb-2015 12:19:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	STD004 HSL			Worklist Smp#:	4
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	3
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16639.D
Lims ID: STD010 HSL
Client ID:
Sample Type: IC Calib Level: 2
Inject. Date: 25-Feb-2015 12:46:30 ALS Bottle#: 4 Worklist Smp#: 5
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: STD010 HSL
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:40 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:02:27

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	96	184876	40.0	40.0	
* 2 Naphthalene-d8	136	6.016	6.022	-0.006	100	701243	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	92	384159	40.0	40.0	
* 4 Phenanthrene-d10	188	9.280	9.280	0.000	97	640383	40.0	40.0	
* 5 Chrysene-d12	240	13.698	13.710	-0.012	96	613004	40.0	40.0	
* 6 Perylene-d12	264	17.892	17.903	-0.011	95	530739	40.0	40.0	
$ 7 2-Fluorophenol	112	3.622	3.622	0.000	92	67803	10.0	10.1	
$ 8 Phenol-d5	99	4.404	4.410	-0.006	98	82351	10.0	9.77	
$ 9 Nitrobenzene-d5	82	5.316	5.322	-0.006	92	72669	10.0	10.1	
$ 10 2-Fluorobiphenyl	172	7.069	7.069	0.000	99	133834	10.0	10.3	
$ 11 2,4,6-Tribromophenol	330	8.569	8.575	-0.006	84	13361	10.0	9.25	
$ 12 Terphenyl-d14	244	11.363	11.369	-0.006	98	125136	10.0	10.3	
13 1,4-Dioxane	88	2.352	2.352	0.000	98	29946	10.0	10.6	
14 N-Nitrosodimethylamine	74	2.575	2.581	-0.006	91	43094	10.0	10.0	
15 Pyridine	79	2.622	2.622	0.000	92	73629	10.0	10.0	
23 Phenol	94	4.416	4.422	-0.006	98	85314	10.0	10.1	
24 Aniline	93	4.493	4.498	-0.005	98	101304	10.0	9.81	
25 Bis(2-chloroethyl)ether	93	4.522	4.528	-0.006	90	69073	10.0	10.4	
26 2-Chlorophenol	128	4.604	4.604	0.000	97	70338	10.0	10.2	
27 1,3-Dichlorobenzene	146	4.745	4.751	-0.006	97	71887	10.0	10.3	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	96	72168	10.0	10.2	
29 Benzyl alcohol	108	4.898	4.904	-0.006	92	41657	10.0	9.59	
30 1,2-Dichlorobenzene	146	4.963	4.963	0.000	96	69610	10.0	10.2	
31 2-Methylphenol	108	4.981	4.981	0.000	94	63348	10.0	10.1	
32 2,2'-oxybis[1-chloropropan	45	5.010	5.010	0.000	93	99866	10.0	10.2	
38 3 & 4 Methylphenol	108	5.122	5.128	-0.006	98	63816	10.0	9.83	
39 3-Methylphenol	108	5.122	5.128	-0.006	98	63816	10.0	9.83	
40 4-Methylphenol	108	5.122	5.128	-0.006	94	63816	10.0	9.83	
41 N-Nitrosodi-n-propylamine	70	5.140	5.145	-0.005	91	45947	10.0	9.75	
42 Acetophenone	105	5.163	5.163	0.000	96	93583	10.0	10.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.281	5.281	0.000	96	30699	10.0	10.1	
44 Nitrobenzene	77	5.340	5.340	0.000	91	71302	10.0	10.3	
46 Isophorone	82	5.545	5.545	0.000	99	115206	10.0	9.79	
48 2-Nitrophenol	139	5.640	5.645	-0.005	90	35195	10.0	10.2	
49 2,4-Dimethylphenol	107	5.634	5.640	-0.006	93	66475	10.0	10.5	
50 Bis(2-chloroethoxy)methane	93	5.734	5.734	0.000	98	80123	10.0	10.4	
52 Benzoic acid	105	5.669	5.728	-0.059	88	89023	20.0	17.9	
53 2,4-Dichlorophenol	162	5.857	5.863	-0.006	97	51694	10.0	9.98	
54 1,2,4-Trichlorobenzene	180	5.951	5.951	0.000	94	53749	10.0	10.2	
57 Naphthalene	128	6.040	6.040	0.000	97	192392	10.0	10.3	
58 4-Chloroaniline	127	6.081	6.081	0.000	96	85639	10.0	10.2	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	98	29013	10.0	10.1	
62 Caprolactam	55	6.410	6.428	-0.018	76	31448	10.0	9.48	
64 4-Chloro-3-methylphenol	107	6.528	6.528	0.000	98	55761	10.0	10.1	
65 2-Methylnaphthalene	142	6.716	6.716	0.000	93	124165	10.0	10.4	
67 1-Methylnaphthalene	142	6.822	6.822	0.000	94	117053	10.0	10.4	
68 Hexachlorocyclopentadiene	237	6.863	6.863	0.000	97	28473	10.0	9.43	
69 1,2,4,5-Tetrachlorobenzene	216	6.881	6.887	-0.005	97	50324	10.0	10.1	
70 2,4,6-Trichlorophenol	196	6.992	6.992	0.000	94	34170	10.0	9.91	
72 2,4,5-Trichlorophenol	196	7.028	7.028	0.000	94	36506	10.0	9.93	
74 1,1'-Biphenyl	154	7.175	7.181	-0.006	95	149463	10.0	10.3	
75 2-Chloronaphthalene	162	7.210	7.216	-0.006	98	113058	10.0	10.4	
77 2-Nitroaniline	65	7.316	7.316	0.000	85	39287	10.0	9.81	
79 Dimethyl phthalate	163	7.457	7.463	-0.006	98	130115	10.0	10.2	
80 1,3-Dinitrobenzene	168	7.522	7.528	-0.006	81	20263	10.0	8.93	
81 2,6-Dinitrotoluene	165	7.539	7.545	-0.006	92	28732	10.0	9.72	
82 Acenaphthylene	152	7.634	7.639	-0.005	98	178896	10.0	10.1	
83 3-Nitroaniline	138	7.728	7.728	0.000	94	37522	10.0	9.82	
84 Acenaphthene	153	7.810	7.810	0.000	95	117997	10.0	10.3	
86 2,4-Dinitrophenol	184	7.822	7.828	-0.006	81	32905	20.0	18.2	
87 4-Nitrophenol	109	7.857	7.869	-0.012	96	40880	20.0	19.8	
89 2,4-Dinitrotoluene	165	7.951	7.951	0.000	91	37208	10.0	9.60	
90 Dibenzofuran	168	7.975	7.981	-0.006	97	164638	10.0	10.2	
92 2,3,4,6-Tetrachlorophenol	232	8.092	8.092	0.000	75	29631	10.0	9.95	
94 Diethyl phthalate	149	8.151	8.157	-0.006	97	134263	10.0	10.3	
96 4-Chlorophenyl phenyl ethe	204	8.298	8.298	0.000	95	60258	10.0	10.3	
98 Fluorene	166	8.322	8.328	-0.006	96	136339	10.0	10.4	
99 4-Nitroaniline	138	8.345	8.357	-0.012	84	37301	10.0	9.98	
100 4,6-Dinitro-2-methylphenol	198	8.357	8.363	-0.006	78	43031	20.0	19.0	
102 N-Nitrosodiphenylamine	169	8.422	8.422	0.000	62	98599	10.0	10.4	
103 Azobenzene	77	8.457	8.463	-0.006	99	144575	10.0	10.0	
104 1,2-Diphenylhydrazine	77	8.457	8.463	-0.006	99	144575	10.1	10.2	
111 4-Bromophenyl phenyl ether	248	8.798	8.798	0.000	73	31038	10.0	10.3	
112 Hexachlorobenzene	284	8.881	8.881	0.000	91	31646	10.0	10.2	
116 Pentachlorophenol	266	9.075	9.080	-0.005	87	38486	20.0	18.8	
119 Phenanthrene	178	9.304	9.310	-0.006	98	185797	10.0	10.5	
120 Anthracene	178	9.357	9.357	0.000	98	188375	10.0	10.4	
122 Carbazole	167	9.516	9.522	-0.006	96	190570	10.0	10.3	
123 Di-n-butyl phthalate	149	9.822	9.822	0.000	100	217656	10.0	9.97	
128 Fluoranthene	202	10.775	10.786	-0.012	99	199608	10.0	10.3	
131 Pyrene	202	11.151	11.157	-0.006	97	206008	10.0	10.1	
136 Famphur	218	12.192	12.198	-0.006	97	66995	10.0	10.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
137 Butyl benzyl phthalate	149	12.292	12.298	-0.006	97	91001	10.0	9.42	
140 3,3'-Dichlorobenzidine	252	13.627	13.639	-0.012	75	57301	10.0	9.46	
141 Benzo[a]anthracene	228	13.663	13.680	-0.017	100	177675	10.0	10.1	
142 Bis(2-ethylhexyl) phthalat	149	13.663	13.668	-0.005	96	118615	10.0	9.40	
143 Chrysene	228	13.763	13.774	-0.011	97	164066	10.0	9.97	
144 Di-n-octyl phthalate	149	15.562	15.574	-0.012	99	176570	10.0	9.28	
146 Benzo[b]fluoranthene	252	16.709	16.727	-0.018	99	151541	10.0	9.77	
147 Benzo[k]fluoranthene	252	16.798	16.815	-0.017	99	145931	10.0	9.16	
148 Benzo[a]pyrene	252	17.709	17.727	-0.018	80	138204	10.0	9.27	
151 Indeno[1,2,3-cd]pyrene	276	21.097	21.127	-0.030	97	116621	10.0	9.32	
152 Dibenz(a,h)anthracene	278	21.186	21.203	-0.017	93	107166	10.0	8.70	
153 Benzo[g,h,i]perylene	276	21.874	21.897	-0.023	94	127555	10.0	9.51	

Reagents:

MS-HSLA010_00015 Amount Added: 200.00 Units: uL

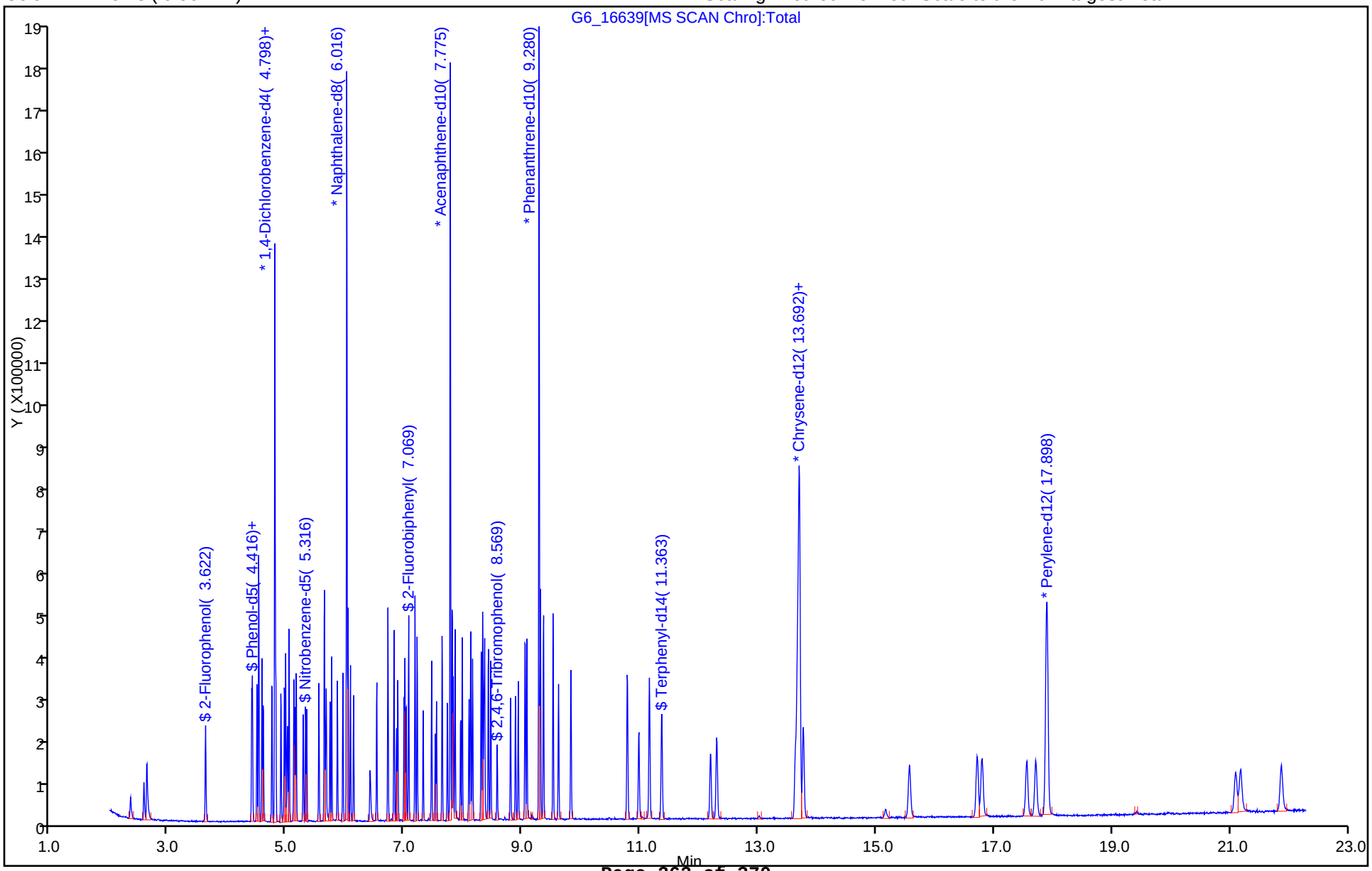
Report Date: 10-Apr-2015 12:09:40 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16639.D

Injection Date:	25-Feb-2015 12:46:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	STD010 HSL			Worklist Smp#:	5
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	4
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16640.D
Lims ID: STD020 HSL
Client ID:
Sample Type: IC Calib Level: 3
Inject. Date: 25-Feb-2015 13:12:30 ALS Bottle#: 5 Worklist Smp#: 6
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: STD020 HSL
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:41 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:04:20

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	96	188896	40.0	40.0	
* 2 Naphthalene-d8	136	6.016	6.022	-0.006	100	722037	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	92	390190	40.0	40.0	
* 4 Phenanthrene-d10	188	9.280	9.280	0.000	97	655697	40.0	40.0	
* 5 Chrysene-d12	240	13.698	13.710	-0.012	96	632287	40.0	40.0	
* 6 Perylene-d12	264	17.892	17.903	-0.011	96	537132	40.0	40.0	
$ 7 2-Fluorophenol	112	3.622	3.622	0.000	93	136584	20.0	19.9	
$ 8 Phenol-d5	99	4.404	4.410	-0.006	99	173137	20.0	20.1	
$ 9 Nitrobenzene-d5	82	5.316	5.322	-0.006	92	146288	20.0	19.8	
$ 10 2-Fluorobiphenyl	172	7.069	7.069	0.000	100	271127	20.0	20.6	
$ 11 2,4,6-Tribromophenol	330	8.569	8.575	-0.006	84	28834	20.0	19.7	
$ 12 Terphenyl-d14	244	11.363	11.369	-0.006	98	250222	20.0	19.9	
13 1,4-Dioxane	88	2.351	2.352	-0.001	95	59534	20.0	20.5	
14 N-Nitrosodimethylamine	74	2.575	2.581	-0.006	89	85710	20.0	19.5	
15 Pyridine	79	2.622	2.622	0.000	90	147856	20.0	19.7	
23 Phenol	94	4.416	4.422	-0.006	98	172136	20.0	19.9	
24 Aniline	93	4.493	4.498	-0.005	97	206891	20.0	19.6	
25 Bis(2-chloroethyl)ether	93	4.522	4.528	-0.006	88	137208	20.0	20.2	
26 2-Chlorophenol	128	4.604	4.604	0.000	97	139729	20.0	19.8	
27 1,3-Dichlorobenzene	146	4.745	4.751	-0.006	97	142995	20.0	20.0	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	94	146967	20.0	20.3	
29 Benzyl alcohol	108	4.904	4.904	0.000	92	86805	20.0	19.6	
30 1,2-Dichlorobenzene	146	4.957	4.963	-0.006	95	140215	20.0	20.2	
31 2-Methylphenol	108	4.981	4.981	0.000	96	127434	20.0	19.9	
32 2,2'-oxybis[1-chloropropan	45	5.010	5.010	0.000	92	199591	20.0	20.0	
38 3 & 4 Methylphenol	108	5.122	5.128	-0.006	97	133429	20.0	20.1	
39 3-Methylphenol	108	5.122	5.128	-0.006	97	133429	20.0	20.1	
40 4-Methylphenol	108	5.122	5.128	-0.006	92	133429	20.0	20.1	
41 N-Nitrosodi-n-propylamine	70	5.140	5.145	-0.005	90	96259	20.0	20.0	
42 Acetophenone	105	5.163	5.163	0.000	96	185111	20.0	19.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.281	5.281	0.000	97	61931	20.0	20.0	
44 Nitrobenzene	77	5.340	5.340	0.000	91	141833	20.0	19.9	
46 Isophorone	82	5.545	5.545	0.000	99	240037	20.0	19.8	
48 2-Nitrophenol	139	5.640	5.645	-0.005	92	71833	20.0	20.2	
49 2,4-Dimethylphenol	107	5.634	5.640	-0.006	92	131383	20.0	20.2	
50 Bis(2-chloroethoxy)methane	93	5.734	5.734	0.000	98	160771	20.0	20.2	
52 Benzoic acid	105	5.681	5.728	-0.047	90	194250	40.0	38.0	
53 2,4-Dichlorophenol	162	5.857	5.863	-0.006	97	106817	20.0	20.0	
54 1,2,4-Trichlorobenzene	180	5.951	5.951	0.000	94	109397	20.0	20.2	
57 Naphthalene	128	6.039	6.040	-0.001	97	390163	20.0	20.3	
58 4-Chloroaniline	127	6.081	6.081	0.000	95	172259	20.0	19.8	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	97	58938	20.0	20.0	
62 Caprolactam	55	6.410	6.428	-0.018	77	65538	20.0	19.2	M
64 4-Chloro-3-methylphenol	107	6.528	6.528	0.000	98	112366	20.0	19.7	
65 2-Methylnaphthalene	142	6.716	6.716	0.000	94	249498	20.0	20.2	
67 1-Methylnaphthalene	142	6.822	6.822	0.000	94	236260	20.0	20.4	
68 Hexachlorocyclopentadiene	237	6.863	6.863	0.000	97	60610	20.0	19.8	
69 1,2,4,5-Tetrachlorobenzene	216	6.881	6.887	-0.005	98	104495	20.0	20.3	
70 2,4,6-Trichlorophenol	196	6.992	6.992	0.000	94	70312	20.0	20.1	
72 2,4,5-Trichlorophenol	196	7.028	7.028	0.000	94	75074	20.0	20.1	
74 1,1'-Biphenyl	154	7.175	7.181	-0.006	95	302324	20.0	20.6	
75 2-Chloronaphthalene	162	7.210	7.216	-0.006	98	227907	20.0	20.6	
77 2-Nitroaniline	65	7.316	7.316	0.000	85	82700	20.0	20.3	
79 Dimethyl phthalate	163	7.457	7.463	-0.006	97	257927	20.0	19.9	
80 1,3-Dinitrobenzene	168	7.522	7.528	-0.006	83	45571	20.0	19.8	
81 2,6-Dinitrotoluene	165	7.539	7.545	-0.006	94	62625	20.0	20.9	
82 Acenaphthylene	152	7.634	7.639	-0.005	99	362680	20.0	20.2	
83 3-Nitroaniline	138	7.728	7.728	0.000	95	77435	20.0	19.9	
84 Acenaphthene	153	7.810	7.810	0.000	95	238340	20.0	20.5	
86 2,4-Dinitrophenol	184	7.822	7.828	-0.006	82	78117	40.0	38.4	
87 4-Nitrophenol	109	7.857	7.869	-0.012	96	82891	40.0	39.6	
89 2,4-Dinitrotoluene	165	7.951	7.951	0.000	91	79927	20.0	20.3	
90 Dibenzofuran	168	7.975	7.981	-0.006	97	334583	20.0	20.5	
92 2,3,4,6-Tetrachlorophenol	232	8.092	8.092	0.000	76	59694	20.0	19.7	
94 Diethyl phthalate	149	8.151	8.157	-0.006	97	273258	20.0	20.7	
96 4-Chlorophenyl phenyl ethe	204	8.298	8.298	0.000	97	121088	20.0	20.5	
98 Fluorene	166	8.322	8.328	-0.006	97	274499	20.0	20.6	
99 4-Nitroaniline	138	8.345	8.357	-0.012	84	76419	20.0	20.1	
100 4,6-Dinitro-2-methylphenol	198	8.357	8.363	-0.006	80	93754	40.0	40.4	
102 N-Nitrosodiphenylamine	169	8.422	8.422	0.000	63	201342	20.0	20.7	
103 Azobenzene	77	8.463	8.463	0.000	99	295740	20.0	20.2	
104 1,2-Diphenylhydrazine	77	8.463	8.463	0.000	100	295740	20.2	20.5	
111 4-Bromophenyl phenyl ether	248	8.798	8.798	0.000	72	62227	20.0	20.3	
112 Hexachlorobenzene	284	8.880	8.881	0.000	91	65622	20.0	20.6	
116 Pentachlorophenol	266	9.075	9.080	-0.005	88	84234	40.0	40.2	
119 Phenanthrene	178	9.304	9.310	-0.006	98	370680	20.0	20.4	
120 Anthracene	178	9.357	9.357	0.000	98	381794	20.0	20.5	
122 Carbazole	167	9.516	9.522	-0.006	96	392545	20.0	20.7	
123 Di-n-butyl phthalate	149	9.822	9.822	0.000	100	453491	20.0	20.3	
128 Fluoranthene	202	10.780	10.786	-0.006	98	405997	20.0	20.4	
131 Pyrene	202	11.151	11.157	-0.006	97	418438	20.0	19.8	
136 Famphur	218	12.192	12.198	-0.006	96	140106	20.0	21.0	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
137 Butyl benzyl phthalate	149	12.292	12.298	-0.006	98	193326	20.0	19.4	
140 3,3'-Dichlorobenzidine	252	13.627	13.639	-0.012	75	120010	20.0	19.2	
141 Benzo[a]anthracene	228	13.668	13.680	-0.012	99	361571	20.0	19.9	
142 Bis(2-ethylhexyl) phthalat	149	13.663	13.668	-0.005	96	252271	20.0	19.4	
143 Chrysene	228	13.763	13.774	-0.011	98	336779	20.0	19.8	
144 Di-n-octyl phthalate	149	15.562	15.574	-0.012	99	392967	20.0	18.2	
146 Benzo[b]fluoranthene	252	16.715	16.727	-0.012	98	312315	20.0	19.9	
147 Benzo[k]fluoranthene	252	16.798	16.815	-0.017	98	313777	20.0	19.5	
148 Benzo[a]pyrene	252	17.703	17.727	-0.024	80	293226	20.0	19.4	
151 Indeno[1,2,3-cd]pyrene	276	21.103	21.127	-0.024	97	244980	20.0	19.0	
152 Dibenz(a,h)anthracene	278	21.180	21.203	-0.023	94	242009	20.0	19.4	
153 Benzo[g,h,i]perylene	276	21.880	21.897	-0.017	94	261176	20.0	19.2	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

MS-HSLA020_00015 Amount Added: 200.00 Units: uL

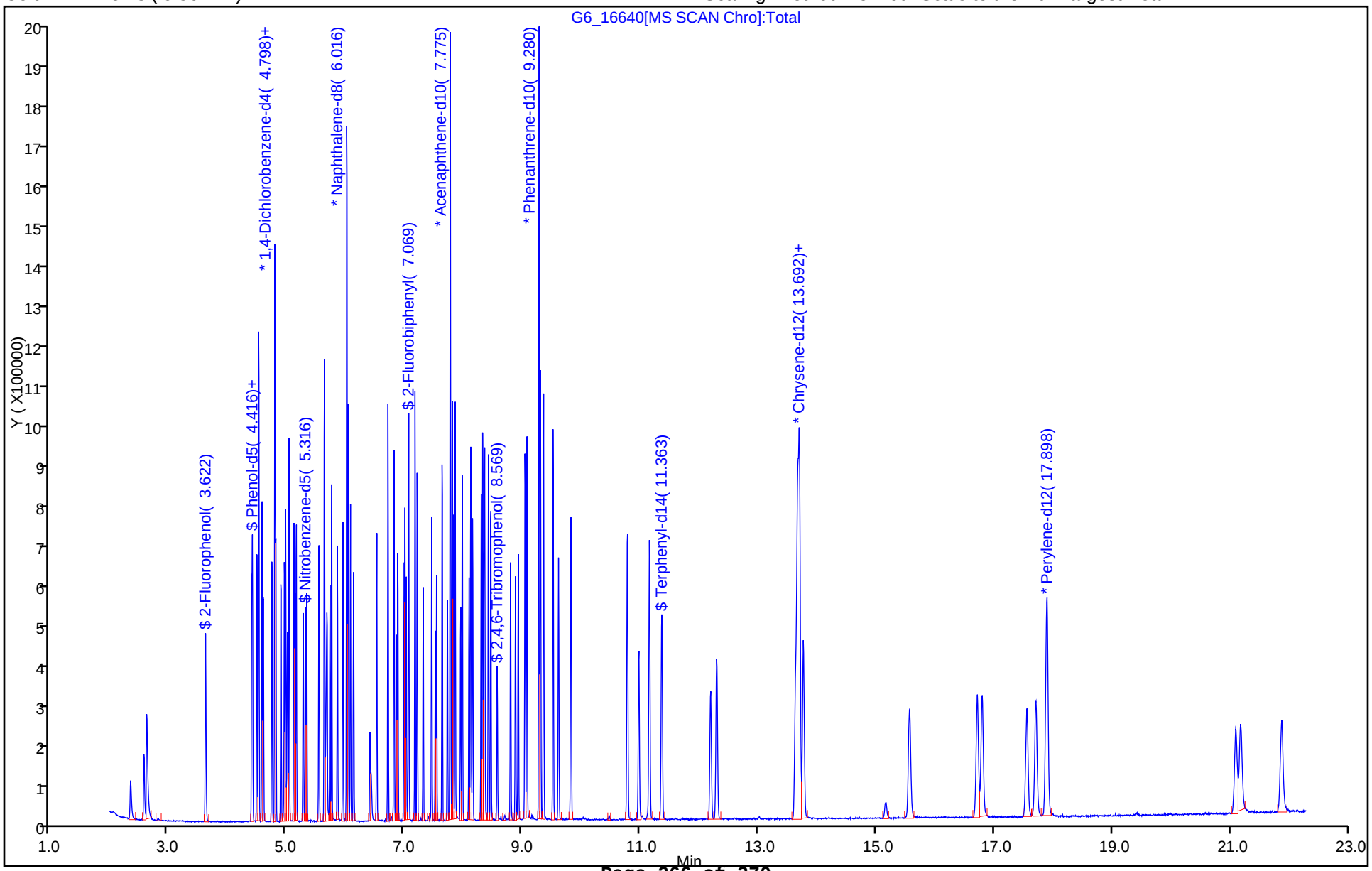
Report Date: 10-Apr-2015 12:09:41 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16640.D

Injection Date:	25-Feb-2015 13:12:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	STD020 HSL			Worklist Smp#:	6
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	5
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



Manual Integration/User Assign Peak Report

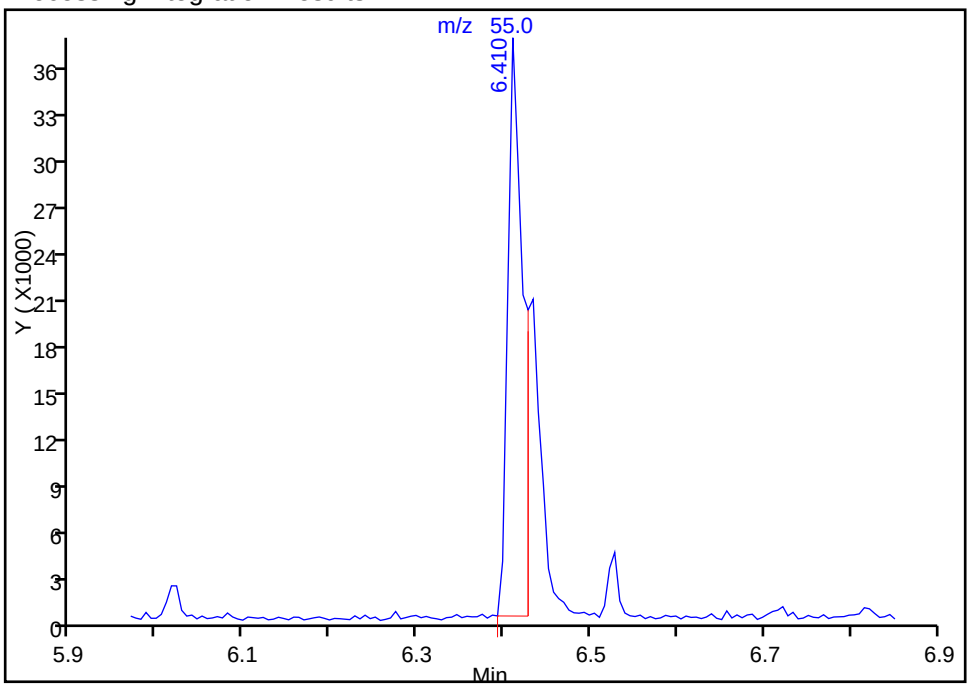
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16640.D				
Injection Date:	25-Feb-2015 13:12:30	Instrument ID:	SMS_G6		
Lims ID:	STD020 HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	5	Worklist Smp#:	6
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

62 Caprolactam, CAS: 105-60-2

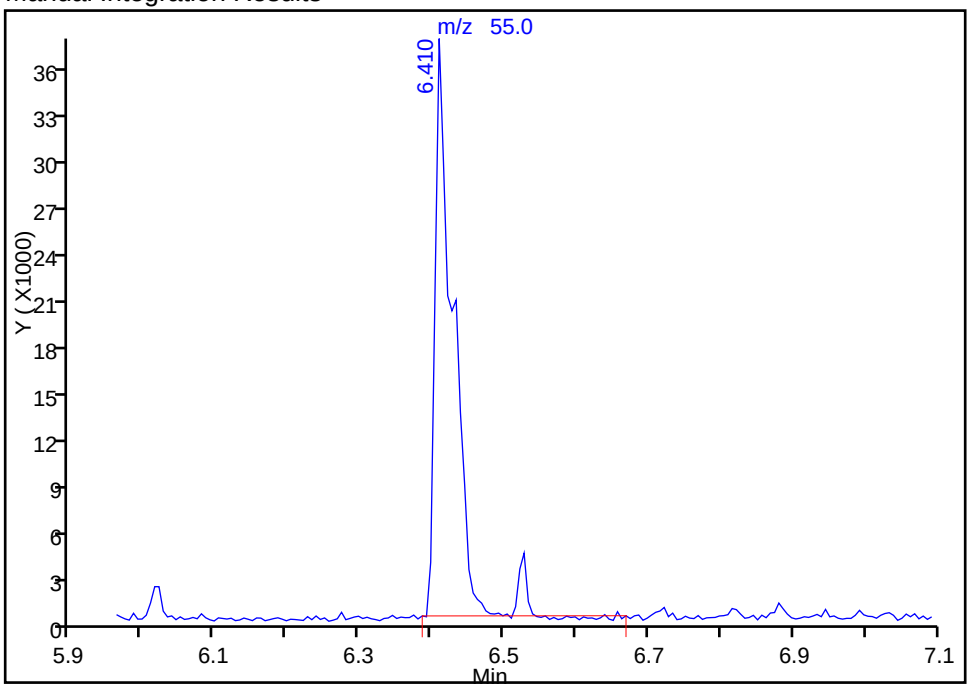
Processing Integration Results

RT: 6.41
Area: 46259
Amount: 18.733378
Amount Units: ug/ml



Manual Integration Results

RT: 6.41
Area: 65538
Amount: 19.189269
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:04:20
Audit Action: Manually Integrated
Audit Reason: Split Peak

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16641.D
Lims ID: STD050 HSL
Client ID:
Sample Type: IC Calib Level: 4
Inject. Date: 25-Feb-2015 13:39:30 ALS Bottle#: 6 Worklist Smp#: 7
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: STD050 HSL
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:42 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:09:09

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	96	188580	40.0	40.0	
* 2 Naphthalene-d8	136	6.016	6.022	-0.006	100	730834	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	93	393471	40.0	40.0	
* 4 Phenanthrene-d10	188	9.280	9.280	0.000	97	661919	40.0	40.0	
* 5 Chrysene-d12	240	13.704	13.710	-0.006	96	630959	40.0	40.0	
* 6 Perylene-d12	264	17.898	17.903	-0.005	95	543830	40.0	40.0	
$ 7 2-Fluorophenol	112	3.622	3.622	0.000	94	342921	50.0	50.0	
$ 8 Phenol-d5	99	4.404	4.410	-0.006	99	432891	50.0	50.4	
$ 9 Nitrobenzene-d5	82	5.322	5.322	0.000	91	377126	50.0	50.3	
$ 10 2-Fluorobiphenyl	172	7.069	7.069	0.000	99	677613	50.0	50.9	
$ 11 2,4,6-Tribromophenol	330	8.575	8.575	0.000	91	75692	50.0	51.2	
$ 12 Terphenyl-d14	244	11.369	11.369	0.000	98	635940	50.0	50.7	
13 1,4-Dioxane	88	2.351	2.352	-0.001	98	142371	50.0	49.2	
14 N-Nitrosodimethylamine	74	2.575	2.581	-0.006	90	222746	50.0	50.7	
15 Pyridine	79	2.622	2.622	0.000	90	375517	50.0	50.1	
23 Phenol	94	4.422	4.422	0.000	99	436251	50.0	50.6	
24 Aniline	93	4.498	4.498	0.000	98	529556	50.0	50.3	
25 Bis(2-chloroethyl)ether	93	4.528	4.528	0.000	89	342205	50.0	50.6	
26 2-Chlorophenol	128	4.604	4.604	0.000	97	356590	50.0	50.7	
27 1,3-Dichlorobenzene	146	4.751	4.751	0.000	97	360797	50.0	50.5	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	93	366601	50.0	50.8	
29 Benzyl alcohol	108	4.904	4.904	0.000	92	224224	50.0	50.6	
30 1,2-Dichlorobenzene	146	4.963	4.963	0.000	96	348029	50.0	50.2	
31 2-Methylphenol	108	4.981	4.981	0.000	96	325292	50.0	50.9	
32 2,2'-oxybis[1-chloropropan	45	5.010	5.010	0.000	93	507190	50.0	50.9	
38 3 & 4 Methylphenol	108	5.122	5.128	-0.006	99	334564	50.0	50.5	
39 3-Methylphenol	108	5.122	5.128	-0.006	99	334564	50.0	50.5	
40 4-Methylphenol	108	5.122	5.128	-0.006	93	334564	50.0	50.5	
41 N-Nitrosodi-n-propylamine	70	5.140	5.145	-0.005	90	246852	50.0	51.4	
42 Acetophenone	105	5.163	5.163	0.000	97	472247	50.0	50.7	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.281	5.281	0.000	96	157602	50.0	50.9	
44 Nitrobenzene	77	5.340	5.340	0.000	90	365089	50.0	50.6	
46 Isophorone	82	5.545	5.545	0.000	99	626824	50.0	51.1	
48 2-Nitrophenol	139	5.639	5.645	-0.006	95	184945	50.0	51.4	
49 2,4-Dimethylphenol	107	5.639	5.640	-0.001	93	335864	50.0	50.9	
50 Bis(2-chloroethoxy)methane	93	5.734	5.734	0.000	98	407948	50.0	50.6	
52 Benzoic acid	105	5.710	5.728	-0.018	91	519854	100.0	100.5	
53 2,4-Dichlorophenol	162	5.863	5.863	0.000	96	269879	50.0	50.0	
54 1,2,4-Trichlorobenzene	180	5.951	5.951	0.000	93	279647	50.0	51.0	
57 Naphthalene	128	6.039	6.040	-0.001	98	977042	50.0	50.3	
58 4-Chloroaniline	127	6.081	6.081	0.000	96	443432	50.0	50.4	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	98	147770	50.0	49.4	
62 Caprolactam	55	6.416	6.428	-0.012	77	171197	50.0	49.5	M
64 4-Chloro-3-methylphenol	107	6.528	6.528	0.000	97	289585	50.0	50.1	
65 2-Methylnaphthalene	142	6.716	6.716	0.000	93	636017	50.0	50.9	
67 1-Methylnaphthalene	142	6.822	6.822	0.000	94	599383	50.0	51.0	
68 Hexachlorocyclopentadiene	237	6.863	6.863	0.000	97	159056	50.0	51.5	
69 1,2,4,5-Tetrachlorobenzene	216	6.886	6.887	0.000	98	266816	50.0	51.3	
70 2,4,6-Trichlorophenol	196	6.992	6.992	0.000	94	180797	50.0	51.2	
72 2,4,5-Trichlorophenol	196	7.028	7.028	0.000	92	189993	50.0	50.5	
74 1,1'-Biphenyl	154	7.175	7.181	-0.006	96	754596	50.0	51.0	
75 2-Chloronaphthalene	162	7.210	7.216	-0.006	98	569839	50.0	51.0	
77 2-Nitroaniline	65	7.316	7.316	0.000	84	210036	50.0	51.2	
79 Dimethyl phthalate	163	7.457	7.463	-0.006	98	649391	50.0	49.6	
80 1,3-Dinitrobenzene	168	7.522	7.528	-0.006	83	119893	50.0	51.6	
81 2,6-Dinitrotoluene	165	7.545	7.545	0.000	95	154313	50.0	51.0	
82 Acenaphthylene	152	7.639	7.639	0.000	99	931810	50.0	51.4	
83 3-Nitroaniline	138	7.728	7.728	0.000	94	204409	50.0	52.2	
84 Acenaphthene	153	7.810	7.810	0.000	95	602051	50.0	51.4	
86 2,4-Dinitrophenol	184	7.828	7.828	0.000	81	217993	100.0	100.7	
87 4-Nitrophenol	109	7.863	7.869	-0.006	95	213891	100.0	101.3	
89 2,4-Dinitrotoluene	165	7.951	7.951	0.000	91	206301	50.0	52.0	
90 Dibenzofuran	168	7.981	7.981	-0.001	97	844896	50.0	51.2	
92 2,3,4,6-Tetrachlorophenol	232	8.092	8.092	0.000	76	157611	50.0	51.7	
94 Diethyl phthalate	149	8.157	8.157	0.000	98	680249	50.0	51.0	
96 4-Chlorophenyl phenyl ethe	204	8.298	8.298	0.000	96	304597	50.0	51.1	
98 Fluorene	166	8.328	8.328	0.000	96	692688	50.0	51.6	
99 4-Nitroaniline	138	8.351	8.357	-0.006	85	197830	50.0	51.7	
100 4,6-Dinitro-2-methylphenol	198	8.363	8.363	0.000	86	249423	100.0	106.5	
102 N-Nitrosodiphenylamine	169	8.422	8.422	0.000	62	498390	50.0	50.8	
103 Azobenzene	77	8.463	8.463	0.000	99	755331	50.0	51.2	
104 1,2-Diphenylhydrazine	77	8.463	8.463	0.000	99	755331	50.5	51.8	
111 4-Bromophenyl phenyl ether	248	8.798	8.798	0.000	72	159165	50.0	51.3	
112 Hexachlorobenzene	284	8.880	8.881	0.000	92	164764	50.0	51.1	
116 Pentachlorophenol	266	9.075	9.080	-0.005	89	224380	100.0	106.0	
119 Phenanthrene	178	9.310	9.310	0.000	98	941975	50.0	51.3	
120 Anthracene	178	9.357	9.357	0.000	98	959498	50.0	51.0	
122 Carbazole	167	9.522	9.522	0.000	96	979659	50.0	51.1	
123 Di-n-butyl phthalate	149	9.822	9.822	0.000	100	1175500	50.0	52.1	
128 Fluoranthene	202	10.780	10.786	-0.006	98	1021800	50.0	50.8	
131 Pyrene	202	11.157	11.157	0.000	97	1068582	50.0	50.7	
136 Famphur	218	12.198	12.198	0.000	97	359540	50.0	54.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
137 Butyl benzyl phthalate	149	12.298	12.298	0.000	98	510447	50.0	51.3	
140 3,3'-Dichlorobenzidine	252	13.633	13.639	-0.006	75	324147	50.0	52.0	
141 Benzo[a]anthracene	228	13.674	13.680	-0.006	99	919851	50.0	50.7	
142 Bis(2-ethylhexyl) phthalat	149	13.662	13.668	-0.006	97	679050	50.0	52.3	
143 Chrysene	228	13.768	13.774	-0.006	98	858164	50.0	50.7	
144 Di-n-octyl phthalate	149	15.568	15.574	-0.006	99	1119188	50.0	49.0	
146 Benzo[b]fluoranthene	252	16.721	16.727	-0.006	99	791973	50.0	49.9	
147 Benzo[k]fluoranthene	252	16.803	16.815	-0.012	98	839735	50.0	51.5	
148 Benzo[a]pyrene	252	17.721	17.727	-0.006	80	792402	50.0	51.9	
151 Indeno[1,2,3-cd]pyrene	276	21.115	21.127	-0.012	97	654992	50.0	50.8	M
152 Dibenz(a,h)anthracene	278	21.197	21.203	-0.006	94	660134	50.0	52.3	
153 Benzo[g,h,i]perylene	276	21.891	21.897	-0.006	96	698128	50.0	50.8	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

MS-HSLA050_00015 Amount Added: 200.00 Units: uL

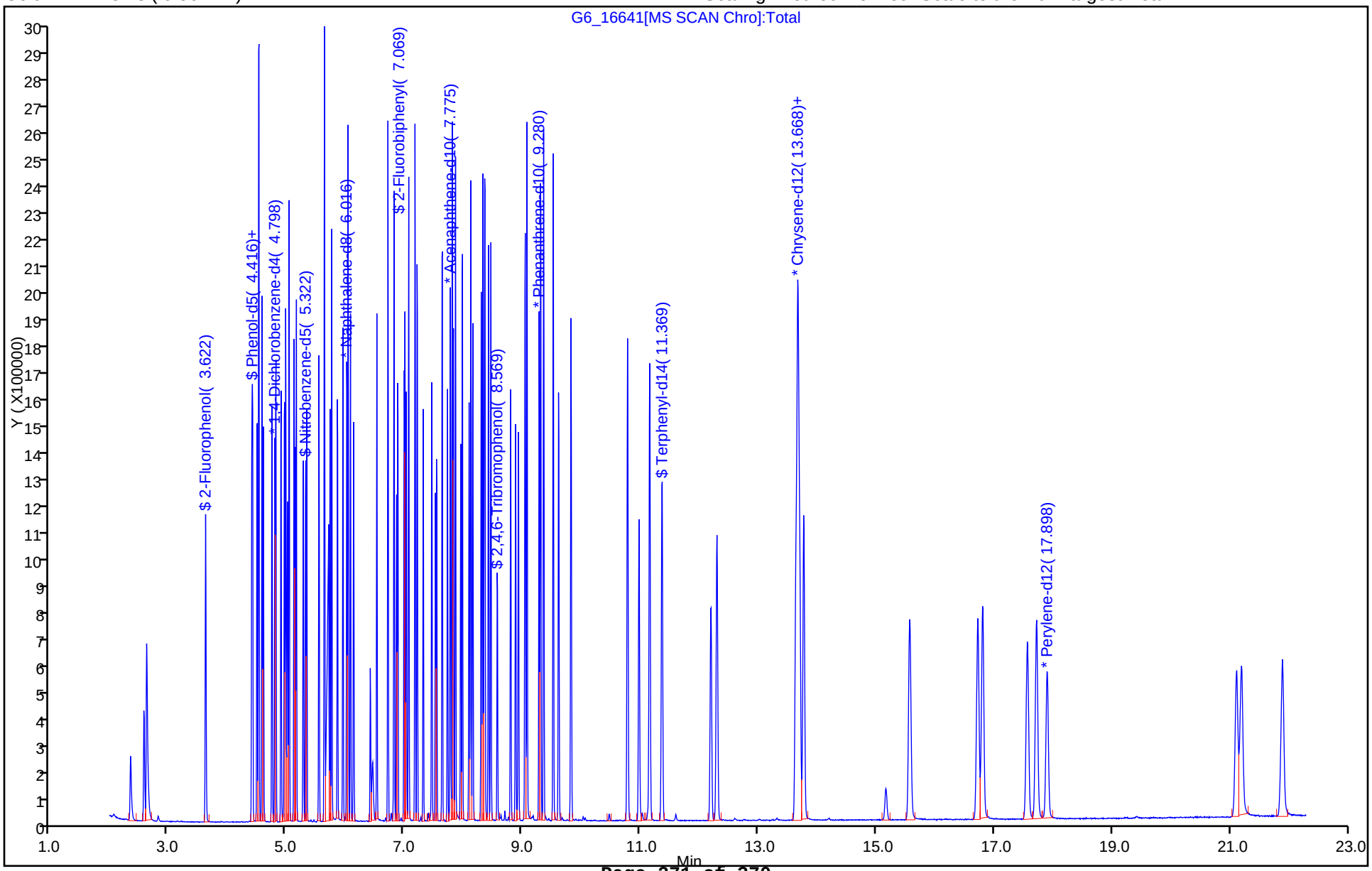
Report Date: 10-Apr-2015 12:09:42 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16641.D

Injection Date:	25-Feb-2015 13:39:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	STD050 HSL			Worklist Smp#:	7
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	6
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



Manual Integration/User Assign Peak Report

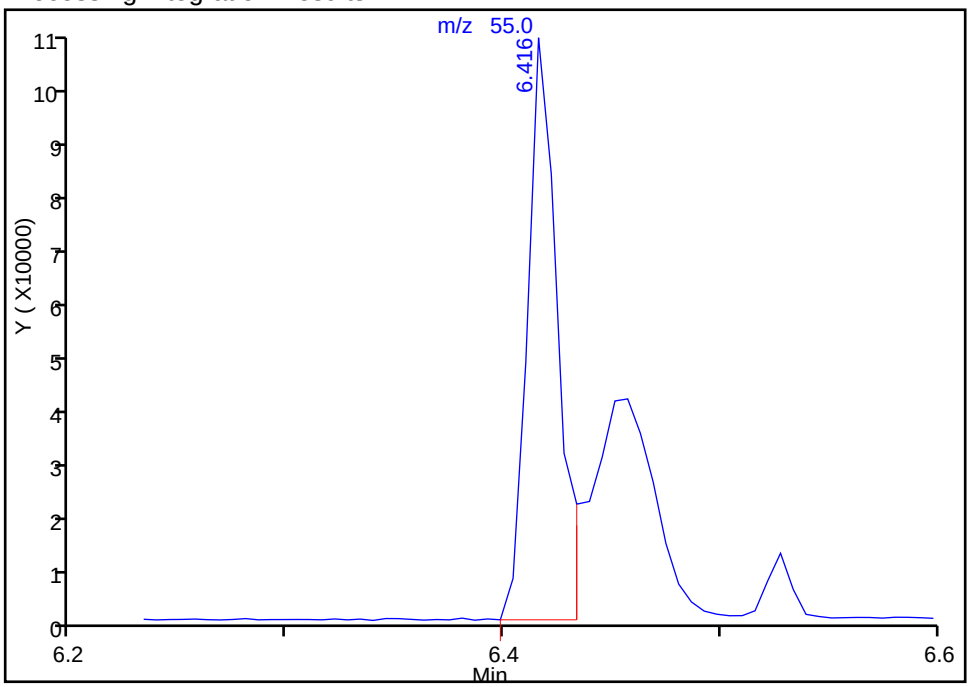
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16641.D				
Injection Date:	25-Feb-2015 13:39:30	Instrument ID:	SMS_G6		
Lims ID:	STD050 HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	6	Worklist Smp#:	7
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

62 Caprolactam, CAS: 105-60-2

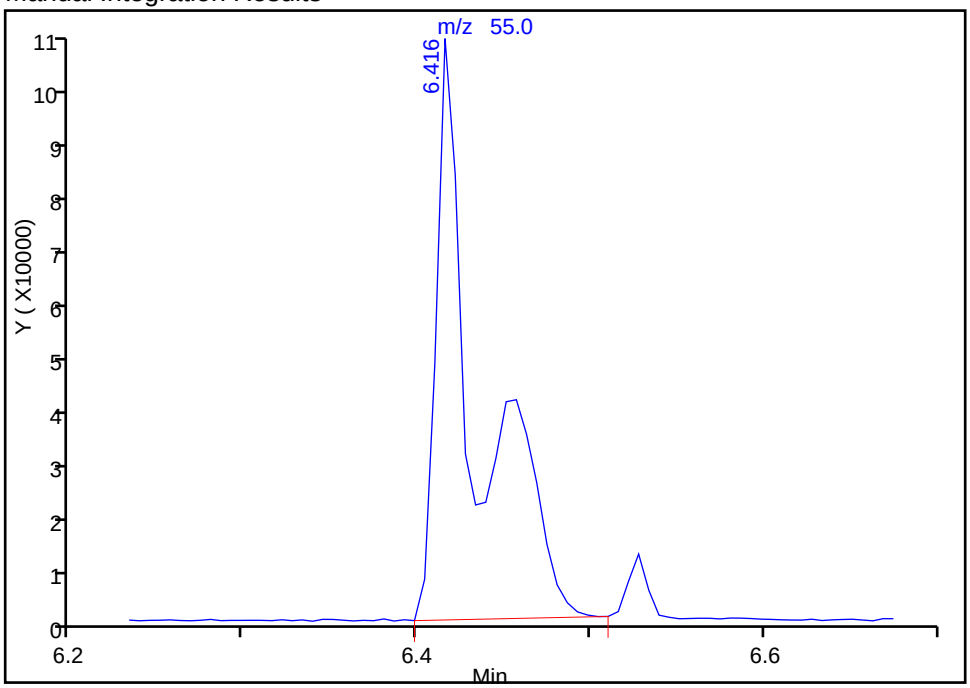
Processing Integration Results

RT: 6.42
Area: 99614
Amount: 38.000554
Amount Units: ug/ml



Manual Integration Results

RT: 6.42
Area: 171197
Amount: 49.522448
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:09:09
Audit Action: Manually Integrated
Audit Reason: Split Peak

Manual Integration/User Assign Peak Report

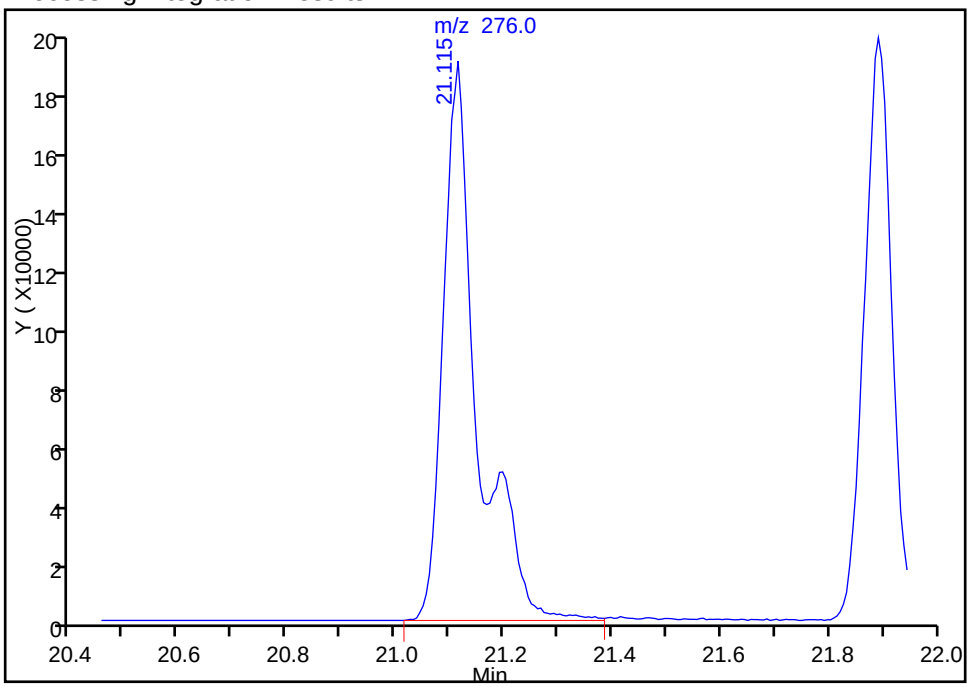
TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16641.D
Injection Date: 25-Feb-2015 13:39:30 Instrument ID: SMS_G6
Lims ID: STD050 HSL
Client ID:
Operator ID: KIEKELD ALS Bottle#: 6 Worklist Smp#: 7
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Method: SMS_G6_8270D Limit Group: MSSV - 8270D
Column: VF-5ms (0.50 mm) Detector MS SCAN

151 Indeno[1,2,3-cd]pyrene, CAS: 193-39-5

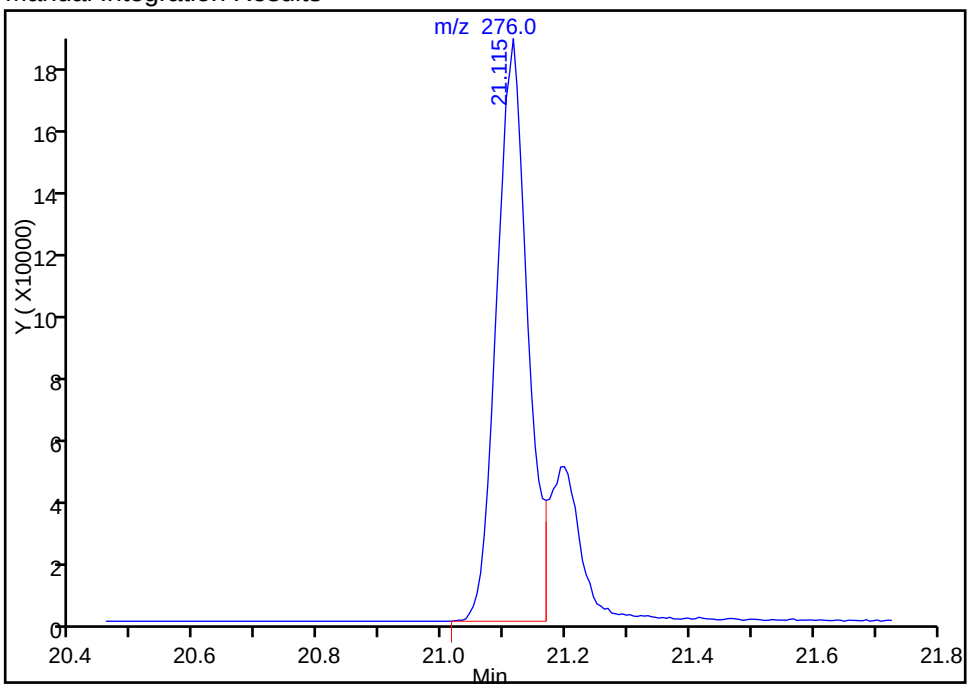
Processing Integration Results

RT: 21.11
Area: 825882
Amount: 56.977890
Amount Units: ug/ml



Manual Integration Results

RT: 21.11
Area: 654992
Amount: 50.847595
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:09:09
Audit Action: Split an Integrated Peak
Audit Reason: Shouldering

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16642.D
Lims ID: STD120 HSL
Client ID:
Sample Type: IC Calib Level: 6
Inject. Date: 25-Feb-2015 14:06:30 ALS Bottle#: 7 Worklist Smp#: 8
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: STD120 HSL
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:43 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:12:35

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	96	184348	40.0	40.0	
* 2 Naphthalene-d8	136	6.022	6.022	0.000	100	702195	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	91	392054	40.0	40.0	
* 4 Phenanthrene-d10	188	9.281	9.280	0.000	97	657834	40.0	40.0	
* 5 Chrysene-d12	240	13.710	13.710	0.000	96	610902	40.0	40.0	
* 6 Perylene-d12	264	17.903	17.903	0.000	95	528824	40.0	40.0	
$ 7 2-Fluorophenol	112	3.622	3.622	0.000	94	808477	120.0	120.6	
$ 8 Phenol-d5	99	4.410	4.410	0.000	99	1009270	120.0	120.1	
$ 9 Nitrobenzene-d5	82	5.322	5.322	0.000	92	875604	120.0	121.6	
$ 10 2-Fluorobiphenyl	172	7.069	7.069	0.000	100	1545378	120.0	116.6	
$ 11 2,4,6-Tribromophenol	330	8.575	8.575	0.000	87	184203	120.0	124.9	
$ 12 Terphenyl-d14	244	11.369	11.369	0.000	98	1459232	120.0	120.2	
13 1,4-Dioxane	88	2.352	2.352	0.000	99	333966	120.0	118.1	
14 N-Nitrosodimethylamine	74	2.581	2.581	0.000	90	518985	120.0	120.8	
15 Pyridine	79	2.622	2.622	0.000	90	891222	120.0	121.7	
23 Phenol	94	4.422	4.422	0.000	99	999282	120.0	118.6	
24 Aniline	93	4.498	4.498	0.000	98	1234749	120.0	119.9	
25 Bis(2-chloroethyl)ether	93	4.528	4.528	0.000	90	802000	120.0	121.3	
26 2-Chlorophenol	128	4.604	4.604	0.000	97	824383	120.0	119.9	
27 1,3-Dichlorobenzene	146	4.751	4.751	0.000	97	832464	120.0	119.3	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	93	837561	120.0	118.8	
29 Benzyl alcohol	108	4.904	4.904	0.000	93	532820	120.0	123.0	
30 1,2-Dichlorobenzene	146	4.963	4.963	0.000	96	802221	120.0	118.3	
31 2-Methylphenol	108	4.987	4.981	0.006	94	749809	120.0	120.0	
32 2,2'-oxybis[1-chloropropan	45	5.010	5.010	0.000	93	1164573	120.0	119.6	
38 3 & 4 Methylphenol	108	5.128	5.128	0.000	98	777835	120.0	120.1	
39 3-Methylphenol	108	5.128	5.128	0.000	98	777835	120.0	120.1	
40 4-Methylphenol	108	5.128	5.128	0.000	94	777835	120.0	120.1	
41 N-Nitrosodi-n-propylamine	70	5.145	5.145	0.000	92	569051	120.0	121.1	
42 Acetophenone	105	5.163	5.163	0.000	96	1084969	120.0	119.2	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.281	5.281	0.000	97	361563	120.0	119.5	
44 Nitrobenzene	77	5.340	5.340	0.000	92	842823	120.0	121.5	
46 Isophorone	82	5.551	5.545	0.006	99	1427939	120.0	121.2	
48 2-Nitrophenol	139	5.645	5.645	0.000	96	415910	120.0	120.3	
49 2,4-Dimethylphenol	107	5.640	5.640	0.000	94	751706	120.0	118.7	
50 Bis(2-chloroethoxy)methane	93	5.734	5.734	0.000	98	920059	120.0	118.8	
52 Benzoic acid	105	5.745	5.728	0.017	90	1285043	240.0	258.5	
53 2,4-Dichlorophenol	162	5.863	5.863	0.000	96	627452	120.0	121.0	
54 1,2,4-Trichlorobenzene	180	5.951	5.951	0.000	94	632193	120.0	119.9	
57 Naphthalene	128	6.040	6.040	0.000	98	2232749	120.0	119.5	
58 4-Chloroaniline	127	6.081	6.081	0.000	96	1011967	120.0	119.8	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	97	341381	120.0	118.9	
62 Caprolactam	55	6.434	6.428	0.006	78	433839	120.0	130.6	M
64 4-Chloro-3-methylphenol	107	6.534	6.528	0.006	98	673785	120.0	121.3	
65 2-Methylnaphthalene	142	6.716	6.716	0.000	93	1433043	120.0	119.4	
67 1-Methylnaphthalene	142	6.822	6.822	0.000	94	1344403	120.0	119.1	
68 Hexachlorocyclopentadiene	237	6.863	6.863	0.000	97	376192	120.0	122.1	
69 1,2,4,5-Tetrachlorobenzene	216	6.887	6.887	0.001	98	588950	120.0	117.9	
70 2,4,6-Trichlorophenol	196	6.992	6.992	0.000	95	416933	120.0	118.4	
72 2,4,5-Trichlorophenol	196	7.028	7.028	0.000	93	446976	120.0	119.2	
74 1,1'-Biphenyl	154	7.181	7.181	0.000	95	1699215	120.0	115.2	
75 2-Chloronaphthalene	162	7.216	7.216	0.000	98	1303430	120.0	117.2	
77 2-Nitroaniline	65	7.316	7.316	0.000	84	495382	120.0	121.3	
79 Dimethyl phthalate	163	7.463	7.463	0.000	98	1494143	120.0	114.5	
80 1,3-Dinitrobenzene	168	7.528	7.528	0.000	83	295917	120.0	127.8	
81 2,6-Dinitrotoluene	165	7.545	7.545	0.000	92	358122	120.0	118.7	
82 Acenaphthylene	152	7.639	7.639	0.000	99	2147354	120.0	118.9	
83 3-Nitroaniline	138	7.734	7.728	0.006	93	467084	120.0	119.7	
84 Acenaphthene	153	7.810	7.810	0.000	95	1360779	120.0	116.7	
86 2,4-Dinitrophenol	184	7.834	7.828	0.006	83	538794	240.0	245.2	
87 4-Nitrophenol	109	7.875	7.869	0.006	95	498510	240.0	237.0	
89 2,4-Dinitrotoluene	165	7.957	7.951	0.006	92	481458	120.0	121.7	
90 Dibenzofuran	168	7.981	7.981	0.000	97	1924090	120.0	117.1	
92 2,3,4,6-Tetrachlorophenol	232	8.098	8.092	0.006	75	371083	120.0	122.1	
94 Diethyl phthalate	149	8.157	8.157	0.000	98	1542842	120.0	116.1	
96 4-Chlorophenyl phenyl ethe	204	8.304	8.298	0.006	95	689660	120.0	116.0	
98 Fluorene	166	8.328	8.328	0.000	96	1563412	120.0	117.0	
99 4-Nitroaniline	138	8.363	8.357	0.006	85	452856	120.0	118.7	
100 4,6-Dinitro-2-methylphenol	198	8.369	8.363	0.006	75	578792	240.0	248.7	
102 N-Nitrosodiphenylamine	169	8.428	8.422	0.006	62	1152861	120.0	118.3	
103 Azobenzene	77	8.463	8.463	0.000	99	1738928	120.0	118.4	
104 1,2-Diphenylhydrazine	77	8.463	8.463	0.000	100	1738928	121.3	119.7	
111 4-Bromophenyl phenyl ether	248	8.798	8.798	0.000	71	370159	120.0	120.1	
112 Hexachlorobenzene	284	8.886	8.881	0.006	92	374109	120.0	116.8	
116 Pentachlorophenol	266	9.081	9.080	0.001	90	519082	240.0	246.6	
119 Phenanthrene	178	9.310	9.310	0.000	98	2117512	120.0	116.1	
120 Anthracene	178	9.363	9.357	0.006	98	2189501	120.0	117.1	
122 Carbazole	167	9.522	9.522	0.000	96	2245142	120.0	117.8	
123 Di-n-butyl phthalate	149	9.822	9.822	0.000	100	2671086	120.0	119.2	
128 Fluoranthene	202	10.786	10.786	0.000	98	2363681	120.0	118.3	
131 Pyrene	202	11.157	11.157	0.000	97	2459885	120.0	120.6	
136 Famphur	218	12.198	12.198	0.000	96	767713	120.0	119.2	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
137 Butyl benzyl phthalate	149	12.298	12.298	0.000	98	1219463	120.0	126.6	
140 3,3'-Dichlorobenzidine	252	13.639	13.639	0.000	76	756387	120.0	125.3	
141 Benzo[a]anthracene	228	13.680	13.680	0.000	99	2102748	120.0	119.8	
142 Bis(2-ethylhexyl) phthalat	149	13.663	13.668	-0.005	97	1589107	120.0	126.3	
143 Chrysene	228	13.780	13.774	0.006	98	1981581	120.0	120.9	
144 Di-n-octyl phthalate	149	15.568	15.574	-0.006	99	2802646	120.0	124.1	
146 Benzo[b]fluoranthene	252	16.733	16.727	0.006	99	1901054	120.0	123.1	
147 Benzo[k]fluoranthene	252	16.815	16.815	0.000	98	1968536	120.0	124.1	
148 Benzo[a]pyrene	252	17.733	17.727	0.006	80	1881464	120.0	126.7	
151 Indeno[1,2,3-cd]pyrene	276	21.127	21.127	0.000	96	1584058	120.0	127.0	M
152 Dibenz(a,h)anthracene	278	21.209	21.203	0.006	95	1573114	120.0	128.1	
153 Benzo[g,h,i]perylene	276	21.909	21.897	0.012	94	1678491	120.0	125.6	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

MS-HSLA120_00015 Amount Added: 200.00 Units: uL

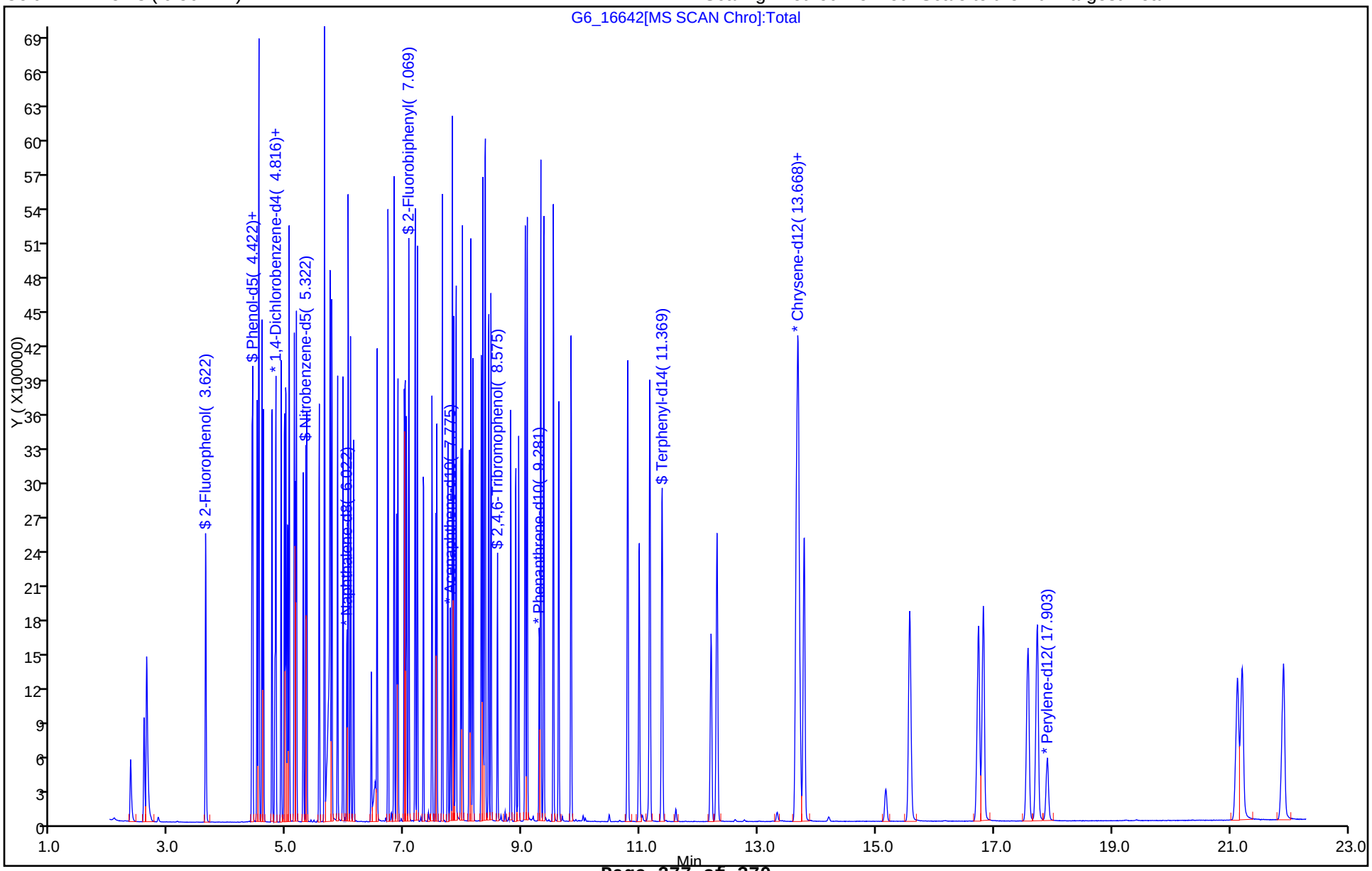
Report Date: 10-Apr-2015 12:09:43 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16642.D

Injection Date:	25-Feb-2015 14:06:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	STD120 HSL			Worklist Smp#:	8
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	7
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



Manual Integration/User Assign Peak Report

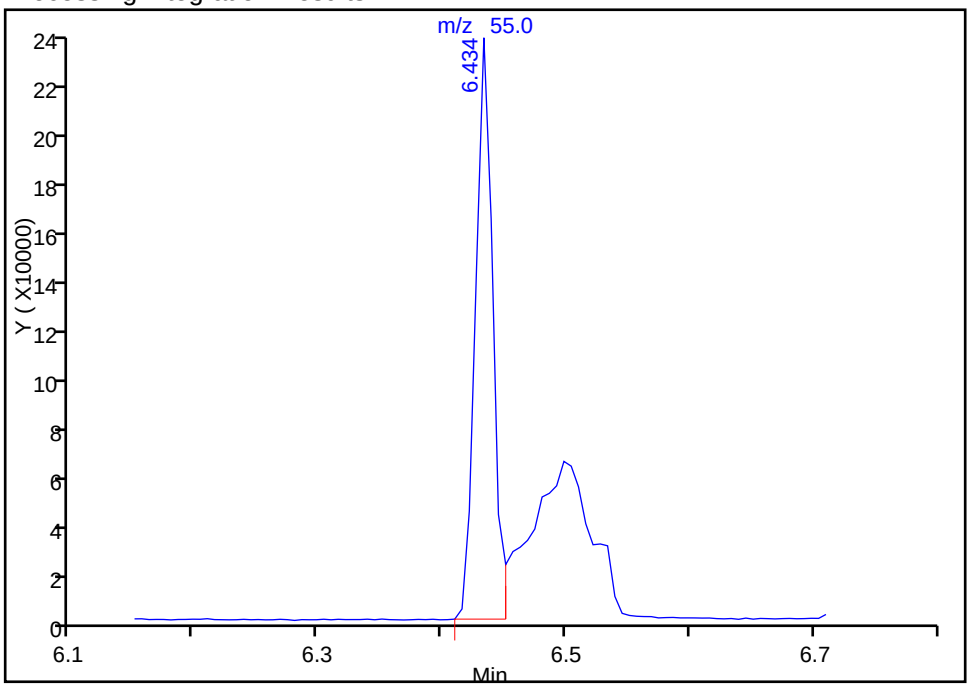
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16642.D				
Injection Date:	25-Feb-2015 14:06:30	Instrument ID:	SMS_G6		
Lims ID:	STD120 HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	7	Worklist Smp#:	8
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

62 Caprolactam, CAS: 105-60-2

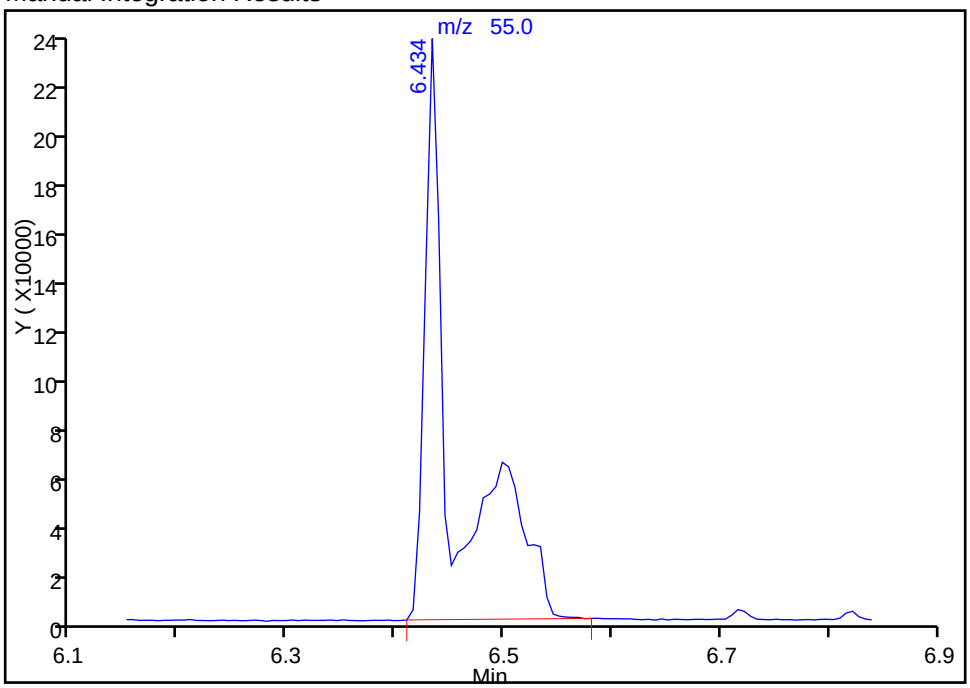
Processing Integration Results

RT: 6.43
Area: 226919
Amount: 84.337564
Amount Units: ug/ml



Manual Integration Results

RT: 6.43
Area: 433839
Amount: 130.6158
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:12:35
Audit Action: Manually Integrated
Audit Reason: Split Peak

Manual Integration/User Assign Peak Report

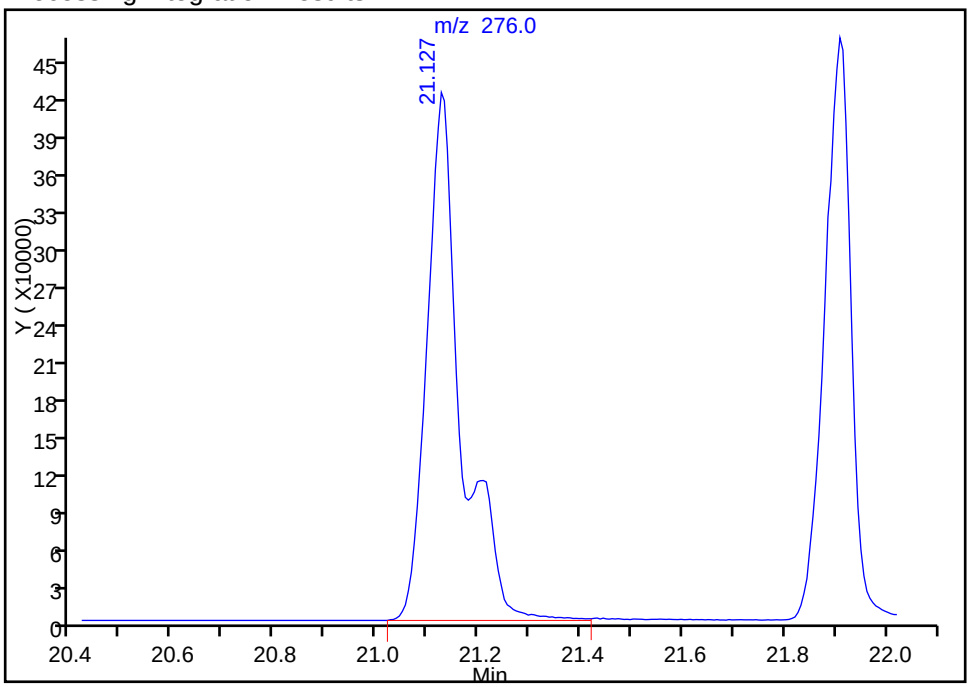
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16642.D				
Injection Date:	25-Feb-2015 14:06:30	Instrument ID:	SMS_G6		
Lims ID:	STD120 HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	7	Worklist Smp#:	8
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

151 Indeno[1,2,3-cd]pyrene, CAS: 193-39-5

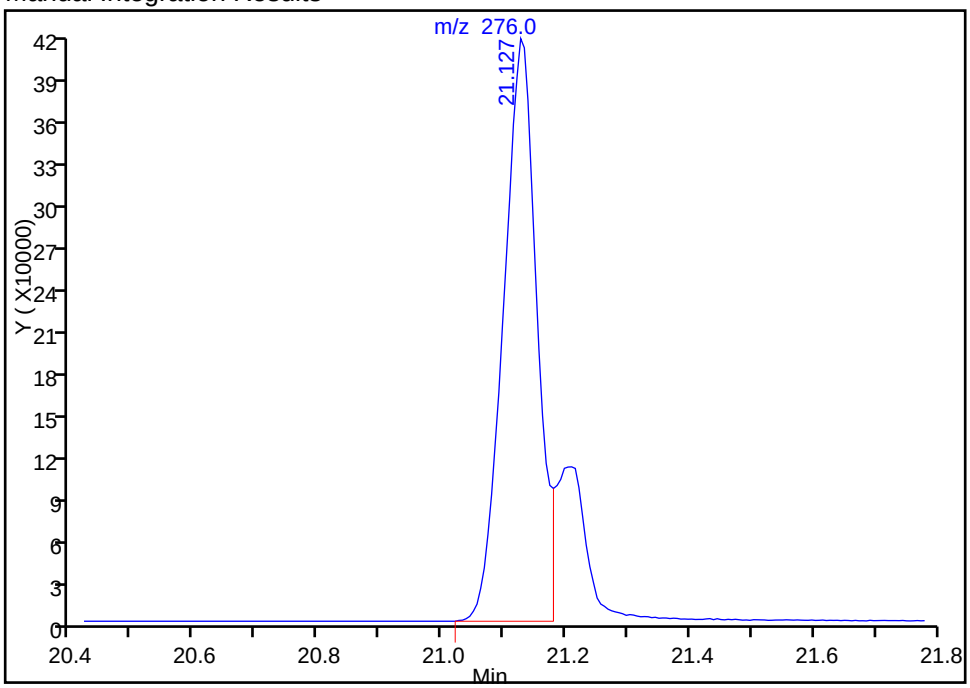
Processing Integration Results

RT: 21.13
Area: 1960031
Amount: 143.9043
Amount Units: ug/ml



Manual Integration Results

RT: 21.13
Area: 1584058
Amount: 127.0092
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:12:35
Audit Action: Split an Integrated Peak
Audit Reason: Shouldering

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16643.D
Lims ID: STD160 HSL
Client ID:
Sample Type: IC Calib Level: 7
Inject. Date: 25-Feb-2015 14:32:30 ALS Bottle#: 8 Worklist Smp#: 9
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: STD160 HSL
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:43 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:14:19

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.799	4.798	0.001	96	186214	40.0	40.0	
* 2 Naphthalene-d8	136	6.022	6.022	0.000	100	703500	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	93	384037	40.0	40.0	
* 4 Phenanthrene-d10	188	9.286	9.280	0.006	97	648058	40.0	40.0	
* 5 Chrysene-d12	240	13.716	13.710	0.006	96	607719	40.0	40.0	
* 6 Perylene-d12	264	17.909	17.903	0.006	95	521860	40.0	40.0	
$ 7 2-Fluorophenol	112	3.622	3.622	0.000	94	1061373	160.0	156.8	
$ 8 Phenol-d5	99	4.410	4.410	0.000	99	1329218	160.0	156.6	
$ 9 Nitrobenzene-d5	82	5.322	5.322	0.000	92	1140831	160.0	158.2	
$ 10 2-Fluorobiphenyl	172	7.075	7.069	0.006	100	1959377	160.0	150.9	
$ 11 2,4,6-Tribromophenol	330	8.581	8.575	0.006	92	236631	160.0	163.8	
$ 12 Terphenyl-d14	244	11.375	11.369	0.006	98	1898972	160.0	157.2	
13 1,4-Dioxane	88	2.352	2.352	0.000	98	439534	160.0	153.9	
14 N-Nitrosodimethylamine	74	2.581	2.581	0.000	91	689099	160.0	158.8	
15 Pyridine	79	2.622	2.622	0.000	91	1173437	160.0	158.6	
23 Phenol	94	4.428	4.422	0.006	99	1321311	160.0	155.3	
24 Aniline	93	4.499	4.498	0.001	98	1657573	160.0	159.3	
25 Bis(2-chloroethyl)ether	93	4.528	4.528	0.000	90	988186	160.0	147.9	
26 2-Chlorophenol	128	4.604	4.604	0.000	98	1078332	160.0	155.2	
27 1,3-Dichlorobenzene	146	4.751	4.751	0.000	97	1080606	160.0	153.3	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	94	1094435	160.0	153.7	
29 Benzyl alcohol	108	4.904	4.904	0.000	93	699135	160.0	159.7	
30 1,2-Dichlorobenzene	146	4.963	4.963	0.000	96	1044311	160.0	152.5	
31 2-Methylphenol	108	4.987	4.981	0.006	95	980672	160.0	155.4	
32 2,2'-oxybis[1-chloropropan	45	5.016	5.010	0.006	93	1512052	160.0	153.8	
38 3 & 4 Methylphenol	108	5.134	5.128	0.006	98	1031059	160.0	157.7	
39 3-Methylphenol	108	5.134	5.128	0.006	98	1031059	160.0	157.7	
40 4-Methylphenol	108	5.134	5.128	0.006	93	1031059	160.0	157.7	
41 N-Nitrosodi-n-propylamine	70	5.151	5.145	0.006	90	741748	160.0	156.3	
42 Acetophenone	105	5.169	5.163	0.006	98	1407057	160.0	153.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.281	5.281	0.000	97	469286	160.0	153.5	
44 Nitrobenzene	77	5.346	5.340	0.006	90	1089690	160.0	156.8	
46 Isophorone	82	5.551	5.545	0.006	99	1881926	160.0	159.5	
48 2-Nitrophenol	139	5.646	5.645	0.001	93	534875	160.0	154.4	
49 2,4-Dimethylphenol	107	5.640	5.640	0.000	93	981303	160.0	154.6	
50 Bis(2-chloroethoxy)methane	93	5.740	5.734	0.006	98	1210771	160.0	156.1	
52 Benzoic acid	105	5.757	5.728	0.029	96	1796945	320.0	360.9	
53 2,4-Dichlorophenol	162	5.863	5.863	0.000	97	820144	160.0	157.8	
54 1,2,4-Trichlorobenzene	180	5.957	5.951	0.006	94	814451	160.0	154.2	
57 Naphthalene	128	6.046	6.040	0.006	97	2872188	160.0	153.5	
58 4-Chloroaniline	127	6.087	6.081	0.006	96	1326050	160.0	156.7	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	97	451504	160.0	156.9	
62 Caprolactam	55	6.445	6.428	0.017	78	567672	160.0	170.6	M
64 4-Chloro-3-methylphenol	107	6.534	6.528	0.006	98	879210	160.0	158.0	
65 2-Methylnaphthalene	142	6.722	6.716	0.006	93	1850886	160.0	153.9	
67 1-Methylnaphthalene	142	6.822	6.822	0.000	94	1735161	160.0	153.4	
68 Hexachlorocyclopentadiene	237	6.863	6.863	0.000	97	489123	160.0	162.1	
69 1,2,4,5-Tetrachlorobenzene	216	6.887	6.887	0.001	98	773416	160.0	154.5	
70 2,4,6-Trichlorophenol	196	6.998	6.992	0.006	94	540716	160.0	156.8	
72 2,4,5-Trichlorophenol	196	7.034	7.028	0.006	93	570902	160.0	155.4	
74 1,1'-Biphenyl	154	7.181	7.181	0.000	95	2210281	160.0	152.9	
75 2-Chloronaphthalene	162	7.216	7.216	0.000	98	1663115	160.0	152.6	
77 2-Nitroaniline	65	7.322	7.316	0.006	84	646189	160.0	161.5	
79 Dimethyl phthalate	163	7.469	7.463	0.006	98	1926232	160.0	150.7	
80 1,3-Dinitrobenzene	168	7.534	7.528	0.006	85	383728	160.0	169.2	
81 2,6-Dinitrotoluene	165	7.551	7.545	0.006	95	471768	160.0	159.7	
82 Acenaphthylene	152	7.640	7.639	0.001	99	2757941	160.0	155.8	
83 3-Nitroaniline	138	7.740	7.728	0.012	94	619865	160.0	162.2	
84 Acenaphthene	153	7.810	7.810	0.000	95	1750014	160.0	153.2	
86 2,4-Dinitrophenol	184	7.834	7.828	0.006	83	703654	320.0	325.9	
87 4-Nitrophenol	109	7.881	7.869	0.012	94	644961	320.0	313.1	
89 2,4-Dinitrotoluene	165	7.963	7.951	0.012	93	625683	160.0	161.5	
90 Dibenzofuran	168	7.987	7.981	0.006	97	2465817	160.0	153.2	
92 2,3,4,6-Tetrachlorophenol	232	8.098	8.092	0.006	76	475470	160.0	159.7	
94 Diethyl phthalate	149	8.163	8.157	0.006	98	1980010	160.0	152.1	
96 4-Chlorophenyl phenyl ethe	204	8.304	8.298	0.006	95	895954	160.0	153.9	
98 Fluorene	166	8.334	8.328	0.006	96	1991029	160.0	152.1	
99 4-Nitroaniline	138	8.369	8.357	0.012	85	595856	160.0	159.5	
100 4,6-Dinitro-2-methylphenol	198	8.375	8.363	0.012	84	761080	320.0	331.9	
102 N-Nitrosodiphenylamine	169	8.428	8.422	0.006	63	1464837	160.0	152.6	
103 Azobenzene	77	8.469	8.463	0.006	99	2228985	160.0	154.9	
104 1,2-Diphenylhydrazine	77	8.469	8.463	0.006	100	2228985	161.8	156.6	
111 4-Bromophenyl phenyl ether	248	8.804	8.798	0.006	72	477097	160.0	157.1	
112 Hexachlorobenzene	284	8.886	8.881	0.006	92	486080	160.0	154.1	
116 Pentachlorophenol	266	9.081	9.080	0.001	90	669774	320.0	323.0	
119 Phenanthrene	178	9.310	9.310	0.000	98	2727512	160.0	151.7	
120 Anthracene	178	9.363	9.357	0.006	98	2822765	160.0	153.3	
122 Carbazole	167	9.528	9.522	0.006	96	2890531	160.0	154.0	
123 Di-n-butyl phthalate	149	9.828	9.822	0.006	100	3456767	160.0	156.5	
128 Fluoranthene	202	10.786	10.786	0.000	98	3068105	160.0	155.8	
131 Pyrene	202	11.163	11.157	0.006	97	3191260	160.0	157.3	
136 Famphur	218	12.204	12.198	0.006	96	941175	160.0	147.0	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
137 Butyl benzyl phthalate	149	12.304	12.298	0.006	98	1597755	160.0	166.8	
140 3,3'-Dichlorobenzidine	252	13.651	13.639	0.012	75	987057	160.0	164.4	
141 Benzo[a]anthracene	228	13.686	13.680	0.006	99	2746652	160.0	157.3	
142 Bis(2-ethylhexyl) phthalat	149	13.674	13.668	0.006	97	2060022	160.0	164.6	
143 Chrysene	228	13.786	13.774	0.012	98	2595420	160.0	159.1	
144 Di-n-octyl phthalate	149	15.580	15.574	0.006	99	3712823	160.0	164.8	
146 Benzo[b]fluoranthene	252	16.745	16.727	0.018	99	2549534	160.0	167.2	
147 Benzo[k]fluoranthene	252	16.833	16.815	0.018	98	2568840	160.0	164.1	
148 Benzo[a]pyrene	252	17.745	17.727	0.018	80	2481467	160.0	169.3	
151 Indeno[1,2,3-cd]pyrene	276	21.145	21.127	0.018	96	2084773	160.0	168.0	M
152 Dibenz(a,h)anthracene	278	21.221	21.203	0.018	96	2075077	160.0	171.2	
153 Benzo[g,h,i]perylene	276	21.927	21.897	0.030	93	2188108	160.0	165.9	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

MS-HSLA160_00015 Amount Added: 200.00 Units: uL

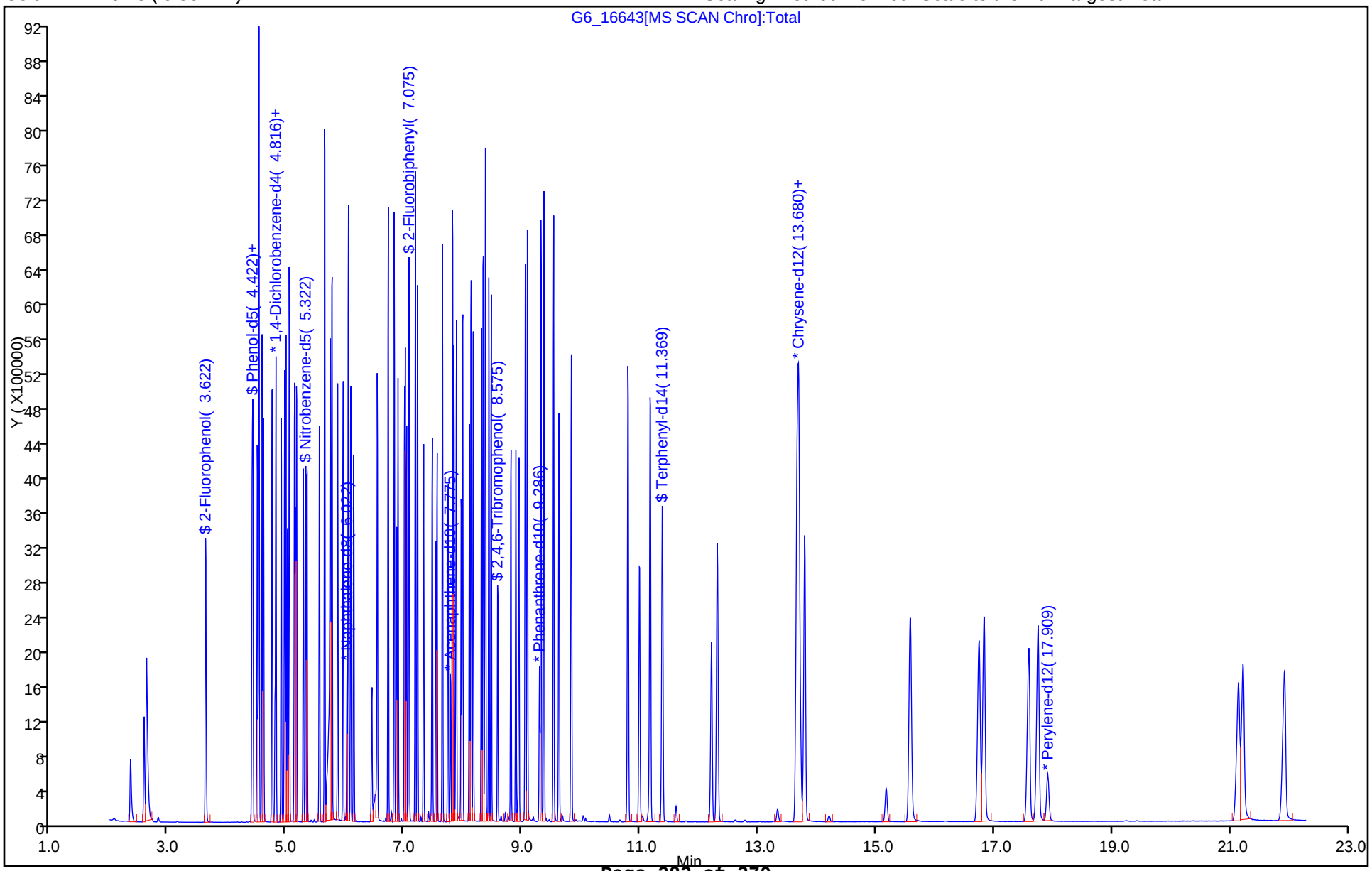
Report Date: 10-Apr-2015 12:09:44 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16643.D

Injection Date:	25-Feb-2015 14:32:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	STD160 HSL			Worklist Smp#:	9
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	8
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



Manual Integration/User Assign Peak Report

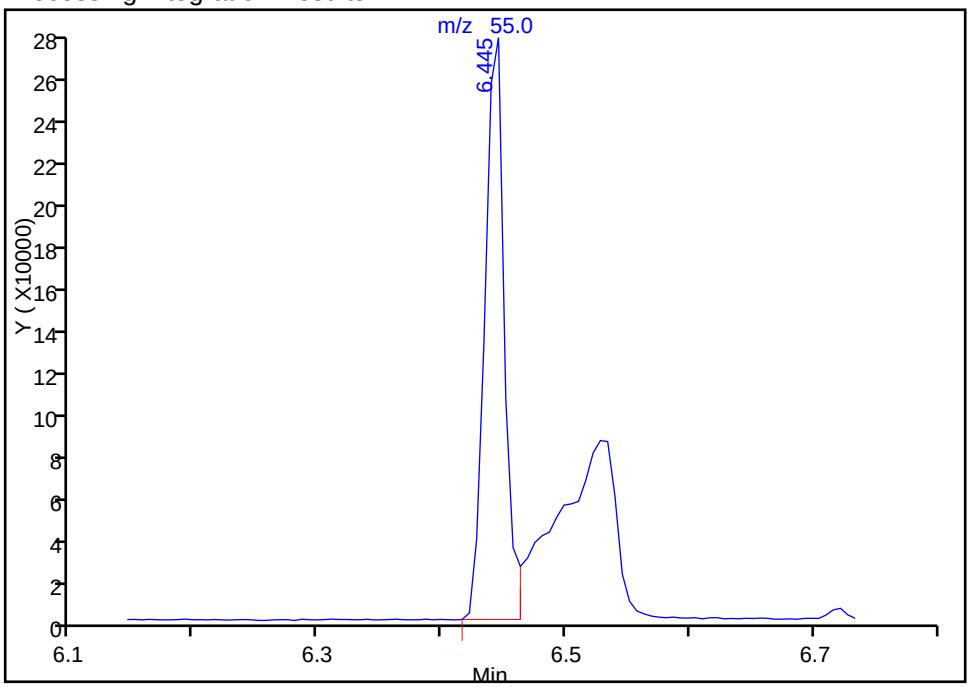
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16643.D				
Injection Date:	25-Feb-2015 14:32:30	Instrument ID:	SMS_G6		
Lims ID:	STD160 HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	8	Worklist Smp#:	9
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

62 Caprolactam, CAS: 105-60-2

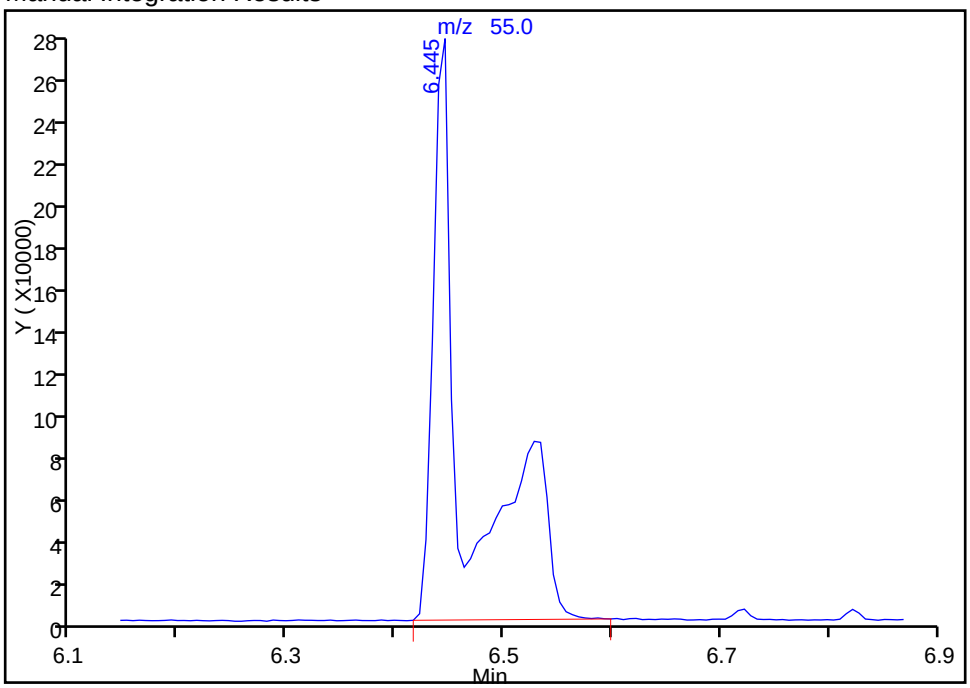
Processing Integration Results

RT: 6.45
Area: 301406
Amount: 103.5209
Amount Units: ug/ml



Manual Integration Results

RT: 6.45
Area: 567672
Amount: 170.5918
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:14:19
Audit Action: Manually Integrated
Audit Reason: Split Peak

Manual Integration/User Assign Peak Report

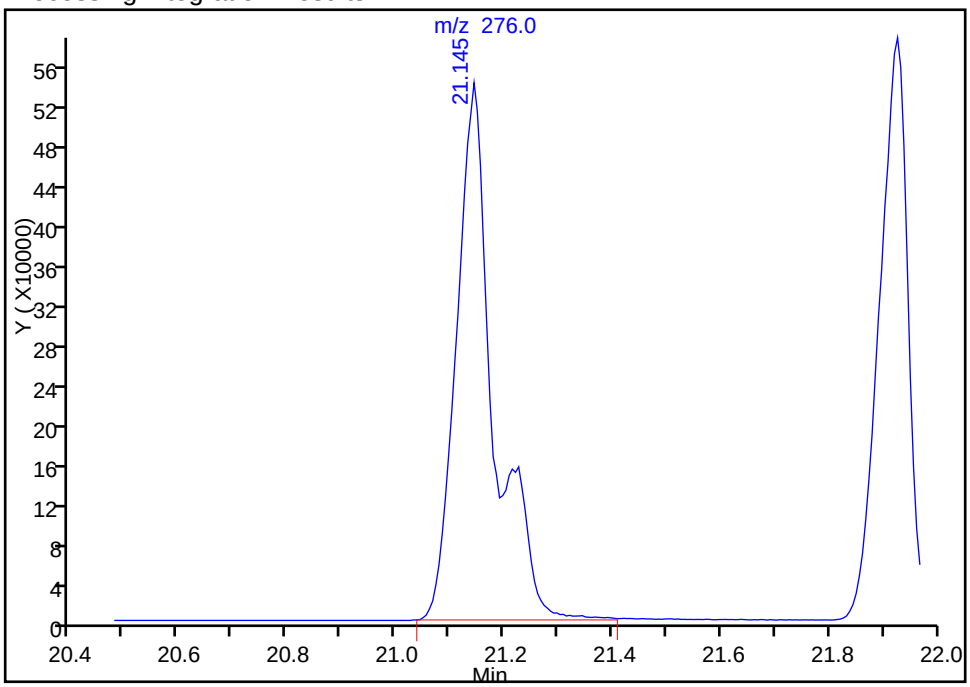
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16643.D				
Injection Date:	25-Feb-2015 14:32:30	Instrument ID:	SMS_G6		
Lims ID:	STD160 HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	8	Worklist Smp#:	9
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

151 Indeno[1,2,3-cd]pyrene, CAS: 193-39-5

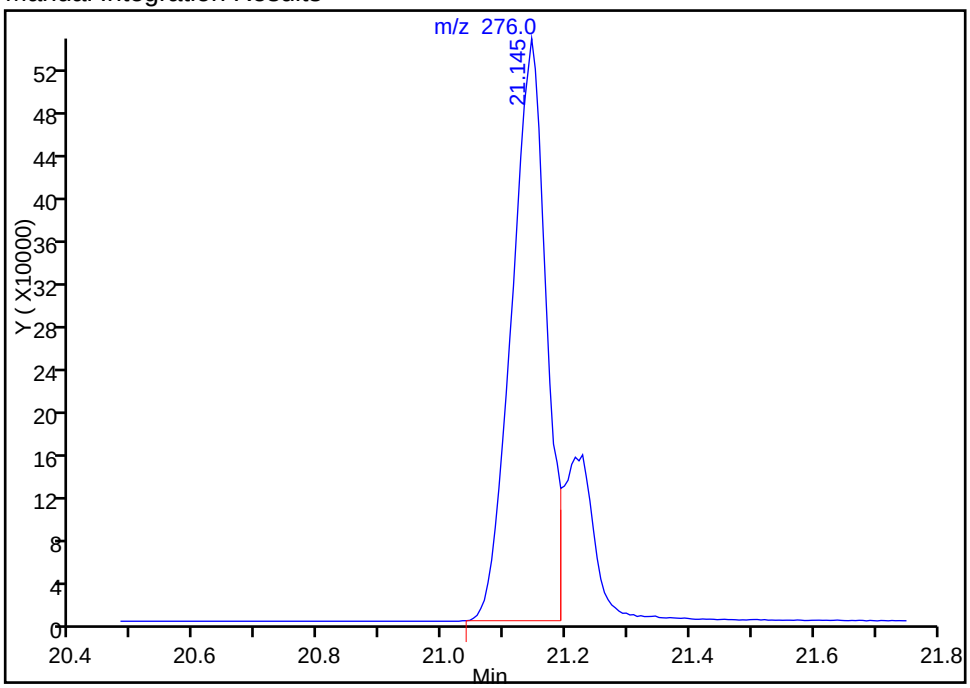
Processing Integration Results

RT: 21.14
Area: 2591657
Amount: 196.9373
Amount Units: ug/ml



Manual Integration Results

RT: 21.14
Area: 2084773
Amount: 168.0318
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:14:19
Audit Action: Split an Integrated Peak
Audit Reason: Shouldering

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D
Lims ID: STD200 HSL
Client ID:
Sample Type: IC Calib Level: 8
Inject. Date: 25-Feb-2015 14:59:30 ALS Bottle#: 9 Worklist Smp#: 10
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: STD200 HSL
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:44 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:22:35

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	96	189664	40.0	40.0	
* 2 Naphthalene-d8	136	6.022	6.022	0.000	100	731189	40.0	40.0	
* 3 Acenaphthene-d10	164	7.781	7.775	0.006	93	395282	40.0	40.0	
* 4 Phenanthrene-d10	188	9.286	9.280	0.006	97	663475	40.0	40.0	
* 5 Chrysene-d12	240	13.721	13.710	0.011	96	626472	40.0	40.0	
* 6 Perylene-d12	264	17.909	17.903	0.006	95	533835	40.0	40.0	
$ 7 2-Fluorophenol	112	3.628	3.622	0.006	94	1340808	200.0	194.4	
$ 8 Phenol-d5	99	4.416	4.410	0.006	99	1686187	200.0	195.1	
$ 9 Nitrobenzene-d5	82	5.328	5.322	0.006	91	1438315	200.0	191.8	
$ 10 2-Fluorobiphenyl	172	7.075	7.069	0.006	99	2471944	200.0	185.0	
$ 11 2,4,6-Tribromophenol	330	8.580	8.575	0.005	91	308954	200.0	207.8	
$ 12 Terphenyl-d14	244	11.374	11.369	0.005	98	2411505	200.0	193.7	
13 1,4-Dioxane	88	2.351	2.352	-0.001	98	558095	200.0	191.8	
14 N-Nitrosodimethylamine	74	2.581	2.581	0.000	91	878936	200.0	198.9	
15 Pyridine	79	2.622	2.622	0.000	91	1482135	200.0	196.6	
23 Phenol	94	4.428	4.422	0.006	99	1676043	200.0	193.4	
24 Aniline	93	4.504	4.498	0.006	98	2067510	200.0	195.1	
25 Bis(2-chloroethyl)ether	93	4.528	4.528	0.000	92	1252467	200.0	184.1	
26 2-Chlorophenol	128	4.604	4.604	0.000	97	1354936	200.0	191.5	
27 1,3-Dichlorobenzene	146	4.751	4.751	0.000	97	1361408	200.0	189.6	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	92	1353318	200.0	186.6	
29 Benzyl alcohol	108	4.910	4.904	0.006	93	883290	200.0	198.1	
30 1,2-Dichlorobenzene	146	4.963	4.963	0.000	95	1312433	200.0	188.1	
31 2-Methylphenol	108	4.987	4.981	0.006	95	1246269	200.0	193.9	
32 2,2'-oxybis[1-chloropropan	45	5.016	5.010	0.006	93	1907157	200.0	190.4	
38 3 & 4 Methylphenol	108	5.134	5.128	0.006	98	1296981	200.0	194.7	
39 3-Methylphenol	108	5.134	5.128	0.006	98	1296981	200.0	194.7	
40 4-Methylphenol	108	5.134	5.128	0.006	93	1296981	200.0	194.7	
41 N-Nitrosodi-n-propylamine	70	5.151	5.145	0.006	91	929339	200.0	192.2	
42 Acetophenone	105	5.169	5.163	0.006	97	1777329	200.0	189.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.281	5.281	0.000	97	593205	200.0	190.6	
44 Nitrobenzene	77	5.345	5.340	0.005	95	1376078	200.0	190.5	
46 Isophorone	82	5.557	5.545	0.012	99	2388530	200.0	194.7	
48 2-Nitrophenol	139	5.645	5.645	0.000	90	672106	200.0	186.7	
49 2,4-Dimethylphenol	107	5.645	5.640	0.005	93	1213524	200.0	184.0	
50 Bis(2-chloroethoxy)methane	93	5.739	5.734	0.005	98	1508491	200.0	187.1	
52 Benzoic acid	105	5.775	5.728	0.047	89	2237473	400.0	432.3	
53 2,4-Dichlorophenol	162	5.869	5.863	0.006	96	1026716	200.0	190.1	
54 1,2,4-Trichlorobenzene	180	5.957	5.951	0.006	94	1032581	200.0	188.1	
57 Naphthalene	128	6.045	6.040	0.005	97	3606766	200.0	185.4	
58 4-Chloroaniline	127	6.087	6.081	0.005	96	1671870	200.0	190.0	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	98	556376	200.0	186.0	
62 Caprolactam	55	6.451	6.428	0.023	77	732007	200.0	211.6	M
64 4-Chloro-3-methylphenol	107	6.539	6.528	0.011	98	1115868	200.0	193.0	
65 2-Methylnaphthalene	142	6.722	6.716	0.006	93	2311519	200.0	185.0	
67 1-Methylnaphthalene	142	6.828	6.822	0.006	95	2161606	200.0	183.9	
68 Hexachlorocyclopentadiene	237	6.869	6.863	0.006	96	621756	200.0	200.2	
69 1,2,4,5-Tetrachlorobenzene	216	6.886	6.887	0.000	98	984311	200.0	189.2	
70 2,4,6-Trichlorophenol	196	6.998	6.992	0.006	94	681085	200.0	191.9	
72 2,4,5-Trichlorophenol	196	7.034	7.028	0.006	93	728184	200.0	192.6	
74 1,1'-Biphenyl	154	7.181	7.181	0.000	96	2739578	200.0	184.1	
75 2-Chloronaphthalene	162	7.222	7.216	0.006	98	2078262	200.0	185.3	
77 2-Nitroaniline	65	7.322	7.316	0.006	85	812592	200.0	197.3	
79 Dimethyl phthalate	163	7.469	7.463	0.006	98	2440056	200.0	185.4	
80 1,3-Dinitrobenzene	168	7.533	7.528	0.005	85	495821	200.0	212.4	
81 2,6-Dinitrotoluene	165	7.551	7.545	0.006	94	595647	200.0	195.8	
82 Acenaphthylene	152	7.645	7.639	0.006	99	3454854	200.0	189.7	
83 3-Nitroaniline	138	7.739	7.728	0.011	94	793558	200.0	201.8	
84 Acenaphthene	153	7.816	7.810	0.006	95	2161568	200.0	183.8	
86 2,4-Dinitrophenol	184	7.839	7.828	0.011	82	903939	400.0	406.0	
87 4-Nitrophenol	109	7.886	7.869	0.017	94	814776	400.0	384.3	
89 2,4-Dinitrotoluene	165	7.963	7.951	0.012	92	798404	200.0	200.2	
90 Dibenzofuran	168	7.986	7.981	0.005	97	3062264	200.0	184.9	
92 2,3,4,6-Tetrachlorophenol	232	8.098	8.092	0.006	76	615581	200.0	200.8	
94 Diethyl phthalate	149	8.163	8.157	0.006	98	2499587	200.0	186.6	
96 4-Chlorophenyl phenyl ethe	204	8.304	8.298	0.006	96	1124576	200.0	187.6	
98 Fluorene	166	8.333	8.328	0.005	95	2502594	200.0	185.7	
99 4-Nitroaniline	138	8.375	8.357	0.018	84	738699	200.0	192.1	
100 4,6-Dinitro-2-methylphenol	198	8.380	8.363	0.017	84	958230	400.0	408.2	
102 N-Nitrosodiphenylamine	169	8.433	8.422	0.011	62	1845140	200.0	187.7	
103 Azobenzene	77	8.469	8.463	0.006	99	2812289	200.0	189.9	
104 1,2-Diphenylhydrazine	77	8.469	8.463	0.006	100	2812289	202.2	192.0	
111 4-Bromophenyl phenyl ether	248	8.804	8.798	0.006	72	598149	200.0	192.4	
112 Hexachlorobenzene	284	8.886	8.881	0.006	92	623678	200.0	193.1	
116 Pentachlorophenol	266	9.086	9.080	0.006	92	859647	400.0	405.0	
119 Phenanthrene	178	9.316	9.310	0.006	98	3407975	200.0	185.2	
120 Anthracene	178	9.369	9.357	0.012	98	3510820	200.0	186.2	
122 Carbazole	167	9.527	9.522	0.005	96	3612710	200.0	187.9	
123 Di-n-butyl phthalate	149	9.827	9.822	0.005	100	4377647	200.0	193.6	
128 Fluoranthene	202	10.792	10.786	0.006	98	3867766	200.0	191.9	
131 Pyrene	202	11.169	11.157	0.012	97	4045386	200.0	193.4	
136 Famphur	218	12.210	12.198	0.012	97	1098625	200.0	166.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
137 Butyl benzyl phthalate	149	12.304	12.298	0.006	98	2035238	200.0	206.1	
140 3,3'-Dichlorobenzidine	252	13.657	13.639	0.018	78	1242385	200.0	200.7	
141 Benzo[a]anthracene	228	13.692	13.680	0.012	99	3473559	200.0	192.9	
142 Bis(2-ethylhexyl) phthalat	149	13.674	13.668	0.006	97	2595812	200.0	201.2	
143 Chrysene	228	13.792	13.774	0.018	98	3298874	200.0	196.2	
144 Di-n-octyl phthalate	149	15.574	15.574	0.000	99	4782051	200.0	205.5	
146 Benzo[b]fluoranthene	252	16.751	16.727	0.024	99	3151061	200.0	202.1	
147 Benzo[k]fluoranthene	252	16.839	16.815	0.024	98	3328296	200.0	207.8	
148 Benzo[a]pyrene	252	17.750	17.727	0.023	81	3149998	200.0	210.1	
151 Indeno[1,2,3-cd]pyrene	276	21.156	21.127	0.029	97	2687797	200.0	210.2	M
152 Dibenz(a,h)anthracene	278	21.233	21.203	0.030	95	2616064	200.0	211.0	
153 Benzo[g,h,i]perylene	276	21.938	21.897	0.041	94	2789274	200.0	206.8	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

MS-HSLA200_00015 Amount Added: 200.00 Units: uL

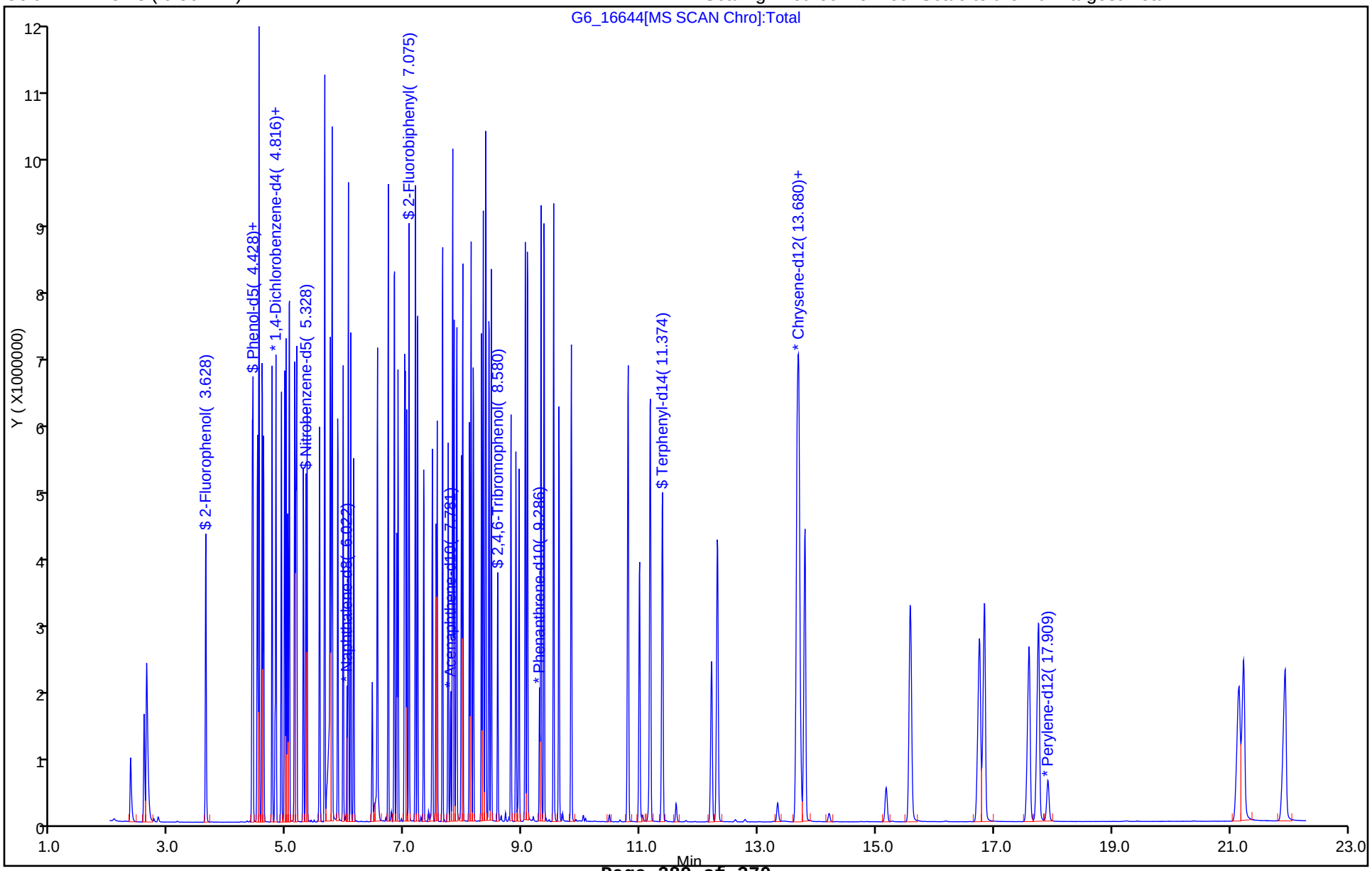
Report Date: 10-Apr-2015 12:09:45 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Injection Date:	25-Feb-2015 14:59:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	STD200 HSL			Worklist Smp#:	10
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	9
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



Manual Integration/User Assign Peak Report

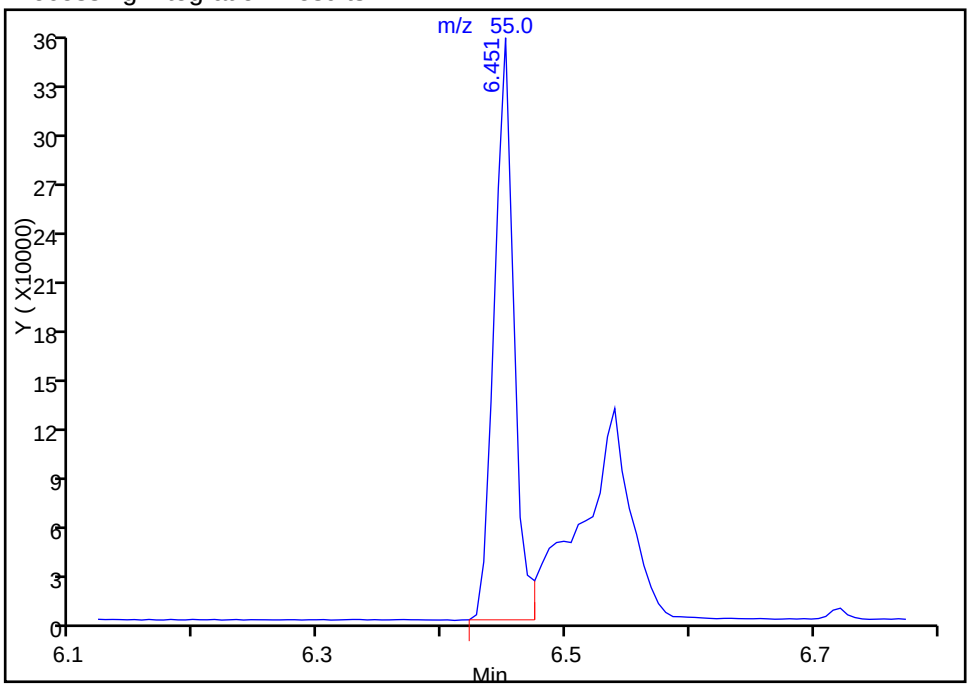
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D				
Injection Date:	25-Feb-2015 14:59:30	Instrument ID:	SMS_G6		
Lims ID:	STD200 HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	9	Worklist Smp#:	10
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

62 Caprolactam, CAS: 105-60-2

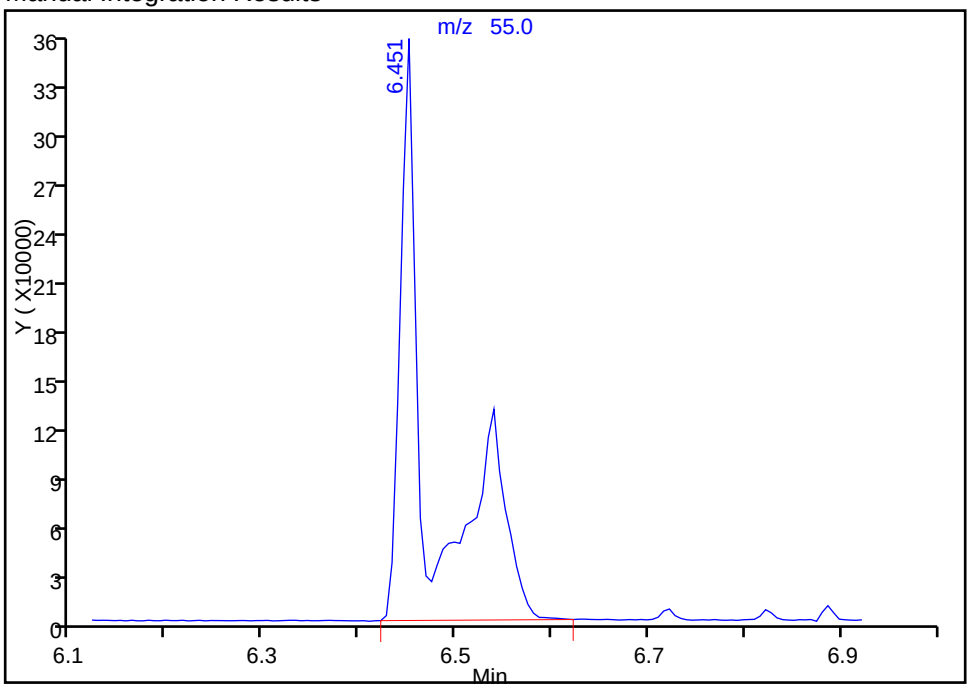
Processing Integration Results

RT: 6.45
Area: 385948
Amount: 119.0334
Amount Units: ug/ml



Manual Integration Results

RT: 6.45
Area: 732007
Amount: 211.6461
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:22:35
Audit Action: Manually Integrated
Audit Reason: Split Peak

Manual Integration/User Assign Peak Report

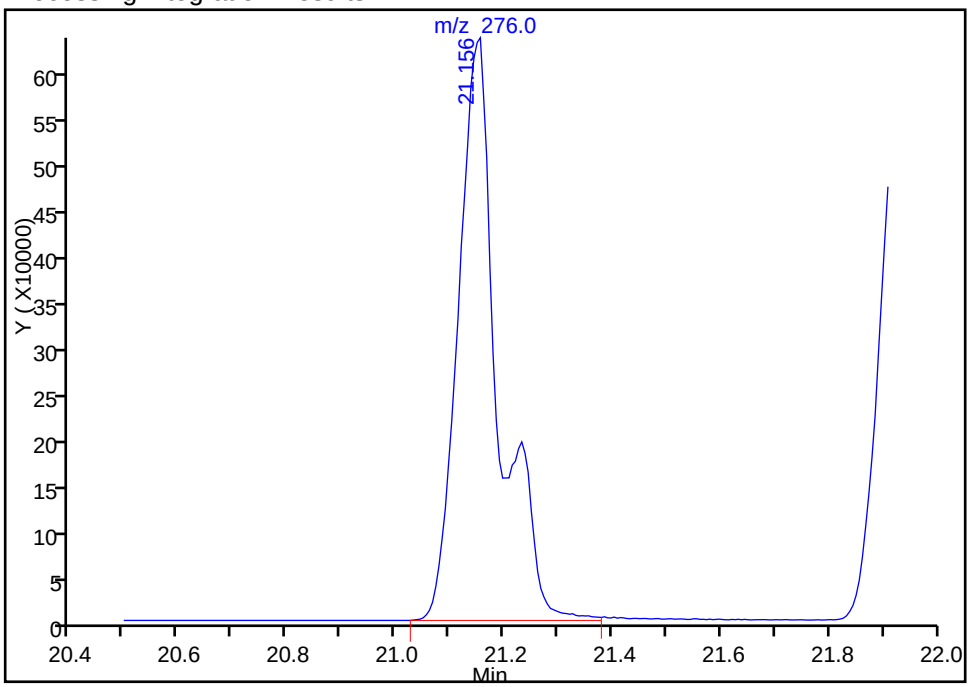
TestAmerica Denver

Data File:	\\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D				
Injection Date:	25-Feb-2015 14:59:30	Instrument ID:	SMS_G6		
Lims ID:	STD200 HSL				
Client ID:					
Operator ID:	KIEKELD	ALS Bottle#:	9	Worklist Smp#:	10
Injection Vol:	0.5 ul	Dil. Factor:	1.0000		
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		
Column:	VF-5ms (0.50 mm)	Detector	MS SCAN		

151 Indeno[1,2,3-cd]pyrene, CAS: 193-39-5

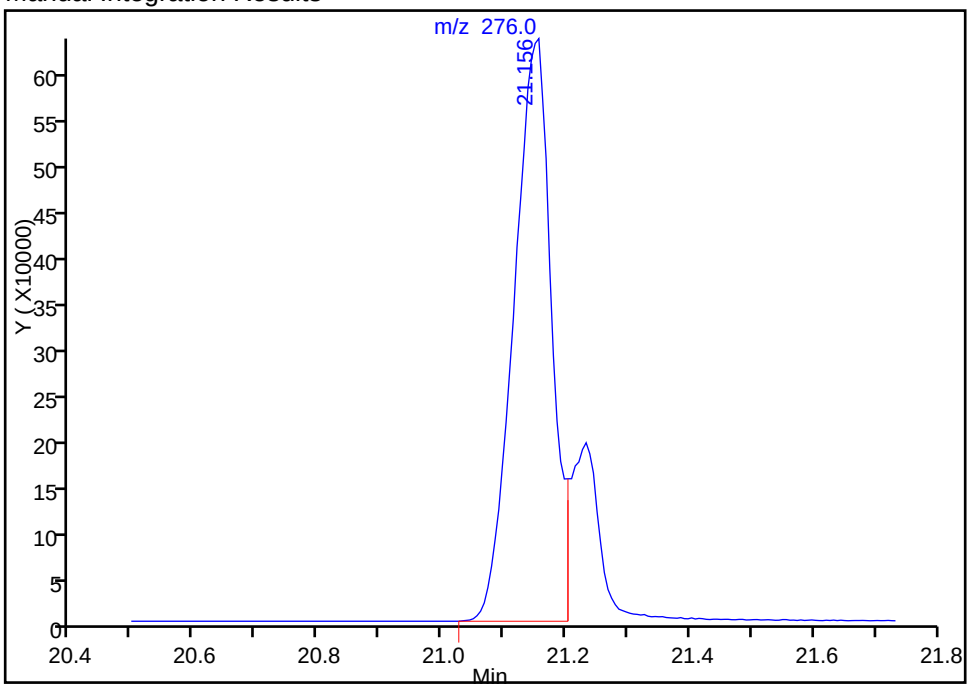
Processing Integration Results

RT: 21.16
Area: 3276282
Amount: 249.0016
Amount Units: ug/ml



Manual Integration Results

RT: 21.16
Area: 2687797
Amount: 210.1505
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:22:35
Audit Action: Split an Integrated Peak
Audit Reason: Shouldering

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: ICV 280-272059/11 Calibration Date: 02/25/2015 15:26

Instrument ID: SMS_G6 Calib Start Date: 02/25/2015 11:53

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 02/25/2015 14:59

Lab File ID: G6_16645.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dioxane	Ave	0.6137	0.5829		95.0	100	-5.0	30.0
N-Nitrosodimethylamine	Ave	0.9321	0.8898		95.5	100	-4.5	30.0
Pyridine	Ave	1.590	1.591		100	100	0.1	30.0
Phenol	Ave	1.828	1.782	0.8000	97.5	100	-2.5	30.0
Aniline	Ave	2.235	2.065		92.4	100	-7.6	30.0
Bis(2-chloroethyl)ether	Ave	1.435	1.326	0.7000	92.4	100	-7.6	30.0
2-Chlorophenol	Ave	1.492	1.435	0.8000	96.2	100	-3.8	30.0
1,3-Dichlorobenzene	Ave	1.514	1.471		97.1	100	-2.9	30.0
1,4-Dichlorobenzene	Ave	1.530	1.484		97.0	100	-3.0	30.0
Benzyl alcohol	Ave	0.9402	0.9405		100	100	0.0	30.0
1,2-Dichlorobenzene	Ave	1.471	1.406		95.6	100	-4.4	30.0
2-Methylphenol	Ave	1.355	1.303	0.7000	96.1	100	-3.9	30.0
bis (2-chloroisopropyl) ether	Ave	2.112	2.069	0.0100	70.5	72.0	-2.1	30.0
3 & 4 Methylphenol	Ave	1.405	1.360		96.8	100	-3.2	30.0
3-Methylphenol	Ave	1.405	1.360		96.8	100	-3.2	30.0
4-Methylphenol	Ave	1.405	1.360	0.6000	96.8	100	-3.2	30.0
N-Nitrosodi-n-propylamine	Ave	1.020	1.004	0.5000	98.5	100	-1.5	30.0
Acetophenone	Ave	1.974	1.893	0.0100	95.9	100	-4.1	30.0
Hexachloroethane	Ave	0.6565	0.6469	0.3000	98.5	100	-1.5	30.0
Nitrobenzene	Ave	0.3951	0.3825		96.8	100	-3.2	30.0
Isophorone	Ave	0.6711	0.7100	0.4000	106	100	5.8	30.0
2,4-Dimethylphenol	Ave	0.3608	0.3058	0.2000	84.7	100	-15.3	30.0
2-Nitrophenol	Ave	0.1969	0.1879	0.1000	95.4	100	-4.6	30.0
Bis(2-chloroethoxy)methane	Ave	0.4411	0.4105	0.3000	93.1	100	-6.9	30.0
Benzoic acid	Ave	0.2831	0.3036		214	200	7.2	30.0
2,4-Dichlorophenol	Ave	0.2955	0.2834	0.2000	95.9	100	-4.1	30.0
1,2,4-Trichlorobenzene	Ave	0.3003	0.2921		97.3	100	-2.7	30.0
Naphthalene	Ave	1.064	1.022	0.7000	96.0	100	-4.0	30.0
4-Chloroaniline	Ave	0.4813	0.4432	0.0100	92.1	100	-7.9	30.0
Hexachlorobutadiene	Ave	0.1636	0.1576	0.0100	96.3	100	-3.7	30.0
4-Chloro-3-methylphenol	Ave	0.3163	0.3048	0.2000	96.4	100	-3.6	30.0
2-Methylnaphthalene	Ave	0.6836	0.6500	0.4000	95.1	100	-4.9	30.0
1-Methylnaphthalene	Ave	0.6430	0.6099		94.9	100	-5.1	30.0
Hexachlorocyclopentadiene	Ave	0.3143	0.3438	0.0500	109	100	9.4	30.0
1,2,4,5-Tetrachlorobenzene	Ave	0.2846	0.2770	0.0100	97.3	100	-2.7	30.0
2,4,6-Trichlorophenol	Ave	0.3592	0.3534	0.2000	98.4	100	-1.6	30.0
2,4,5-Trichlorophenol	Ave	0.3826	0.3767	0.2000	98.5	100	-1.5	30.0
1,1'-Biphenyl	Ave	1.506	1.446		96.1	100	-3.9	30.0
2-Chloronaphthalene	Ave	1.135	1.087	0.8000	95.8	100	-4.2	30.0
2-Nitroaniline	Ave	0.4168	0.4117	0.0100	98.8	100	-1.2	30.0

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: ICV 280-272059/11 Calibration Date: 02/25/2015 15:26

Instrument ID: SMS_G6 Calib Start Date: 02/25/2015 11:53

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 02/25/2015 14:59

Lab File ID: G6_16645.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dimethyl phthalate	Ave	1.332	1.233	0.0100	92.6	100	-7.4	30.0
1,3-Dinitrobenzene	Ave	0.2362	0.2353		99.6	100	-0.4	30.0
2,6-Dinitrotoluene	Ave	0.3078	0.3024	0.2000	98.3	100	-1.7	30.0
Acenaphthylene	Ave	1.843	1.852	0.9000	100	100	0.5	30.0
3-Nitroaniline	Ave	0.3980	0.3797	0.0100	95.4	100	-4.6	30.0
Acenaphthene	Ave	1.190	1.174	0.9000	98.7	100	-1.3	30.0
2,4-Dinitrophenol	Lin2		0.2239	0.0100	200	200	0.1	30.0
4-Nitrophenol	Ave	0.2146	0.2082	0.0100	194	200	-3.0	30.0
2,4-Dinitrotoluene	Ave	0.4036	0.4040	0.2000	100	100	0.1	30.0
Dibenzofuran	Ave	1.676	1.618	0.8000	96.5	100	-3.5	30.0
2,3,4,6-Tetrachlorophenol	Ave	0.3102	0.3099	0.0100	99.9	100	-0.0	30.0
Diethyl phthalate	Ave	1.356	1.312	0.0100	96.8	100	-3.2	30.0
4-Chlorophenyl phenyl ether	Ave	0.6065	0.5809	0.4000	95.8	100	-4.2	30.0
Fluorene	Ave	1.364	1.335	0.9000	97.9	100	-2.1	30.0
4-Nitroaniline	Ave	0.3891	0.3770	0.0100	96.9	100	-3.1	30.0
4,6-Dinitro-2-methylphenol	Ave	0.1415	0.1482	0.0100	209	200	4.7	30.0
N-Nitrosodiphenylamine	Ave	0.5925	0.5590	0.0100	94.3	100	-5.7	30.0
1,2-Diphenylhydrazine(as Azobenzene)	Ave	1.482	1.440		98.2	101	-2.9	30.0
Azobenzene	Ave	1.498	1.456		97.1	100	-2.9	30.0
4-Bromophenyl phenyl ether	Ave	0.1874	0.1831	0.1000	97.7	100	-2.3	30.0
Hexachlorobenzene	Ave	0.1947	0.1887	0.1000	96.9	100	-3.1	30.0
Pentachlorophenol	Ave	0.1280	0.1337	0.0500	209	200	4.5	30.0
Phenanthrene	Ave	1.109	1.061	0.7000	95.6	100	-4.4	30.0
Anthracene	Ave	1.137	1.092	0.7000	96.0	100	-4.0	30.0
Carbazole	Ave	1.159	1.094	0.0100	94.4	100	-5.6	30.0
Di-n-butyl phthalate	Ave	1.363	1.336	0.0100	98.0	100	-2.0	30.0
Fluoranthene	Ave	1.215	1.172	0.6000	96.5	100	-3.5	30.0
Pyrene	Ave	1.336	1.309	0.6000	98.0	100	-2.0	30.0
Butyl benzyl phthalate	Ave	0.6306	0.6514	0.0100	103	100	3.3	30.0
Bis(2-ethylhexyl) phthalate	Ave	0.8237	0.8366	0.0100	102	100	1.6	30.0
Benzo[a]anthracene	Ave	1.150	1.121	0.8000	97.5	100	-2.5	30.0
Chrysene	Ave	1.074	1.079	0.7000	100	100	0.5	30.0
Di-n-octyl phthalate	Lin2		1.489	0.0100	101	100	1.0	30.0
Benzo[b]fluoranthene	Ave	1.169	1.152	0.7000	98.6	100	-1.4	30.0
Benzo[k]fluoranthene	Ave	1.200	1.201	0.7000	100	100	0.1	30.0
Benzo[a]pyrene	Ave	1.123	1.105	0.7000	98.4	100	-1.6	30.0
Indeno[1,2,3-cd]pyrene	Ave	0.8166	0.7800	0.5000	95.5	100	-4.5	30.0
Dibenz(a,h)anthracene	Ave	0.9288	0.9593	0.4000	103	100	3.3	30.0
Benzo[g,h,i]perylene	Ave	1.011	1.017	0.5000	101	100	0.7	30.0

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16645.D
Lims ID: ICV HSL 1
Client ID:
Sample Type: ICV
Inject. Date: 25-Feb-2015 15:26:30 ALS Bottle#: 10 Worklist Smp#: 11
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: ICV HSL 1
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist:

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:44 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:24:38

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	96	188037	40.0	40.0	
* 2 Naphthalene-d8	136	6.022	6.022	0.000	100	718467	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	93	393887	40.0	40.0	
* 4 Phenanthrene-d10	188	9.286	9.280	0.006	97	663955	40.0	40.0	
* 5 Chrysene-d12	240	13.710	13.710	0.000	96	625452	40.0	40.0	
* 6 Perylene-d12	264	17.909	17.903	0.006	95	534138	40.0	40.0	
13 1,4-Dioxane	88	2.352	2.352	0.000	98	273993	100.0	95.0	
14 N-Nitrosodimethylamine	74	2.581	2.581	0.000	91	418285	100.0	95.5	
15 Pyridine	79	2.622	2.622	0.000	91	748145	100.0	100.1	
23 Phenol	94	4.422	4.422	0.000	99	837710	100.0	97.5	
24 Aniline	93	4.499	4.498	0.000	98	970675	100.0	92.4	
25 Bis(2-chloroethyl)ether	93	4.528	4.528	0.000	88	623551	100.0	92.4	
26 2-Chlorophenol	128	4.604	4.604	0.000	97	674582	100.0	96.2	
27 1,3-Dichlorobenzene	146	4.751	4.751	0.000	97	691355	100.0	97.1	
28 1,4-Dichlorobenzene	146	4.816	4.816	0.000	92	697797	100.0	97.0	
29 Benzyl alcohol	108	4.904	4.904	0.000	93	442115	100.0	100.0	
30 1,2-Dichlorobenzene	146	4.963	4.963	0.000	96	660939	100.0	95.6	
31 2-Methylphenol	108	4.987	4.981	0.006	94	612489	100.0	96.1	
32 2,2'-oxybis[1-chloropropan	45	5.016	5.010	0.006	94	700124	72.0	70.5	
38 3 & 4 Methylphenol	108	5.128	5.128	0.000	98	639290	100.0	96.8	
39 3-Methylphenol	108	5.128	5.128	0.000	98	639290	100.0	96.8	
40 4-Methylphenol	108	5.128	5.128	0.000	93	639290	100.0	96.8	
41 N-Nitrosodi-n-propylamine	70	5.146	5.145	0.001	91	471958	100.0	98.5	
42 Acetophenone	105	5.163	5.163	0.000	96	889654	100.0	95.9	
43 Hexachloroethane	117	5.281	5.281	0.000	96	304123	100.0	98.5	
44 Nitrobenzene	77	5.340	5.340	0.000	92	687043	100.0	96.8	
46 Isophorone	82	5.551	5.545	0.006	99	1275320	100.0	105.8	
48 2-Nitrophenol	139	5.646	5.645	0.001	91	337503	100.0	95.4	
49 2,4-Dimethylphenol	107	5.640	5.640	0.000	89	549179	100.0	84.7	
50 Bis(2-chloroethoxy)methane	93	5.734	5.734	0.000	98	737363	100.0	93.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
52 Benzoic acid	105	5.740	5.728	0.012	62	1090730	200.0	214.5	
53 2,4-Dichlorophenol	162	5.863	5.863	0.000	97	509014	100.0	95.9	
54 1,2,4-Trichlorobenzene	180	5.957	5.951	0.006	94	524624	100.0	97.3	
57 Naphthalene	128	6.045	6.040	0.005	97	1834953	100.0	96.0	
58 4-Chloroaniline	127	6.081	6.081	0.000	96	795990	100.0	92.1	
59 Hexachlorobutadiene	225	6.134	6.134	0.000	98	283045	100.0	96.3	
64 4-Chloro-3-methylphenol	107	6.528	6.528	0.000	98	547506	100.0	96.4	
65 2-Methylnaphthalene	142	6.722	6.716	0.006	93	1167552	100.0	95.1	
67 1-Methylnaphthalene	142	6.822	6.822	0.000	94	1095488	100.0	94.9	
68 Hexachlorocyclopentadiene	237	6.863	6.863	0.000	95	338550	100.0	109.4	
69 1,2,4,5-Tetrachlorobenzene	216	6.887	6.887	0.001	98	497548	100.0	97.3	
70 2,4,6-Trichlorophenol	196	6.992	6.992	0.000	95	347961	100.0	98.4	
72 2,4,5-Trichlorophenol	196	7.028	7.028	0.000	93	370987	100.0	98.5	
74 1,1'-Biphenyl	154	7.181	7.181	0.000	95	1424130	100.0	96.1	
75 2-Chloronaphthalene	162	7.216	7.216	0.000	98	1070439	100.0	95.8	
77 2-Nitroaniline	65	7.322	7.316	0.006	85	405449	100.0	98.8	
79 Dimethyl phthalate	163	7.463	7.463	0.000	98	1213708	100.0	92.6	
80 1,3-Dinitrobenzene	168	7.528	7.528	0.000	82	231683	100.0	99.6	
81 2,6-Dinitrotoluene	165	7.545	7.545	0.000	93	297820	100.0	98.3	
82 Acenaphthylene	152	7.639	7.639	0.000	99	1823356	100.0	100.5	
83 3-Nitroaniline	138	7.734	7.728	0.006	94	373849	100.0	95.4	
84 Acenaphthene	153	7.810	7.810	0.000	95	1156147	100.0	98.7	
86 2,4-Dinitrophenol	184	7.834	7.828	0.006	82	440859	200.0	200.3	
87 4-Nitrophenol	109	7.869	7.869	0.000	95	410060	200.0	194.1	
89 2,4-Dinitrotoluene	165	7.957	7.951	0.006	91	397792	100.0	100.1	
90 Dibenzofuran	168	7.981	7.981	0.000	97	1592968	100.0	96.5	
92 2,3,4,6-Tetrachlorophenol	232	8.098	8.092	0.006	75	305136	100.0	99.9	
94 Diethyl phthalate	149	8.163	8.157	0.006	97	1292089	100.0	96.8	
96 4-Chlorophenyl phenyl ethe	204	8.304	8.298	0.006	95	572039	100.0	95.8	
98 Fluorene	166	8.328	8.328	0.000	96	1315039	100.0	97.9	
99 4-Nitroaniline	138	8.357	8.357	0.000	85	371247	100.0	96.9	
100 4,6-Dinitro-2-methylphenol	198	8.369	8.363	0.006	83	491977	200.0	209.4	
102 N-Nitrosodiphenylamine	169	8.428	8.422	0.006	63	927943	100.0	94.3	
103 Azobenzene	77	8.463	8.463	0.000	99	1433394	100.0	97.1	
104 1,2-Diphenylhydrazine	77	8.463	8.463	0.000	100	1433394	101.1	98.2	
111 4-Bromophenyl phenyl ether	248	8.804	8.798	0.006	73	303952	100.0	97.7	
112 Hexachlorobenzene	284	8.886	8.881	0.006	92	313262	100.0	96.9	
116 Pentachlorophenol	266	9.081	9.080	0.001	89	443743	200.0	208.9	
119 Phenanthrene	178	9.310	9.310	0.000	98	1760445	100.0	95.6	
120 Anthracene	178	9.363	9.357	0.006	98	1812001	100.0	96.0	
122 Carbazole	167	9.528	9.522	0.006	96	1816093	100.0	94.4	
123 Di-n-butyl phthalate	149	9.828	9.822	0.006	100	2217309	100.0	98.0	
128 Fluoranthene	202	10.786	10.786	0.000	98	1946063	100.0	96.5	
131 Pyrene	202	11.163	11.157	0.006	97	2046147	100.0	98.0	
137 Butyl benzyl phthalate	149	12.304	12.298	0.006	98	1018562	100.0	103.3	
141 Benzo[a]anthracene	228	13.680	13.680	0.000	99	1753483	100.0	97.5	
142 Bis(2-ethylhexyl) phthalat	149	13.674	13.668	0.006	97	1308106	100.0	101.6	
143 Chrysene	228	13.780	13.774	0.006	98	1686807	100.0	100.5	
144 Di-n-octyl phthalate	149	15.574	15.574	0.000	99	2328363	100.0	101.0	
146 Benzo[b]fluoranthene	252	16.733	16.727	0.006	99	1538895	100.0	98.6	
147 Benzo[k]fluoranthene	252	16.821	16.815	0.006	98	1604191	100.0	100.1	
148 Benzo[a]pyrene	252	17.733	17.727	0.006	80	1475603	100.0	98.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
151 Indeno[1,2,3-cd]pyrene	276	21.127	21.127	0.000	98	1219599	100.0	95.5	M
152 Dibenz(a,h)anthracene	278	21.215	21.203	0.012	94	1281000	100.0	103.3	
153 Benzo[g,h,i]perylene	276	21.915	21.897	0.018	96	1358418	100.0	100.7	

QC Flag Legend

Review Flags

M - Manually Integrated

Reagents:

MS-HSLB1B3SSV_00022 Amount Added: 200.00 Units: uL

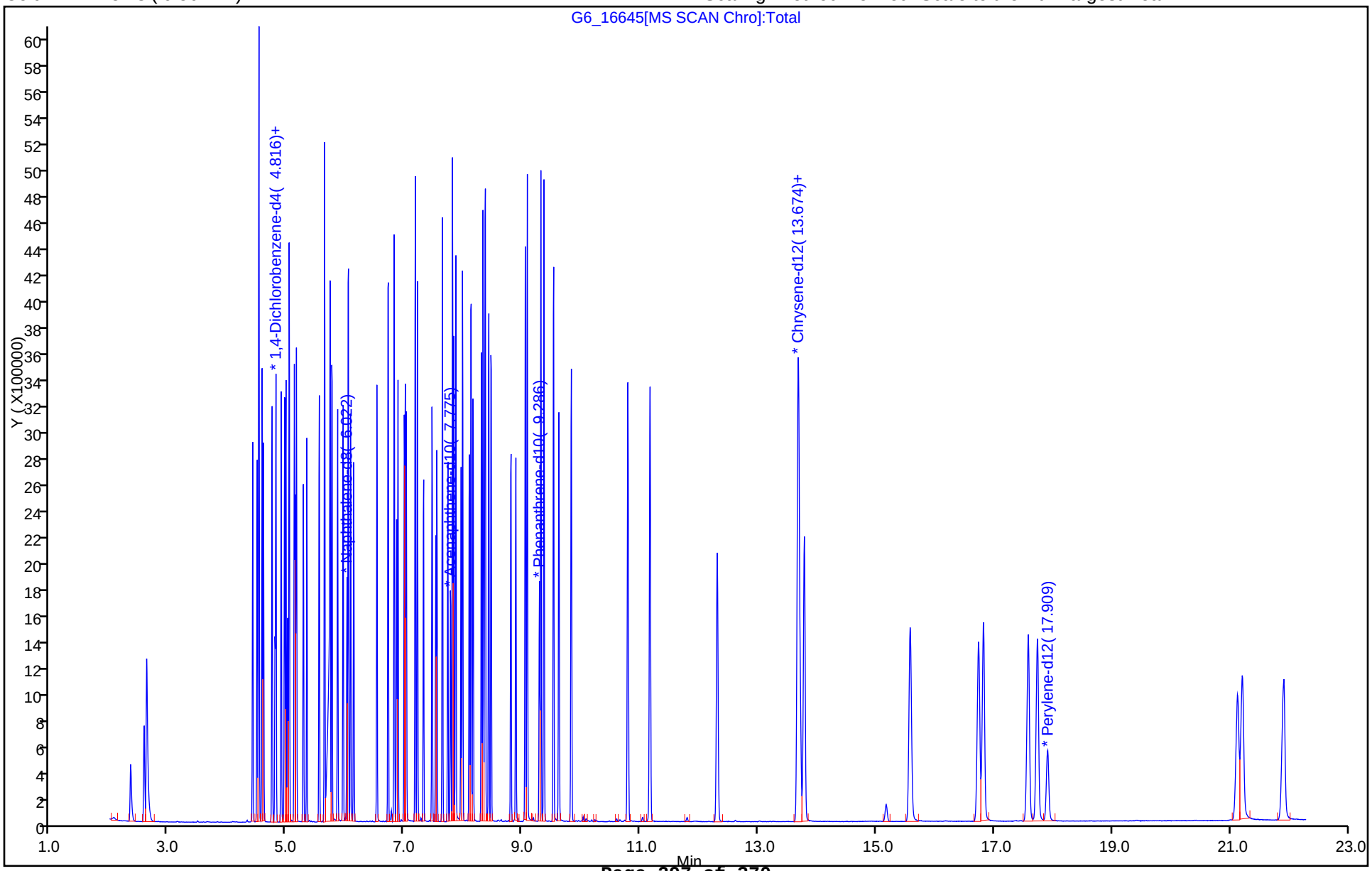
Report Date: 10-Apr-2015 12:09:46 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16645.D

Injection Date:	25-Feb-2015 15:26:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	ICV HSL 1			Worklist Smp#:	11
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	10
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



Manual Integration/User Assign Peak Report

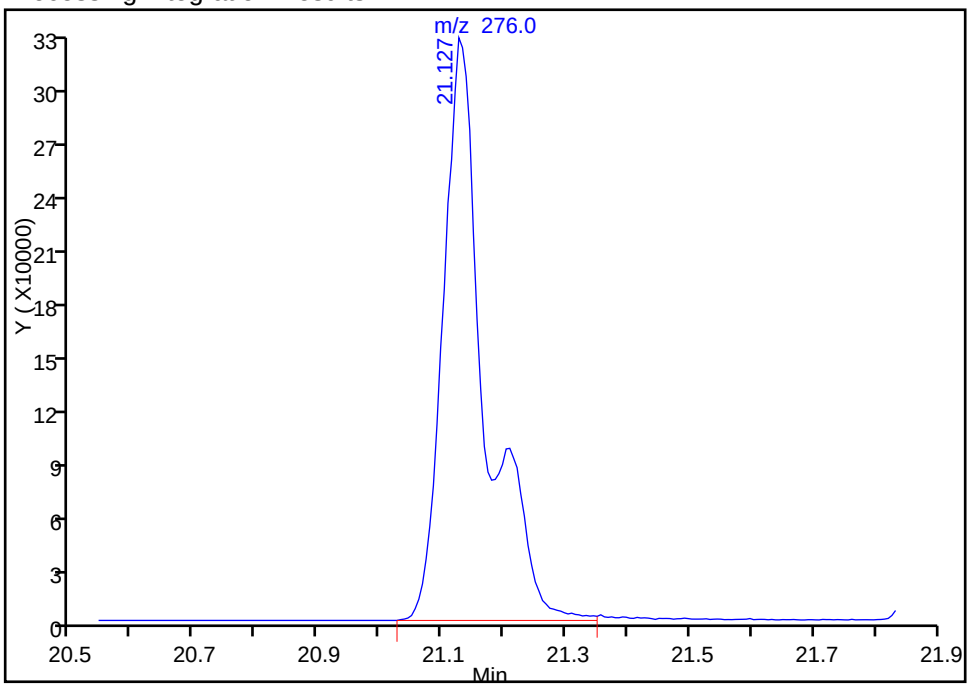
TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16645.D
Injection Date: 25-Feb-2015 15:26:30 Instrument ID: SMS_G6
Lims ID: ICV HSL 1
Client ID:
Operator ID: KIEKELD ALS Bottle#: 10 Worklist Smp#: 11
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Method: SMS_G6_8270D Limit Group: MSSV - 8270D
Column: VF-5ms (0.50 mm) Detector MS SCAN

151 Indeno[1,2,3-cd]pyrene, CAS: 193-39-5

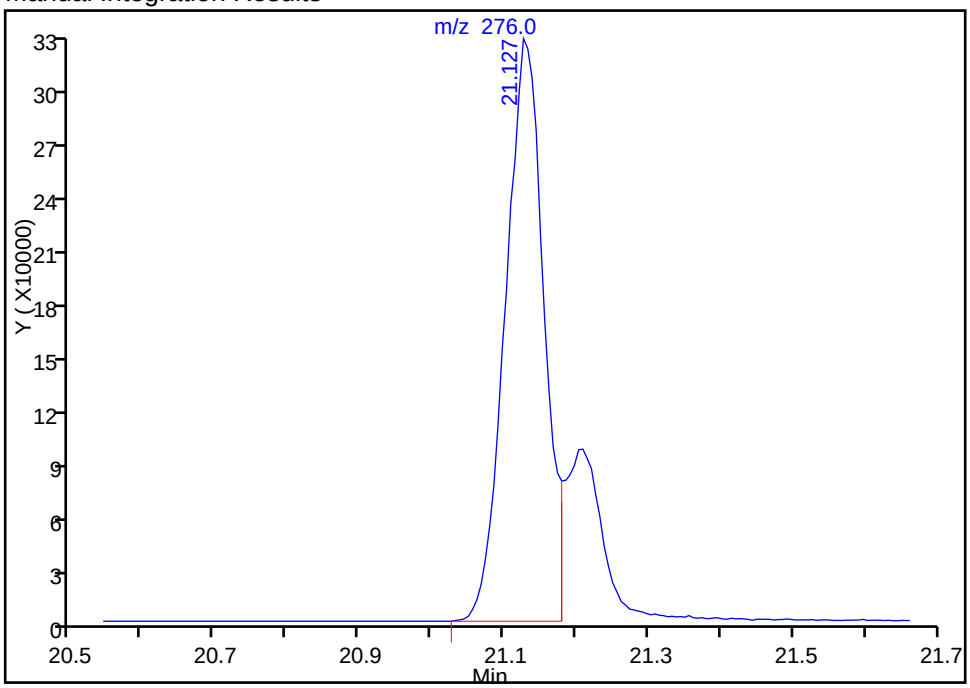
Processing Integration Results

RT: 21.13
Area: 1550066
Amount: 121.3925
Amount Units: ug/ml



Manual Integration Results

RT: 21.13
Area: 1219599
Amount: 95.512153
Amount Units: ug/ml



Reviewer: hoeflera, 10-Apr-2015 09:24:38
Audit Action: Split an Integrated Peak
Audit Reason: Shouldering

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: ICV 280-272059/12 Calibration Date: 02/25/2015 15:52

Instrument ID: SMS_G6 Calib Start Date: 05/14/2014 14:05

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 05/14/2014 17:05

Lab File ID: G6_16646.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Benzidine	Ave	0.6743			50.0	100		30.0

Report Date: 10-Apr-2015 12:09:46 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16646.D
Lims ID: ICV HSL 2
Client ID:
Sample Type: ICV
Inject. Date: 25-Feb-2015 15:52:30 ALS Bottle#: 11 Worklist Smp#: 12
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: ICV HSL 2
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist:

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:44 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:25:41

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	97	185167	40.0	40.0	
* 2 Naphthalene-d8	136	6.022	6.022	0.000	100	715813	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	93	400054	40.0	40.0	
* 4 Phenanthrene-d10	188	9.280	9.280	0.000	97	674469	40.0	40.0	
* 5 Chrysene-d12	240	13.698	13.710	-0.012	96	649859	40.0	40.0	
* 6 Perylene-d12	264	17.898	17.903	-0.005	95	546537	40.0	40.0	
62 Caprolactam	55	6.416	6.428	-0.012	78	352151	100.0	104.0	
129 Benzidine	184		10.466				ND	ND	
140 3,3'-Dichlorobenzidine	252	13.639	13.639	0.000	76	707262	100.0	110.1	

QC Flag Legend

Processing Flags
ND - Not Detected or Marked ND

Reagents:

MS-HSLB2SSV_00020 Amount Added: 200.00 Units: uL

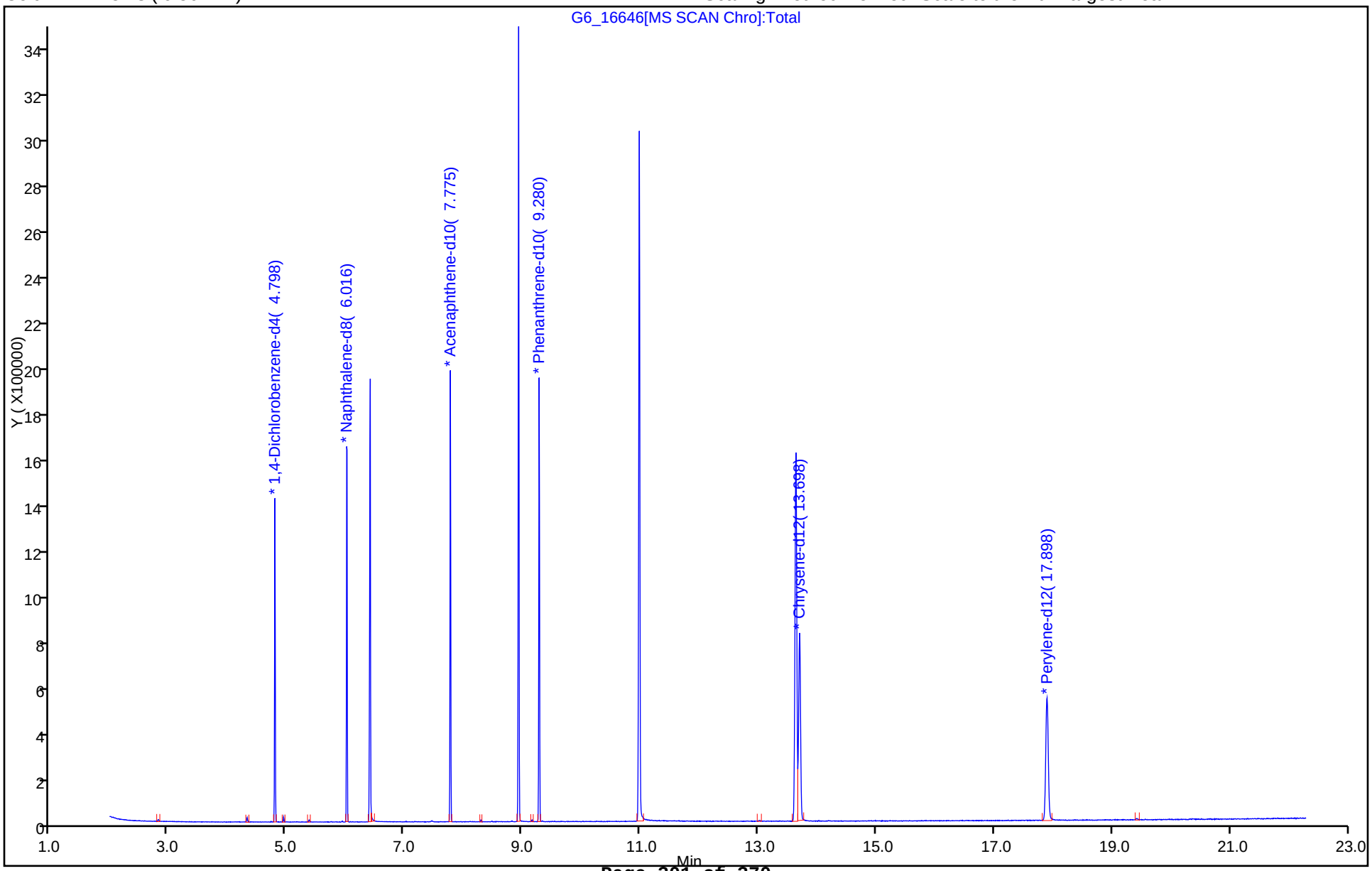
Report Date: 10-Apr-2015 12:09:46 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16646.D

Injection Date:	25-Feb-2015 15:52:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	ICV HSL 2			Worklist Smp#:	12
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	11
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: ICV 280-272059/12 Calibration Date: 02/25/2015 15:52

Instrument ID: SMS_G6 Calib Start Date: 02/25/2015 11:53

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 02/25/2015 14:59

Lab File ID: G6_16646.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Caprolactam	Ave	0.1892	0.1968		104	100	4.0	30.0
3,3'-Dichlorobenzidine	Ave	0.3953	0.4353	0.0100	110	100	10.1	30.0

Report Date: 10-Apr-2015 12:09:46 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16646.D
Lims ID: ICV HSL 2
Client ID:
Sample Type: ICV
Inject. Date: 25-Feb-2015 15:52:30 ALS Bottle#: 11 Worklist Smp#: 12
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: ICV HSL 2
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist:

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:44 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:25:41

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	97	185167	40.0	40.0	
* 2 Naphthalene-d8	136	6.022	6.022	0.000	100	715813	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	93	400054	40.0	40.0	
* 4 Phenanthrene-d10	188	9.280	9.280	0.000	97	674469	40.0	40.0	
* 5 Chrysene-d12	240	13.698	13.710	-0.012	96	649859	40.0	40.0	
* 6 Perylene-d12	264	17.898	17.903	-0.005	95	546537	40.0	40.0	
62 Caprolactam	55	6.416	6.428	-0.012	78	352151	100.0	104.0	
129 Benzidine	184		10.466				ND	ND	
140 3,3'-Dichlorobenzidine	252	13.639	13.639	0.000	76	707262	100.0	110.1	

QC Flag Legend

Processing Flags
ND - Not Detected or Marked ND

Reagents:

MS-HSLB2SSV_00020 Amount Added: 200.00 Units: uL

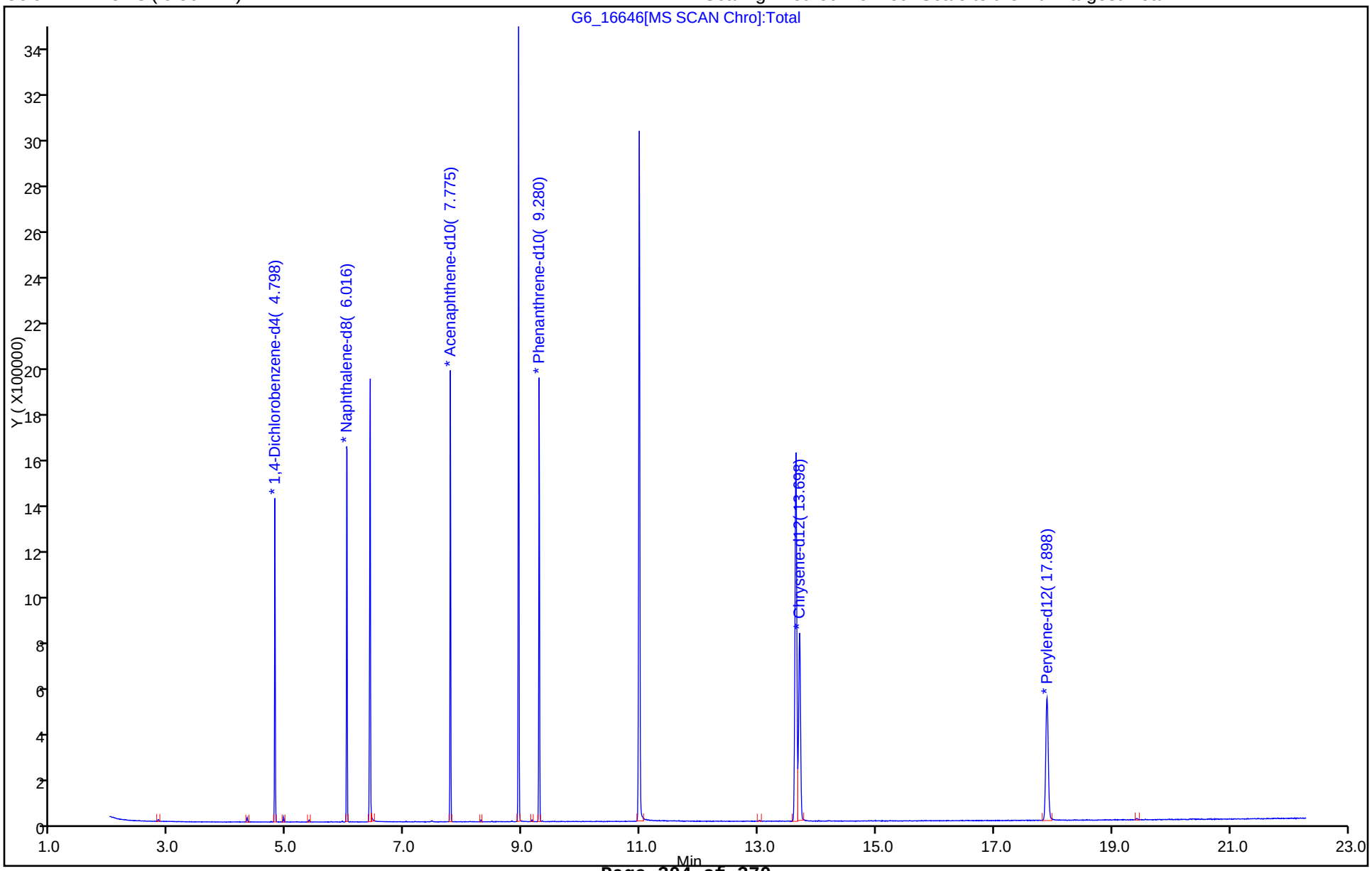
Report Date: 10-Apr-2015 12:09:46 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16646.D

Injection Date:	25-Feb-2015 15:52:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	ICV HSL 2			Worklist Smp#:	12
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	11
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: ICV 280-272059/13 Calibration Date: 02/25/2015 16:19

Instrument ID: SMS_G6 Calib Start Date: 02/25/2015 11:53

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 02/25/2015 14:59

Lab File ID: G6_16647.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Famphur	Ave	0.4215	0.4268		101	100	1.2	30.0

Report Date: 10-Apr-2015 12:09:47 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16647.D
Lims ID: ICV FAM
Client ID:
Sample Type: ICV
Inject. Date: 25-Feb-2015 16:19:30 ALS Bottle#: 12 Worklist Smp#: 13
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: ICV FAM
Operator ID: KIEKELD Instrument ID: SMS_G6
Sublist:

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:44 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: hoeflera Date: 10-Apr-2015 09:26:30

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.798	4.798	0.000	97	186013	40.0	40.0	
* 2 Naphthalene-d8	136	6.016	6.022	-0.006	100	710514	40.0	40.0	
* 3 Acenaphthene-d10	164	7.775	7.775	0.000	92	389552	40.0	40.0	
* 4 Phenanthrene-d10	188	9.281	9.280	0.001	98	665613	40.0	40.0	
* 5 Chrysene-d12	240	13.704	13.710	-0.006	96	633484	40.0	40.0	
* 6 Perylene-d12	264	17.904	17.903	0.001	95	533431	40.0	40.0	
136 Famphur	218	12.204	12.198	0.006	97	675896	100.0	101.2	

Reagents:

MS-FAMSSV_100_00013 Amount Added: 200.00 Units: uL

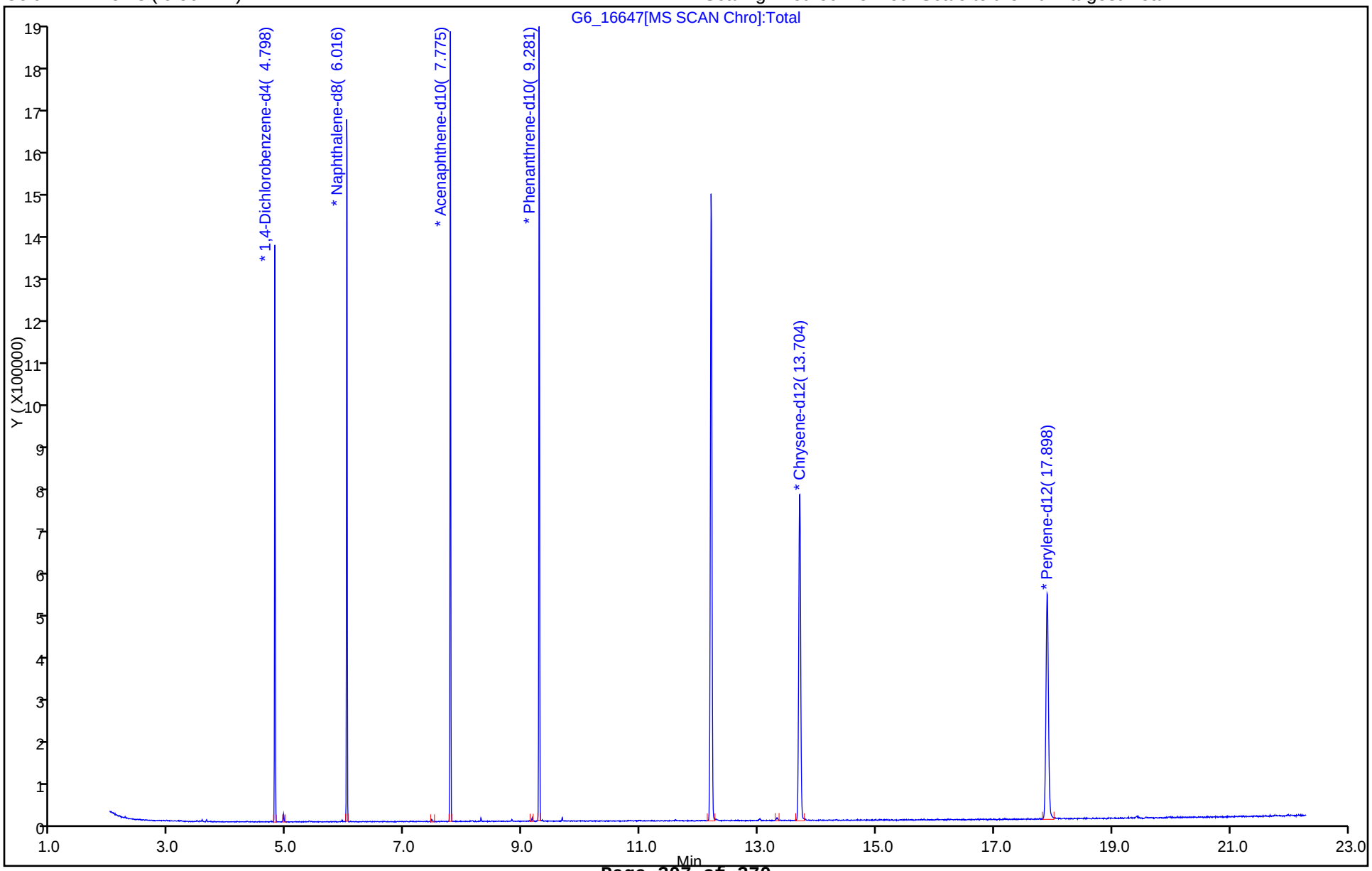
Report Date: 10-Apr-2015 12:09:47 Chrom Revision: 2.2 13-Mar-2015 11:20:44

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16647.D

Injection Date:	25-Feb-2015 16:19:30	Instrument ID:	SMS_G6	Operator ID:	KIEKELD
Lims ID:	ICV FAM			Worklist Smp#:	13
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	12
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: CCV 280-273380/3 Calibration Date: 04/18/2015 15:20

Instrument ID: SMS_G6 Calib Start Date: 05/14/2014 14:05

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 05/14/2014 17:05

Lab File ID: G6_17414.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Benzidine	Ave	0.6743			50.0	80.0		

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17414.D
Lims ID: CCV HSL
Client ID:
Sample Type: CCV
Inject. Date: 18-Apr-2015 15:20:30 ALS Bottle#: 2 Worklist Smp#: 3
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: CCV HSL
Operator ID: HOEFLERA Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:11 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: hoeflera Date: 18-Apr-2015 16:20:23

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	96	216765	40.0	40.0	
* 2 Naphthalene-d8	136	5.951	5.951	0.000	100	822108	40.0	40.0	
* 3 Acenaphthene-d10	164	7.710	7.710	0.000	92	444568	40.0	40.0	
* 4 Phenanthrene-d10	188	9.210	9.210	0.000	97	728379	40.0	40.0	
* 5 Chrysene-d12	240	13.533	13.533	0.000	96	645212	40.0	40.0	
* 6 Perylene-d12	264	17.662	17.662	0.000	95	551738	40.0	40.0	
$ 7 2-Fluorophenol	112	3.493	3.493	0.000	92	662801	80.0	84.1	
$ 8 Phenol-d5	99	4.334	4.334	0.000	98	786514	80.0	79.6	
$ 9 Nitrobenzene-d5	82	5.245	5.245	0.000	91	686359	80.0	81.4	
$ 10 2-Fluorobiphenyl	172	7.004	7.004	0.000	100	1217564	80.0	81.0	
$ 11 2,4,6-Tribromophenol	330	8.504	8.504	0.000	87	142497	80.0	85.2	
$ 12 Terphenyl-d14	244	11.239	11.239	0.000	98	1088368	80.0	84.9	
13 1,4-Dioxane	88	2.040	2.040	0.000	97	290423	80.0	87.3	
14 N-Nitrosodimethylamine	74	2.305	2.305	0.000	92	443847	80.0	87.9	
15 Pyridine	79	2.352	2.352	0.000	91	720818	80.0	83.7	
23 Phenol	94	4.346	4.346	0.000	99	780029	80.0	78.7	
24 Aniline	93	4.404	4.404	0.000	99	964440	80.0	79.6	
25 Bis(2-chloroethyl)ether	93	4.434	4.434	0.000	93	607225	80.0	78.1	
26 2-Chlorophenol	128	4.516	4.516	0.000	96	660959	80.0	81.7	
27 1,3-Dichlorobenzene	146	4.657	4.657	0.000	97	666683	80.0	81.2	
28 1,4-Dichlorobenzene	146	4.728	4.728	0.000	94	677917	80.0	81.8	
29 Benzyl alcohol	108	4.828	4.828	0.000	94	419923	80.0	82.4	
30 1,2-Dichlorobenzene	146	4.875	4.875	0.000	96	647653	80.0	81.2	
31 2-Methylphenol	108	4.916	4.916	0.000	95	583382	80.0	79.4	
32 2,2'-oxybis[1-chloropropan	45	4.934	4.934	0.000	94	869271	80.0	75.9	
38 3 & 4 Methylphenol	108	5.063	5.063	0.000	96	598915	80.0	78.7	
39 3-Methylphenol	108	5.063	5.063	0.000	96	598915	80.0	78.7	
40 4-Methylphenol	108	5.063	5.063	0.000	96	598915	80.0	78.7	
41 N-Nitrosodi-n-propylamine	70	5.069	5.069	0.000	88	430335	80.0	77.9	
42 Acetophenone	105	5.087	5.087	0.000	98	870713	80.0	81.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.198	5.198	0.000	97	293300	80.0	82.4	
44 Nitrobenzene	77	5.263	5.263	0.000	91	655413	80.0	80.7	
46 Isophorone	82	5.475	5.475	0.000	100	1123098	80.0	81.4	
48 2-Nitrophenol	139	5.569	5.569	0.000	95	340452	80.0	84.1	
49 2,4-Dimethylphenol	107	5.581	5.581	0.000	95	590548	80.0	79.6	
50 Bis(2-chloroethoxy)methane	93	5.669	5.669	0.000	98	736507	80.0	81.2	
52 Benzoic acid	105	5.687	5.687	0.000	89	935797	160.0	160.8	
53 2,4-Dichlorophenol	162	5.798	5.798	0.000	97	503006	80.0	82.8	
54 1,2,4-Trichlorobenzene	180	5.887	5.887	0.000	94	507924	80.0	82.3	
57 Naphthalene	128	5.975	5.975	0.000	97	1782439	80.0	81.5	
58 4-Chloroaniline	127	6.016	6.016	0.000	97	817016	80.0	82.6	
59 Hexachlorobutadiene	225	6.069	6.069	0.000	98	278432	80.0	82.8	
62 Caprolactam	55	6.369	6.369	0.000	79	309525	80.0	79.6	
64 4-Chloro-3-methylphenol	107	6.481	6.481	0.000	96	516463	80.0	79.4	
65 2-Methylnaphthalene	142	6.651	6.651	0.000	93	1132023	80.0	80.6	
67 1-Methylnaphthalene	142	6.757	6.757	0.000	94	1075001	80.0	81.3	
68 Hexachlorocyclopentadiene	237	6.798	6.798	0.000	96	261959	80.0	75.0	
69 1,2,4,5-Tetrachlorobenzene	216	6.822	6.822	0.000	98	474419	80.0	81.1	
70 2,4,6-Trichlorophenol	196	6.934	6.934	0.000	94	334561	80.0	83.8	
72 2,4,5-Trichlorophenol	196	6.969	6.969	0.000	94	358741	80.0	84.4	
74 1,1'-Biphenyl	154	7.110	7.110	0.000	95	1365544	80.0	81.6	
75 2-Chloronaphthalene	162	7.151	7.151	0.000	97	1040634	80.0	82.5	
77 2-Nitroaniline	65	7.251	7.251	0.000	85	372400	80.0	80.4	
79 Dimethyl phthalate	163	7.398	7.398	0.000	98	1169708	80.0	79.0	
80 1,3-Dinitrobenzene	168	7.463	7.463	0.000	91	237829	80.0	90.6	
81 2,6-Dinitrotoluene	165	7.481	7.481	0.000	94	287368	80.0	84.0	
82 Acenaphthylene	152	7.569	7.569	0.000	98	1693447	80.0	82.7	
83 3-Nitroaniline	138	7.663	7.663	0.000	95	374298	80.0	84.6	
84 Acenaphthene	153	7.739	7.739	0.000	95	1070647	80.0	81.0	
86 2,4-Dinitrophenol	184	7.763	7.763	0.000	82	412726	160.0	166.7	
87 4-Nitrophenol	109	7.816	7.816	0.000	94	369345	160.0	154.9	
89 2,4-Dinitrotoluene	165	7.887	7.887	0.000	92	389897	80.0	86.9	
90 Dibenzofuran	168	7.910	7.910	0.000	97	1516485	80.0	81.4	
92 2,3,4,6-Tetrachlorophenol	232	8.028	8.028	0.000	75	289462	80.0	84.0	
94 Diethyl phthalate	149	8.092	8.092	0.000	98	1239779	80.0	82.3	
96 4-Chlorophenyl phenyl ethe	204	8.234	8.234	0.000	93	546791	80.0	81.1	
98 Fluorene	166	8.257	8.257	0.000	96	1241761	80.0	81.9	
99 4-Nitroaniline	138	8.286	8.286	0.000	87	373129	80.0	86.3	
100 4,6-Dinitro-2-methylphenol	198	8.298	8.298	0.000	85	468101	160.0	181.6	
102 N-Nitrosodiphenylamine	169	8.357	8.357	0.000	61	930374	80.0	86.2	
103 Azobenzene	77	8.392	8.392	0.000	100	1326025	80.0	79.6	
104 1,2-Diphenylhydrazine	77	8.392	8.392	0.000	100	1326025	80.9	80.5	
111 4-Bromophenyl phenyl ether	248	8.728	8.728	0.000	70	288713	80.0	84.6	
112 Hexachlorobenzene	284	8.810	8.810	0.000	92	296937	80.0	83.8	
116 Pentachlorophenol	266	9.010	9.010	0.000	89	378615	160.0	162.5	
119 Phenanthrene	178	9.233	9.233	0.000	98	1673042	80.0	82.8	
120 Anthracene	178	9.286	9.286	0.000	98	1733749	80.0	83.8	
122 Carbazole	167	9.445	9.445	0.000	95	1760082	80.0	83.4	
123 Di-n-butyl phthalate	149	9.739	9.739	0.000	100	2112316	80.0	85.1	
128 Fluoranthene	202	10.675	10.675	0.000	98	1791559	80.0	81.0	
129 Benzidine	184		10.466				ND	ND	
131 Pyrene	202	11.039	11.039	0.000	97	1855215	80.0	86.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
136 Famphur	218	12.051	12.051	0.000	97	594971	80.0	87.5	
137 Butyl benzyl phthalate	149	12.151	12.151	0.000	97	928375	80.0	91.3	
140 3,3'-Dichlorobenzidine	252	13.468	13.468	0.000	75	559891	80.0	87.8	
141 Benzo[a]anthracene	228	13.504	13.504	0.000	99	1533206	80.0	82.7	
142 Bis(2-ethylhexyl) phthalat	149	13.492	13.492	0.000	97	1195540	80.0	90.0	
143 Chrysene	228	13.598	13.598	0.000	98	1461997	80.0	84.4	
144 Di-n-octyl phthalate	149	15.362	15.362	0.000	99	2125005	80.0	89.6	
146 Benzo[b]fluoranthene	252	16.509	16.509	0.000	99	1376089	80.0	85.4	
147 Benzo[k]fluoranthene	252	16.592	16.592	0.000	99	1396697	80.0	84.4	
148 Benzo[a]pyrene	252	17.492	17.492	0.000	80	1347455	80.0	87.0	
151 Indeno[1,2,3-cd]pyrene	276	20.862	20.862	0.000	96	1096109	80.0	83.2	M
152 Dibenz(a,h)anthracene	278	20.939	20.939	0.000	94	1111518	80.0	86.8	
153 Benzo[g,h,i]perylene	276	21.633	21.633	0.000	95	1166748	80.0	83.7	

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

MS-HSLACCV080_00038 Amount Added: 200.00 Units: uL

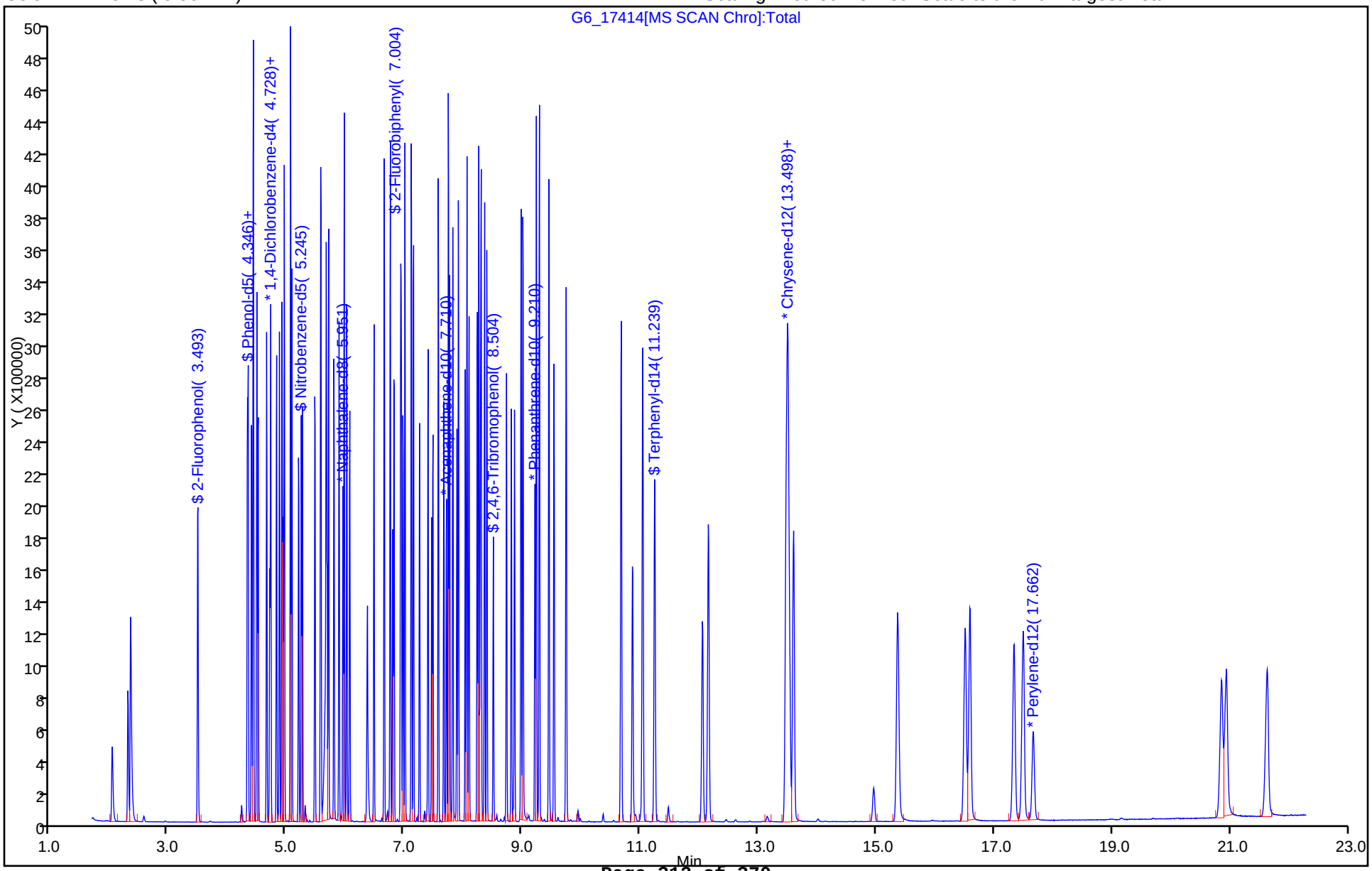
Report Date: 20-Apr-2015 07:09:12 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17414.D

Injection Date:	18-Apr-2015 15:20:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	CCV HSL			Worklist Smp#:	3
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	2
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: CCV 280-273380/3 Calibration Date: 04/18/2015 15:20

Instrument ID: SMS_G6 Calib Start Date: 02/25/2015 11:53

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 02/25/2015 14:59

Lab File ID: G6_17414.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
1,4-Dioxane	Ave	0.6137	0.6699		87.3	80.0	9.2	20.0
N-Nitrosodimethylamine	Ave	0.9321	1.024		87.9	80.0	9.8	20.0
Pyridine	Ave	1.590	1.663		83.7	80.0	4.6	20.0
Phenol	Ave	1.828	1.799	0.8000	78.7	80.0	-1.6	20.0
Aniline	Ave	2.235	2.225		79.6	80.0	-0.4	20.0
Bis(2-chloroethyl)ether	Ave	1.435	1.401	0.7000	78.1	80.0	-2.4	20.0
2-Chlorophenol	Ave	1.492	1.525	0.8000	81.7	80.0	2.2	20.0
1,3-Dichlorobenzene	Ave	1.514	1.538		81.2	80.0	1.5	20.0
1,4-Dichlorobenzene	Ave	1.530	1.564		81.8	80.0	2.2	20.0
Benzyl alcohol	Ave	0.9402	0.9686		82.4	80.0	3.0	20.0
1,2-Dichlorobenzene	Ave	1.471	1.494		81.2	80.0	1.5	20.0
2-Methylphenol	Ave	1.355	1.346	0.7000	79.4	80.0	-0.7	20.0
bis (2-chloroisopropyl) ether	Ave	2.112	2.005	0.0100	75.9	80.0	-5.1	20.0
3 & 4 Methylphenol	Ave	1.405	1.381		78.7	80.0	-1.7	20.0
3-Methylphenol	Ave	1.405	1.381		78.7	80.0	-1.7	20.0
4-Methylphenol	Ave	1.405	1.381	0.6000	78.7	80.0	-1.7	20.0
N-Nitrosodi-n-propylamine	Ave	1.020	0.9926	0.5000	77.9	80.0	-2.6	20.0
Acetophenone	Ave	1.974	2.008	0.0100	81.4	80.0	1.7	20.0
Hexachloroethane	Ave	0.6565	0.6765	0.3000	82.4	80.0	3.0	20.0
Nitrobenzene	Ave	0.3951	0.3986		80.7	80.0	0.9	20.0
Isophorone	Ave	0.6711	0.6831	0.4000	81.4	80.0	1.8	20.0
2-Nitrophenol	Ave	0.1969	0.2071	0.1000	84.1	80.0	5.2	20.0
2,4-Dimethylphenol	Ave	0.3608	0.3592	0.2000	79.6	80.0	-0.5	20.0
Bis(2-chloroethoxy)methane	Ave	0.4411	0.4479	0.3000	81.2	80.0	1.6	20.0
Benzoic acid	Ave	0.2831	0.2846		161	160	0.5	20.0
2,4-Dichlorophenol	Ave	0.2955	0.3059	0.2000	82.8	80.0	3.5	20.0
1,2,4-Trichlorobenzene	Ave	0.3003	0.3089		82.3	80.0	2.9	20.0
Naphthalene	Ave	1.064	1.084	0.7000	81.5	80.0	1.9	20.0
4-Chloroaniline	Ave	0.4813	0.4969	0.0100	82.6	80.0	3.3	20.0
Hexachlorobutadiene	Ave	0.1636	0.1693	0.0100	82.8	80.0	3.5	20.0
Caprolactam	Ave	0.1892	0.1883		79.6	80.0	-0.5	20.0
4-Chloro-3-methylphenol	Ave	0.3163	0.3141	0.2000	79.4	80.0	-0.7	20.0
2-Methylnaphthalene	Ave	0.6836	0.6885	0.4000	80.6	80.0	0.7	20.0
1-Methylnaphthalene	Ave	0.6430	0.6538		81.3	80.0	1.7	20.0
Hexachlorocyclopentadiene	Ave	0.3143	0.2946	0.0500	75.0	80.0	-6.3	20.0
1,2,4,5-Tetrachlorobenzene	Ave	0.2846	0.2885	0.0100	81.1	80.0	1.4	20.0
2,4,6-Trichlorophenol	Ave	0.3592	0.3763	0.2000	83.8	80.0	4.8	20.0
2,4,5-Trichlorophenol	Ave	0.3826	0.4035	0.2000	84.4	80.0	5.5	20.0
1,1'-Biphenyl	Ave	1.506	1.536		81.6	80.0	2.0	20.0
2-Chloronaphthalene	Ave	1.135	1.170	0.8000	82.5	80.0	3.1	20.0

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: CCV 280-273380/3 Calibration Date: 04/18/2015 15:20

Instrument ID: SMS_G6 Calib Start Date: 02/25/2015 11:53

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 02/25/2015 14:59

Lab File ID: G6_17414.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
2-Nitroaniline	Ave	0.4168	0.4188	0.0100	80.4	80.0	0.5	20.0
Dimethyl phthalate	Ave	1.332	1.316	0.0100	79.0	80.0	-1.2	20.0
1,3-Dinitrobenzene	Ave	0.2362	0.2675		90.6	80.0	13.2	20.0
2,6-Dinitrotoluene	Ave	0.3078	0.3232	0.2000	84.0	80.0	5.0	20.0
Acenaphthylene	Ave	1.843	1.905	0.9000	82.7	80.0	3.3	20.0
3-Nitroaniline	Ave	0.3980	0.4210	0.0100	84.6	80.0	5.8	20.0
Acenaphthene	Ave	1.190	1.204	0.9000	81.0	80.0	1.2	20.0
2,4-Dinitrophenol	Lin2		0.2321	0.0100	167	160	4.2	20.0
4-Nitrophenol	Ave	0.2146	0.2077	0.0100	155	160	-3.2	20.0
2,4-Dinitrotoluene	Ave	0.4036	0.4385	0.2000	86.9	80.0	8.7	20.0
Dibenzofuran	Ave	1.676	1.706	0.8000	81.4	80.0	1.8	20.0
2,3,4,6-Tetrachlorophenol	Ave	0.3102	0.3256	0.0100	84.0	80.0	5.0	20.0
Diethyl phthalate	Ave	1.356	1.394	0.0100	82.3	80.0	2.9	20.0
4-Chlorophenyl phenyl ether	Ave	0.6065	0.6150	0.4000	81.1	80.0	1.4	20.0
Fluorene	Ave	1.364	1.397	0.9000	81.9	80.0	2.4	20.0
4-Nitroaniline	Ave	0.3891	0.4197	0.0100	86.3	80.0	7.8	20.0
4,6-Dinitro-2-methylphenol	Ave	0.1415	0.1607	0.0100	182	160	13.5	20.0
N-Nitrosodiphenylamine	Ave	0.5925	0.6387	0.0100	86.2	80.0	7.8	20.0
1,2-Diphenylhydrazine(as Azobenzene)	Ave	1.482	1.475		80.5	80.9	-0.5	20.0
Azobenzene	Ave	1.498	1.491		79.6	80.0	-0.5	20.0
4-Bromophenyl phenyl ether	Ave	0.1874	0.1982	0.1000	84.6	80.0	5.7	20.0
Hexachlorobenzene	Ave	0.1947	0.2038	0.1000	83.8	80.0	4.7	20.0
Pentachlorophenol	Ave	0.1280	0.1300	0.0500	162	160	1.5	20.0
Phenanthrene	Ave	1.109	1.148	0.7000	82.8	80.0	3.5	20.0
Anthracene	Ave	1.137	1.190	0.7000	83.8	80.0	4.7	20.0
Carbazole	Ave	1.159	1.208	0.0100	83.4	80.0	4.3	20.0
Di-n-butyl phthalate	Ave	1.363	1.450	0.0100	85.1	80.0	6.4	20.0
Fluoranthene	Ave	1.215	1.230	0.6000	81.0	80.0	1.2	20.0
Pyrene	Ave	1.336	1.438	0.6000	86.1	80.0	7.6	20.0
Famphur	Ave	0.4215	0.4611		87.5	80.0	9.4	20.0
Butyl benzyl phthalate	Ave	0.6306	0.7194	0.0100	91.3	80.0	14.1	20.0
3,3'-Dichlorobenzidine	Ave	0.3953	0.4339	0.0100	87.8	80.0	9.8	20.0
Bis(2-ethylhexyl) phthalate	Ave	0.8237	0.9265	0.0100	90.0	80.0	12.5	20.0
Benzo[a]anthracene	Ave	1.150	1.188	0.8000	82.7	80.0	3.3	20.0
Chrysene	Ave	1.074	1.133	0.7000	84.4	80.0	5.5	20.0
Di-n-octyl phthalate	Lin2		1.647	0.0100	89.6	80.0	12.0	20.0
Benzo[b]fluoranthene	Ave	1.169	1.247	0.7000	85.4	80.0	6.7	20.0
Benzo[k]fluoranthene	Ave	1.200	1.266	0.7000	84.4	80.0	5.5	20.0
Benzo[a]pyrene	Ave	1.123	1.221	0.7000	87.0	80.0	8.7	20.0
Indeno[1,2,3-cd]pyrene	Ave	0.8166	0.8494	0.5000	83.2	80.0	4.0	20.0

FORM VII
GC/MS SEMI VOA CONTINUING CALIBRATION DATA

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Lab Sample ID: CCV 280-273380/3 Calibration Date: 04/18/2015 15:20

Instrument ID: SMS_G6 Calib Start Date: 02/25/2015 11:53

GC Column: Vf-5MS (30.25) ID: 0.25(mm) Calib End Date: 02/25/2015 14:59

Lab File ID: G6_17414.D Conc. Units: ug/mL

ANALYTE	CURVE TYPE	AVE RRF	RRF	MIN RRF	CALC AMOUNT	SPIKE AMOUNT	%D	MAX %D
Dibenz(a,h)anthracene	Ave	0.9288	1.007	0.4000	86.8	80.0	8.4	20.0
Benzo[g,h,i]perylene	Ave	1.011	1.057	0.5000	83.7	80.0	4.6	20.0
2-Fluorophenol (Surr)	Ave	1.454	1.529		84.1	80.0	5.1	20.0
Phenol-d5 (Surr)	Ave	1.823	1.814		79.6	80.0	-0.5	20.0
Nitrobenzene-d5 (Surr)	Ave	0.4101	0.4174		81.4	80.0	1.8	20.0
2-Fluorobiphenyl	Ave	1.352	1.369		81.0	80.0	1.3	20.0
2,4,6-Tribromophenol (Surr)	Ave	0.1504	0.1603		85.2	80.0	6.5	20.0
Terphenyl-d14 (Surr)	Ave	0.7949	0.8434		84.9	80.0	6.1	20.0

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17414.D
Lims ID: CCV HSL
Client ID:
Sample Type: CCV
Inject. Date: 18-Apr-2015 15:20:30 ALS Bottle#: 2 Worklist Smp#: 3
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: CCV HSL
Operator ID: HOEFLERA Instrument ID: SMS_G6
Sublist: chrom-SMS_G6_8270D*sub6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:11 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: hoeflera Date: 18-Apr-2015 16:20:23

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	96	216765	40.0	40.0	
* 2 Naphthalene-d8	136	5.951	5.951	0.000	100	822108	40.0	40.0	
* 3 Acenaphthene-d10	164	7.710	7.710	0.000	92	444568	40.0	40.0	
* 4 Phenanthrene-d10	188	9.210	9.210	0.000	97	728379	40.0	40.0	
* 5 Chrysene-d12	240	13.533	13.533	0.000	96	645212	40.0	40.0	
* 6 Perylene-d12	264	17.662	17.662	0.000	95	551738	40.0	40.0	
$ 7 2-Fluorophenol	112	3.493	3.493	0.000	92	662801	80.0	84.1	
$ 8 Phenol-d5	99	4.334	4.334	0.000	98	786514	80.0	79.6	
$ 9 Nitrobenzene-d5	82	5.245	5.245	0.000	91	686359	80.0	81.4	
$ 10 2-Fluorobiphenyl	172	7.004	7.004	0.000	100	1217564	80.0	81.0	
$ 11 2,4,6-Tribromophenol	330	8.504	8.504	0.000	87	142497	80.0	85.2	
$ 12 Terphenyl-d14	244	11.239	11.239	0.000	98	1088368	80.0	84.9	
13 1,4-Dioxane	88	2.040	2.040	0.000	97	290423	80.0	87.3	
14 N-Nitrosodimethylamine	74	2.305	2.305	0.000	92	443847	80.0	87.9	
15 Pyridine	79	2.352	2.352	0.000	91	720818	80.0	83.7	
23 Phenol	94	4.346	4.346	0.000	99	780029	80.0	78.7	
24 Aniline	93	4.404	4.404	0.000	99	964440	80.0	79.6	
25 Bis(2-chloroethyl)ether	93	4.434	4.434	0.000	93	607225	80.0	78.1	
26 2-Chlorophenol	128	4.516	4.516	0.000	96	660959	80.0	81.7	
27 1,3-Dichlorobenzene	146	4.657	4.657	0.000	97	666683	80.0	81.2	
28 1,4-Dichlorobenzene	146	4.728	4.728	0.000	94	677917	80.0	81.8	
29 Benzyl alcohol	108	4.828	4.828	0.000	94	419923	80.0	82.4	
30 1,2-Dichlorobenzene	146	4.875	4.875	0.000	96	647653	80.0	81.2	
31 2-Methylphenol	108	4.916	4.916	0.000	95	583382	80.0	79.4	
32 2,2'-oxybis[1-chloropropan	45	4.934	4.934	0.000	94	869271	80.0	75.9	
38 3 & 4 Methylphenol	108	5.063	5.063	0.000	96	598915	80.0	78.7	
39 3-Methylphenol	108	5.063	5.063	0.000	96	598915	80.0	78.7	
40 4-Methylphenol	108	5.063	5.063	0.000	96	598915	80.0	78.7	
41 N-Nitrosodi-n-propylamine	70	5.069	5.069	0.000	88	430335	80.0	77.9	
42 Acetophenone	105	5.087	5.087	0.000	98	870713	80.0	81.4	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.198	5.198	0.000	97	293300	80.0	82.4	
44 Nitrobenzene	77	5.263	5.263	0.000	91	655413	80.0	80.7	
46 Isophorone	82	5.475	5.475	0.000	100	1123098	80.0	81.4	
48 2-Nitrophenol	139	5.569	5.569	0.000	95	340452	80.0	84.1	
49 2,4-Dimethylphenol	107	5.581	5.581	0.000	95	590548	80.0	79.6	
50 Bis(2-chloroethoxy)methane	93	5.669	5.669	0.000	98	736507	80.0	81.2	
52 Benzoic acid	105	5.687	5.687	0.000	89	935797	160.0	160.8	
53 2,4-Dichlorophenol	162	5.798	5.798	0.000	97	503006	80.0	82.8	
54 1,2,4-Trichlorobenzene	180	5.887	5.887	0.000	94	507924	80.0	82.3	
57 Naphthalene	128	5.975	5.975	0.000	97	1782439	80.0	81.5	
58 4-Chloroaniline	127	6.016	6.016	0.000	97	817016	80.0	82.6	
59 Hexachlorobutadiene	225	6.069	6.069	0.000	98	278432	80.0	82.8	
62 Caprolactam	55	6.369	6.369	0.000	79	309525	80.0	79.6	
64 4-Chloro-3-methylphenol	107	6.481	6.481	0.000	96	516463	80.0	79.4	
65 2-Methylnaphthalene	142	6.651	6.651	0.000	93	1132023	80.0	80.6	
67 1-Methylnaphthalene	142	6.757	6.757	0.000	94	1075001	80.0	81.3	
68 Hexachlorocyclopentadiene	237	6.798	6.798	0.000	96	261959	80.0	75.0	
69 1,2,4,5-Tetrachlorobenzene	216	6.822	6.822	0.000	98	474419	80.0	81.1	
70 2,4,6-Trichlorophenol	196	6.934	6.934	0.000	94	334561	80.0	83.8	
72 2,4,5-Trichlorophenol	196	6.969	6.969	0.000	94	358741	80.0	84.4	
74 1,1'-Biphenyl	154	7.110	7.110	0.000	95	1365544	80.0	81.6	
75 2-Chloronaphthalene	162	7.151	7.151	0.000	97	1040634	80.0	82.5	
77 2-Nitroaniline	65	7.251	7.251	0.000	85	372400	80.0	80.4	
79 Dimethyl phthalate	163	7.398	7.398	0.000	98	1169708	80.0	79.0	
80 1,3-Dinitrobenzene	168	7.463	7.463	0.000	91	237829	80.0	90.6	
81 2,6-Dinitrotoluene	165	7.481	7.481	0.000	94	287368	80.0	84.0	
82 Acenaphthylene	152	7.569	7.569	0.000	98	1693447	80.0	82.7	
83 3-Nitroaniline	138	7.663	7.663	0.000	95	374298	80.0	84.6	
84 Acenaphthene	153	7.739	7.739	0.000	95	1070647	80.0	81.0	
86 2,4-Dinitrophenol	184	7.763	7.763	0.000	82	412726	160.0	166.7	
87 4-Nitrophenol	109	7.816	7.816	0.000	94	369345	160.0	154.9	
89 2,4-Dinitrotoluene	165	7.887	7.887	0.000	92	389897	80.0	86.9	
90 Dibenzofuran	168	7.910	7.910	0.000	97	1516485	80.0	81.4	
92 2,3,4,6-Tetrachlorophenol	232	8.028	8.028	0.000	75	289462	80.0	84.0	
94 Diethyl phthalate	149	8.092	8.092	0.000	98	1239779	80.0	82.3	
96 4-Chlorophenyl phenyl ethe	204	8.234	8.234	0.000	93	546791	80.0	81.1	
98 Fluorene	166	8.257	8.257	0.000	96	1241761	80.0	81.9	
99 4-Nitroaniline	138	8.286	8.286	0.000	87	373129	80.0	86.3	
100 4,6-Dinitro-2-methylphenol	198	8.298	8.298	0.000	85	468101	160.0	181.6	
102 N-Nitrosodiphenylamine	169	8.357	8.357	0.000	61	930374	80.0	86.2	
103 Azobenzene	77	8.392	8.392	0.000	100	1326025	80.0	79.6	
104 1,2-Diphenylhydrazine	77	8.392	8.392	0.000	100	1326025	80.9	80.5	
111 4-Bromophenyl phenyl ether	248	8.728	8.728	0.000	70	288713	80.0	84.6	
112 Hexachlorobenzene	284	8.810	8.810	0.000	92	296937	80.0	83.8	
116 Pentachlorophenol	266	9.010	9.010	0.000	89	378615	160.0	162.5	
119 Phenanthrene	178	9.233	9.233	0.000	98	1673042	80.0	82.8	
120 Anthracene	178	9.286	9.286	0.000	98	1733749	80.0	83.8	
122 Carbazole	167	9.445	9.445	0.000	95	1760082	80.0	83.4	
123 Di-n-butyl phthalate	149	9.739	9.739	0.000	100	2112316	80.0	85.1	
128 Fluoranthene	202	10.675	10.675	0.000	98	1791559	80.0	81.0	
129 Benzidine	184		10.466				ND	ND	
131 Pyrene	202	11.039	11.039	0.000	97	1855215	80.0	86.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
136 Famphur	218	12.051	12.051	0.000	97	594971	80.0	87.5	
137 Butyl benzyl phthalate	149	12.151	12.151	0.000	97	928375	80.0	91.3	
140 3,3'-Dichlorobenzidine	252	13.468	13.468	0.000	75	559891	80.0	87.8	
141 Benzo[a]anthracene	228	13.504	13.504	0.000	99	1533206	80.0	82.7	
142 Bis(2-ethylhexyl) phthalat	149	13.492	13.492	0.000	97	1195540	80.0	90.0	
143 Chrysene	228	13.598	13.598	0.000	98	1461997	80.0	84.4	
144 Di-n-octyl phthalate	149	15.362	15.362	0.000	99	2125005	80.0	89.6	
146 Benzo[b]fluoranthene	252	16.509	16.509	0.000	99	1376089	80.0	85.4	
147 Benzo[k]fluoranthene	252	16.592	16.592	0.000	99	1396697	80.0	84.4	
148 Benzo[a]pyrene	252	17.492	17.492	0.000	80	1347455	80.0	87.0	
151 Indeno[1,2,3-cd]pyrene	276	20.862	20.862	0.000	96	1096109	80.0	83.2	M
152 Dibenz(a,h)anthracene	278	20.939	20.939	0.000	94	1111518	80.0	86.8	
153 Benzo[g,h,i]perylene	276	21.633	21.633	0.000	95	1166748	80.0	83.7	

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

MS-HSLACCV080_00038 Amount Added: 200.00 Units: uL

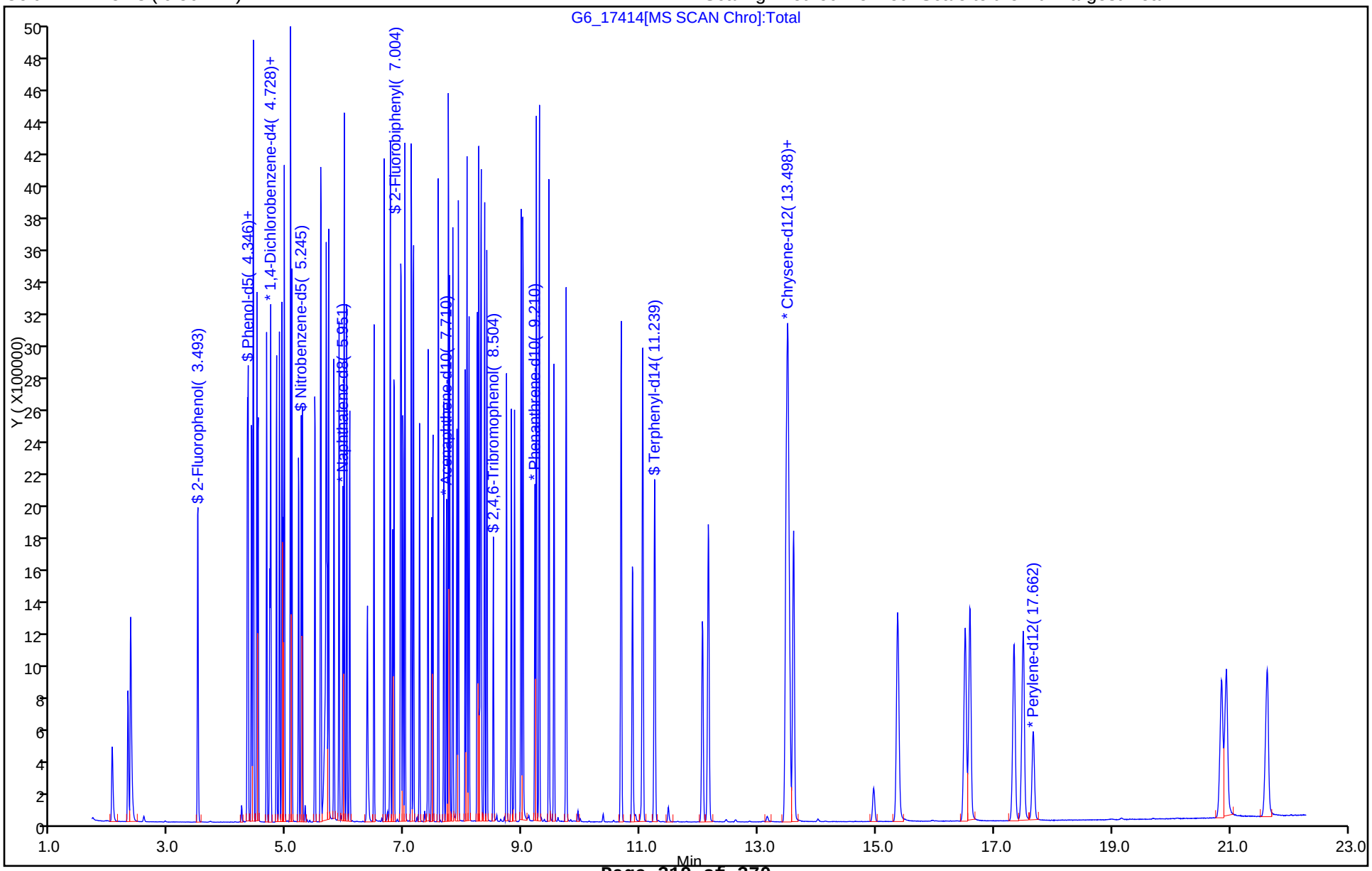
Report Date: 20-Apr-2015 07:09:12 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17414.D

Injection Date:	18-Apr-2015 15:20:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	CCV HSL			Worklist Smp#:	3
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	2
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



Manual Integration/User Assign Peak Report

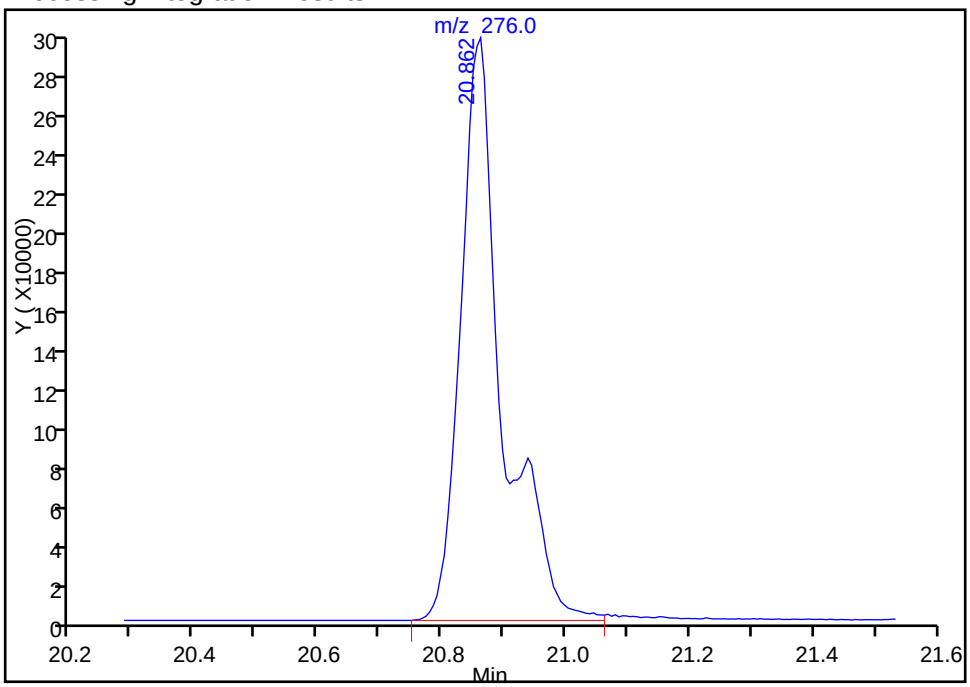
TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17414.D
Injection Date: 18-Apr-2015 15:20:30 Instrument ID: SMS_G6
Lims ID: CCV HSL
Client ID:
Operator ID: HOEFLERA ALS Bottle#: 2 Worklist Smp#: 3
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Method: SMS_G6_8270D Limit Group: MSSV - 8270D
Column: VF-5ms (0.50 mm) Detector MS SCAN

151 Indeno[1,2,3-cd]pyrene, CAS: 193-39-5

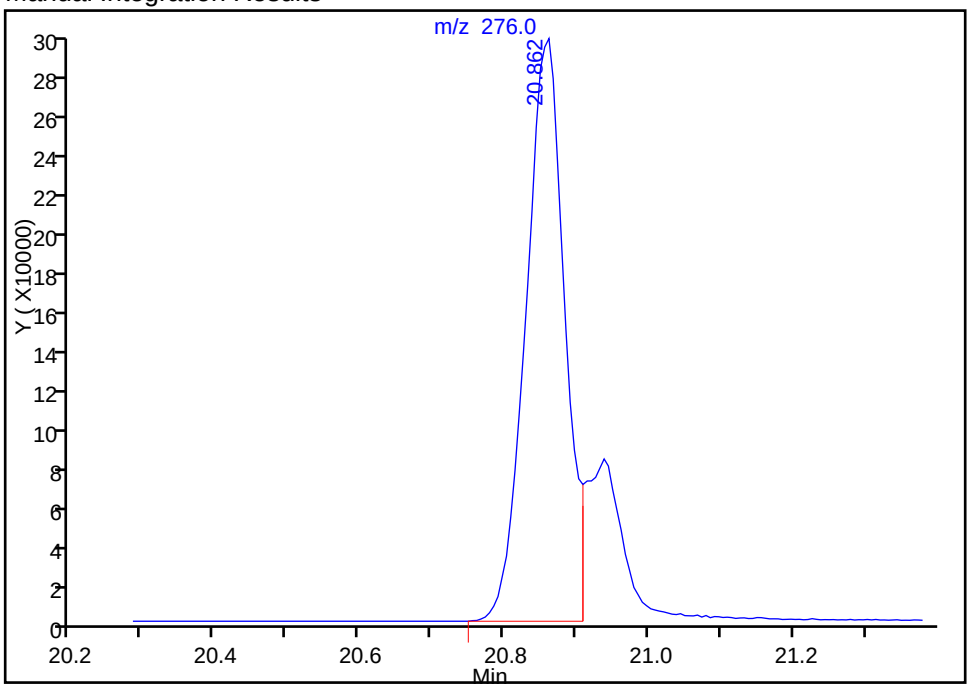
Processing Integration Results

RT: 20.86
Area: 1367322
Amount: 103.8016
Amount Units: ug/ml



Manual Integration Results

RT: 20.86
Area: 1096109
Amount: 83.212174
Amount Units: ug/ml



Reviewer: hoeflera, 18-Apr-2015 16:20:23
Audit Action: Split an Integrated Peak
Audit Reason: Shouldering

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16636.D
Lims ID: DFTPP
Client ID:
Sample Type: DFTPP
Inject. Date: 25-Feb-2015 11:44:30 ALS Bottle#: 1 Worklist Smp#: 2
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: DFTPP
Operator ID: KIEKELD Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 10-Apr-2015 12:09:37 Calib Date: 25-Feb-2015 14:59:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16644.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK023

First Level Reviewer: kiekeld Date: 10-Apr-2015 08:08:49

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
20 Pentachlorophenol_T	266	3.898	3.898	0.000	90	188400	NR	NR	7
35 Benzidine_T	184	5.051	5.051	0.000	100	1315879	NR	NR	7
156 DFTPP									
157 4,4'-DDE	246	5.169	5.169	0.000	58	654	NR	NR	7
158 4,4'-DDD	235	5.487	5.487	0.000	94	10699	NR	NR	7
159 4,4'-DDT	235	5.734	5.734	0.000	97	496436	NR	NR	7

QC Flag Legend

Processing Flags
NR - Missing Quant Standard
7 - Failed Limit of Detection

Reagents:

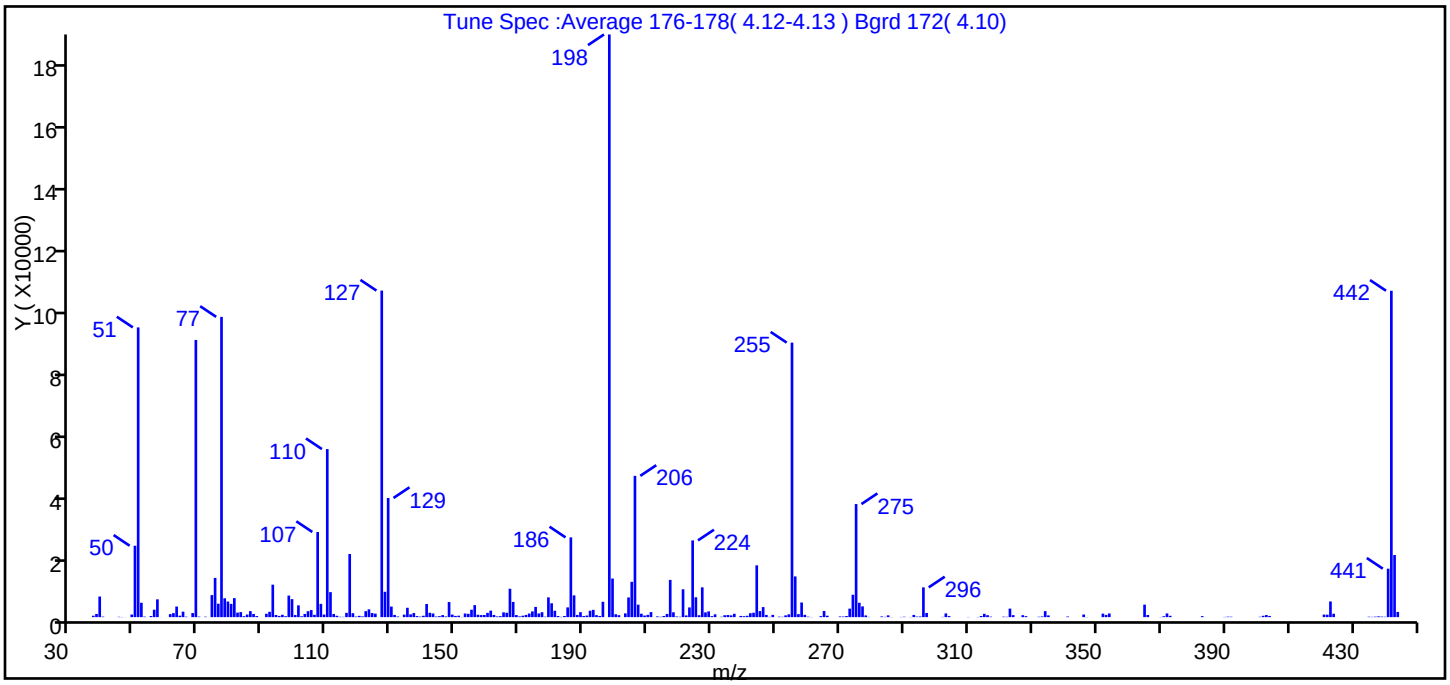
MS-DFTPP_00038 Amount Added: 200.00 Units: uL

MS Tune Report

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16636.D
Injection Date: 25-Feb-2015 11:44:30 Instrument ID: SMS_G6
Lims ID: DFTPP
Client ID:
Operator ID: KIEKELD ALS Bottle#: 1 Worklist Smp#: 2
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Method: SMS_G6_8270D Limit Group: MSSV - 8270D
Tune Method: DFTPP Method 8270

156 DFTPP



m/z	Ion Abundance Criteria	% Relative Abundance
198	Base peak, 100% relative abundance	100.0
51	30-60% of mass 198	49.7
68	<2% of mass 69	0.7 (1.5)
69	Present	47.6
70	<2% of mass 69	0.1 (0.2)
127	40-60% of mass 198	56.0
197	<1% of mass 198	0.0
199	5-9% of mass 198	6.6
275	10-30% of mass 198	19.4
365	>1% of mass 198	2.1
441	Present but less than mass 443	8.3 (78.0)
442	>40% of mass 198	56.0
443	17-23% of mass 442	10.7 (19.0)

Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16636.D\SMS_G6_8270D.rslt\spectra.d
Injection Date: 25-Feb-2015 11:44:30
Spectrum: Tune Spec :Average 176-178(4.12-4.13) Bgrd 172(4.10)
Base Peak: 198.00
Minimum % Base Peak: 0
Number of Points: 285

m/z	Y	m/z	Y	m/z	Y	m/z	Y
37.00	459	121.00	232	194.00	649	277.00	3468
38.00	1028	122.00	1930	195.00	379	278.00	551
39.00	6660	123.00	2519	196.00	4964	279.00	72
40.00	175	124.00	1326	198.00	188224	283.00	229
45.00	111	125.00	1129	199.00	12470	284.00	86
46.00	56	127.00	105488	200.00	938	285.00	577
47.00	18	128.00	8194	201.00	634	286.00	51
49.00	841	129.00	38496	203.00	1213	289.00	66
50.00	23072	130.00	3418	204.00	6363	290.00	132
51.00	93600	131.00	719	205.00	11405	292.00	60
52.00	4637	132.00	207	206.00	45640	293.00	716
53.00	168	133.00	8	207.00	4035	294.00	175
55.00	369	134.00	809	208.00	1101	295.00	204
56.00	2402	135.00	3026	209.00	516	296.00	9622
57.00	5751	136.00	912	210.00	745	297.00	1351
58.00	19	137.00	1332	211.00	1629	302.00	79
61.00	984	138.00	345	213.00	241	303.00	1194
62.00	1301	139.00	162	214.00	105	304.00	346
63.00	3429	140.00	450	215.00	382	310.00	61
64.00	473	141.00	4239	216.00	1029	313.00	58
65.00	1769	142.00	1410	217.00	12030	314.00	373
66.00	142	143.00	1176	218.00	1576	315.00	1075
68.00	1318	144.00	194	219.00	212	316.00	634
69.00	89544	145.00	322	220.00	128	317.00	162
70.00	167	146.00	661	221.00	9001	321.00	150
72.00	147	147.00	230	222.00	550	322.00	108
74.00	7124	148.00	4884	223.00	3140	323.00	2758
75.00	12670	149.00	828	224.00	24792	324.00	651
76.00	4339	150.00	402	225.00	6420	327.00	629
77.00	97000	151.00	465	226.00	621	328.00	325
78.00	6095	152.00	62	227.00	9628	332.00	128
79.00	5037	153.00	1178	228.00	1542	333.00	215
80.00	4279	154.00	1072	229.00	1866	334.00	1943

Report Date: 10-Apr-2015 12:09:37 Chrom Revision: 2.2 13-Mar-2015 11:20:44
Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16636.D\SMS_G6_8270D.rslt\spectra.d
Injection Date: 25-Feb-2015 11:44:30
Spectrum: Tune Spec :Average 176-178(4.12-4.13) Bgrd 172(4.10)
Base Peak: 198.00
Minimum % Base Peak: 0
Number of Points: 285

m/z	Y	m/z	Y	m/z	Y	m/z	Y
81.00	6145	155.00	2405	230.00	331	335.00	561
82.00	1463	156.00	3883	231.00	909	341.00	204
83.00	1631	157.00	773	233.00	124	346.00	858
84.00	302	158.00	659	234.00	617	347.00	53
85.00	841	159.00	698	235.00	689	352.00	1135
86.00	1946	160.00	1385	236.00	574	353.00	731
87.00	1021	161.00	2102	237.00	1086	354.00	1189
88.00	349	162.00	717	239.00	395	355.00	7
89.00	5	163.00	237	240.00	309	365.00	4039
91.00	1091	164.00	319	241.00	444	366.00	704
92.00	1723	165.00	1551	242.00	1272	370.00	57
93.00	10487	166.00	1415	243.00	1423	371.00	292
94.00	690	167.00	9174	244.00	16736	372.00	1201
95.00	398	168.00	4916	245.00	1957	373.00	434
96.00	758	169.00	673	246.00	3250	383.00	328
97.00	297	170.00	245	247.00	721	390.00	70
98.00	6968	171.00	465	248.00	88	391.00	164
99.00	5816	172.00	712	249.00	724	392.00	137
100.00	676	173.00	1157	251.00	163	401.00	116
101.00	3817	174.00	1856	252.00	105	402.00	438
102.00	279	175.00	3296	253.00	569	403.00	684
103.00	1041	176.00	1189	254.00	898	404.00	379
104.00	1921	177.00	1604	255.00	88688	421.00	838
105.00	2282	178.00	135	256.00	13156	422.00	784
106.00	758	179.00	6388	257.00	978	423.00	5066
107.00	27512	180.00	4482	258.00	4733	424.00	1095
108.00	4296	181.00	2051	259.00	747	435.00	136
109.00	843	182.00	372	260.00	153	436.00	96
110.00	54288	183.00	82	261.00	51	437.00	156
111.00	8075	184.00	376	264.00	330	438.00	252
112.00	1035	185.00	3153	265.00	2008	439.00	169
113.00	446	186.00	25768	266.00	463	440.00	121
114.00	128	187.00	7034	270.00	218	441.00	15657
115.00	102	188.00	729	271.00	188	442.00	105440

Report Date: 10-Apr-2015 12:09:37 Chrom Revision: 2.2 13-Mar-2015 11:20:44
Data File: \\Denchrom\ChromData\SMS_G6\20150410-33765.b\G6_16636.D\SMS_G6_8270D.rslt\spectra.d
Injection Date: 25-Feb-2015 11:44:30
Spectrum: Tune Spec :Average 176-178(4.12-4.13) Bgrd 172(4.10)
Base Peak: 198.00
Minimum % Base Peak: 0
Number of Points: 285

m/z	Y	m/z	Y	m/z	Y	m/z	Y
116.00	1390	189.00	1612	272.00	353	443.00	20072
117.00	20384	190.00	354	273.00	2747	444.00	1689
118.00	1268	191.00	522	274.00	7230		
119.00	185	192.00	2080	275.00	36528		
120.00	419	193.00	2355	276.00	4643		

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17413.D
Lims ID: DFTPP
Client ID:
Sample Type: DFTPP
Inject. Date: 18-Apr-2015 15:09:30 ALS Bottle#: 1 Worklist Smp#: 2
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: DFTPP
Operator ID: HOEFLERA Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:10 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: hoeflera Date: 18-Apr-2015 16:00:53

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
20 Pentachlorophenol_T	266	3.822	3.822	0.000	90	155066	NR	NR	7
35 Benzidine_T	184	4.952	4.952	0.000	100	1154158	NR	NR	7
156 DFTPP									
157 4,4'-DDE	246	5.063	5.063	0.000	71	443	NR	NR	7
158 4,4'-DDD	235	5.375	5.375	0.000	94	10074	NR	NR	7
159 4,4'-DDT	235	5.604	5.604	0.000	97	420659	NR	NR	7

QC Flag Legend

Processing Flags
NR - Missing Quant Standard
7 - Failed Limit of Detection

Reagents:

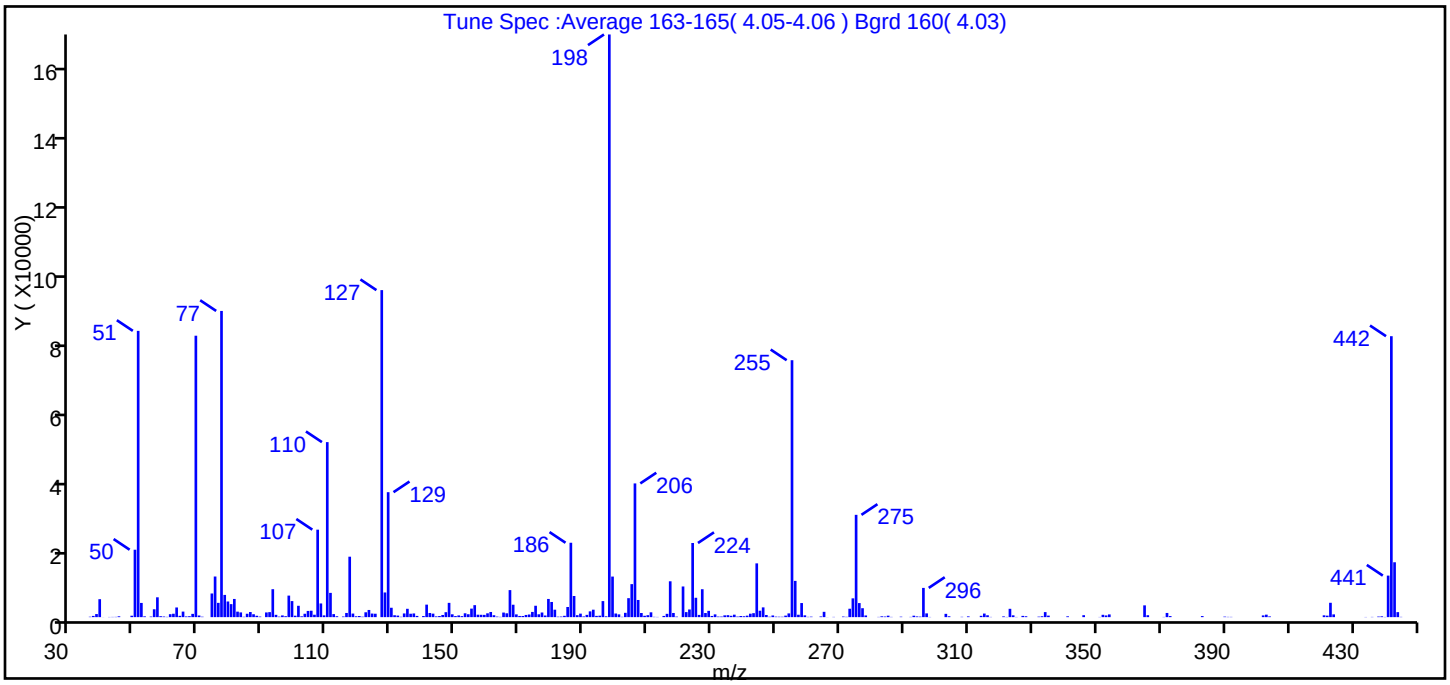
MS-DFTPP_00038 Amount Added: 200.00 Units: uL

MS Tune Report

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17413.D
Injection Date: 18-Apr-2015 15:09:30 Instrument ID: SMS_G6
Lims ID: DFTPP
Client ID:
Operator ID: HOEFLERA ALS Bottle#: 1 Worklist Smp#: 2
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Method: SMS_G6_8270D Limit Group: MSSV - 8270D
Tune Method: DFTPP Method 8270

156 DFTPP



m/z	Ion Abundance Criteria	% Relative Abundance
198	Base peak, 100% relative abundance	100.0
51	30-60% of mass 198	49.1
68	<2% of mass 69	0.6 (1.2)
69	Present	48.3
70	<2% of mass 69	0.3 (0.6)
127	40-60% of mass 198	56.1
197	<1% of mass 198	0.1
199	5-9% of mass 198	7.0
275	10-30% of mass 198	17.6
365	>1% of mass 198	2.0
441	Present but less than mass 443	7.1 (75.7)
442	>40% of mass 198	48.2
443	17-23% of mass 442	9.4 (19.5)

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17413.D\SMS_G6_8270D.rslt\spectra.d
Injection Date: 18-Apr-2015 15:09:30
Spectrum: Tune Spec :Average 163-165(4.05-4.06) Bgrd 160(4.03)
Base Peak: 198.00
Minimum % Base Peak: 0
Number of Points: 289

m/z	Y	m/z	Y	m/z	Y	m/z	Y
36.00	116	118.00	1027	193.00	2136	273.00	2412
37.00	353	119.00	243	194.00	364	274.00	5390
38.00	884	120.00	324	195.00	418	275.00	29328
39.00	5148	121.00	125	196.00	4599	276.00	4033
40.00	19	122.00	1392	197.00	233	277.00	2563
42.00	54	123.00	2058	198.00	167040	278.00	481
43.00	66	124.00	1028	199.00	11632	282.00	66
44.00	75	125.00	979	200.00	1041	283.00	237
45.00	240	127.00	93752	201.00	794	284.00	171
49.00	406	128.00	7074	202.00	90	285.00	420
50.00	19328	129.00	35824	203.00	1238	286.00	68
51.00	82072	130.00	2693	204.00	5440	289.00	133
52.00	4075	131.00	545	205.00	9465	292.00	70
53.00	224	132.00	444	206.00	38336	293.00	414
55.00	168	133.00	99	207.00	4920	294.00	204
56.00	2257	134.00	1058	208.00	1183	295.00	140
57.00	5688	135.00	2392	209.00	394	296.00	8383
58.00	269	136.00	979	210.00	564	297.00	1076
59.00	176	137.00	1058	211.00	1384	298.00	54
60.00	80	138.00	353	212.00	86	303.00	945
61.00	848	140.00	313	213.00	61	304.00	252
62.00	976	141.00	3585	215.00	346	308.00	91
63.00	2772	142.00	1190	216.00	949	310.00	214
64.00	376	143.00	1000	217.00	10279	314.00	380
65.00	1592	144.00	212	218.00	1192	315.00	1026
66.00	195	145.00	318	219.00	149	316.00	536
67.00	268	146.00	620	221.00	8791	317.00	64
68.00	940	147.00	1422	222.00	1443	321.00	212
69.00	80680	148.00	4096	223.00	2198	322.00	56
70.00	468	149.00	815	224.00	21256	323.00	2389
71.00	124	150.00	305	225.00	5553	324.00	516
74.00	6782	151.00	489	226.00	639	325.00	56
75.00	11648	152.00	291	227.00	8004	327.00	360

Report Date: 20-Apr-2015 07:09:10 Chrom Revision: 2.2 09-Apr-2015 10:05:40
Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17413.D\SMS_G6_8270D.rslt\spectra.d
Injection Date: 18-Apr-2015 15:09:30
Spectrum: Tune Spec :Average 163-165(4.05-4.06) Bgrd 160(4.03)
Base Peak: 198.00
Minimum % Base Peak: 0
Number of Points: 289

m/z	Y	m/z	Y	m/z	Y	m/z	Y
76.00	4093	153.00	1107	228.00	1182	328.00	255
77.00	87784	154.00	827	229.00	1805	332.00	143
78.00	6373	155.00	2446	230.00	342	333.00	248
79.00	4487	156.00	3435	231.00	793	334.00	1445
80.00	3730	157.00	699	232.00	163	335.00	436
81.00	5215	158.00	674	233.00	166	341.00	236
82.00	1575	159.00	653	234.00	514	346.00	556
83.00	1356	160.00	1098	235.00	561	351.00	56
84.00	105	161.00	1481	236.00	388	352.00	670
85.00	973	162.00	551	237.00	706	353.00	489
86.00	1466	163.00	156	238.00	161	354.00	780
87.00	832	164.00	75	239.00	327	365.00	3378
88.00	397	165.00	1313	240.00	254	366.00	551
89.00	160	166.00	1103	241.00	425	371.00	58
91.00	1348	167.00	7751	242.00	986	372.00	1200
92.00	1437	168.00	3561	243.00	1137	373.00	307
93.00	8011	169.00	790	244.00	15415	383.00	307
94.00	661	170.00	339	245.00	1842	390.00	145
95.00	146	171.00	332	246.00	2800	391.00	77
96.00	490	172.00	630	247.00	617	392.00	72
97.00	266	173.00	757	248.00	117	402.00	496
98.00	6176	174.00	1494	249.00	481	403.00	711
99.00	4602	175.00	3265	250.00	145	404.00	215
100.00	411	176.00	816	251.00	121	421.00	521
101.00	3256	177.00	1330	252.00	111	422.00	428
102.00	234	178.00	467	253.00	434	423.00	4117
103.00	969	179.00	5213	254.00	1073	424.00	792
104.00	1780	180.00	4348	255.00	73656	434.00	50
105.00	1844	181.00	2137	256.00	10393	436.00	67
106.00	689	182.00	136	257.00	687	438.00	168
107.00	25080	183.00	168	258.00	4041	439.00	232
108.00	3924	184.00	381	259.00	518	440.00	57
109.00	505	185.00	2912	260.00	82	441.00	11908
110.00	50176	186.00	21328	261.00	139	442.00	80544

Report Date: 20-Apr-2015 07:09:10 Chrom Revision: 2.2 09-Apr-2015 10:05:40
Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17413.D\SMS_G6_8270D.rslt\spectra.d
Injection Date: 18-Apr-2015 15:09:30
Spectrum: Tune Spec :Average 163-165(4.05-4.06) Bgrd 160(4.03)
Base Peak: 198.00
Minimum % Base Peak: 0
Number of Points: 289

m/z	Y	m/z	Y	m/z	Y	m/z	Y
111.00	6956	187.00	6061	264.00	196	443.00	15732
112.00	872	188.00	563	265.00	1526	444.00	1434
113.00	257	189.00	1003	266.00	24	445.00	77
115.00	180	190.00	193	268.00	67		
116.00	1198	191.00	652	271.00	171		
117.00	17328	192.00	1657	272.00	91		

FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Denver | Job No.: 280-67711-2

SDG No.:

Client Sample ID: | Lab Sample ID: MB 280-272870/1-A

Matrix: Water | Lab File ID: G6_17420.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3520C | Date Extracted: 04/15/2015 18:50

Sample wt/vol: 1000(mL) | Date Analyzed: 04/18/2015 17:49

Con. Extract Vol.: 1(mL) | Dilution Factor: 1

Injection Volume: 0.5(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 273380 | Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
105-60-2	Caprolactam	2.5	U	5.0	2.5

CAS NO.	SURROGATE	%REC	Q	LIMITS
118-79-6	2,4,6-Tribromophenol (Surr)	65		48-135
321-60-8	2-Fluorobiphenyl	62		48-135
367-12-4	2-Fluorophenol (Surr)	64		41-135
4165-60-0	Nitrobenzene-d5 (Surr)	65		42-135
4165-62-2	Phenol-d5 (Surr)	64		46-135
1718-51-0	Terphenyl-d14 (Surr)	64		20-135

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17420.D
Lims ID: MB 280-272870/1-A
Client ID:
Sample Type: MB
Inject. Date: 18-Apr-2015 17:49:30 ALS Bottle#: 8 Worklist Smp#: 27
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: MB280-273368_1-A
Operator ID: HOEFLERA Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:13 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: Deconvolution ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: kiekeld Date: 20-Apr-2015 07:04:31

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	96	217388	40.0	40.0	
* 2 Naphthalene-d8	136	5.945	5.951	-0.006	100	813022	40.0	40.0	
* 3 Acenaphthene-d10	164	7.704	7.704	0.000	92	452207	40.0	40.0	
* 4 Phenanthrene-d10	188	9.204	9.210	-0.006	97	769713	40.0	40.0	
* 5 Chrysene-d12	240	13.510	13.516	-0.006	97	727965	40.0	40.0	
* 6 Perylene-d12	264	17.645	17.651	-0.006	95	630046	40.0	40.0	
$ 7 2-Fluorophenol	112	3.487	3.493	-0.006	93	504071	100.0	63.8	
$ 8 Phenol-d5	99	4.328	4.334	-0.006	99	638469	100.0	64.4	
$ 9 Nitrobenzene-d5	82	5.240	5.245	-0.005	90	537720	100.0	64.5	
$ 10 2-Fluorobiphenyl	172	6.998	7.004	-0.006	100	953384	100.0	62.4	
$ 11 2,4,6-Tribromophenol	330	8.498	8.504	-0.006	88	109771	100.0	64.5	
$ 12 Terphenyl-d14	244	11.233	11.239	-0.006	98	924818	100.0	63.9	
13 1,4-Dioxane	88		2.040					ND	
14 N-Nitrosodimethylamine	74		2.305					ND	
15 Pyridine	79		2.352					ND	
16 2-Picoline	93		2.987					ND	
17 N-Nitrosomethylethylamine	88		3.081					ND	
18 Methyl methanesulfonate	80		3.351					ND	
19 N-Nitrosodiethylamine	102		3.704					ND	
20 Pentachlorophenol_T	266		3.822					ND	
21 Ethyl methanesulfonate	79		3.969					ND	
22 Pentachloroethane	117		4.446					ND	
23 Phenol	94		4.346					ND	
24 Aniline	93		4.404					ND	
25 Bis(2-chloroethyl)ether	93		4.434					ND	
26 2-Chlorophenol	128		4.516					ND	
27 1,3-Dichlorobenzene	146		4.657					ND	
28 1,4-Dichlorobenzene	146		4.728					ND	
29 Benzyl alcohol	108		4.828					ND	
30 1,2-Dichlorobenzene	146		4.875					ND	
31 2-Methylphenol	108		4.916					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
32 2,2'-oxybis[1-chloropropan	45		4.934					ND	
33 N-Nitrosopyrrolidine	100		5.075					ND	
34 N-Nitrosomorpholine	116		5.104					ND	
35 Benzidine_T	184		4.952					ND	
36 2-Toluidine	106		5.122					ND	
37 Benzaldehyde	106		5.015					ND	
38 3 & 4 Methylphenol	108		5.063					ND	
39 3-Methylphenol	108		5.063					ND	
40 4-Methylphenol	108		5.063					ND	
41 N-Nitrosodi-n-propylamine	70		5.069					ND	
42 Acetophenone	105		5.087					ND	
43 Hexachloroethane	117		5.198					ND	
44 Nitrobenzene	77		5.263					ND	
45 N-Nitrosopiperidine	114		5.410					ND	
46 Isophorone	82		5.475					ND	
47 o,o',o''-Triethylphosphoro	198		5.622					ND	
48 2-Nitrophenol	139		5.569					ND	
49 2,4-Dimethylphenol	107		5.581					ND	
50 Bis(2-chloroethoxy)methane	93		5.669					ND	
51 alpha,alpha-Dimethyl phene	58		5.792					ND	
52 Benzoic acid	105		5.687					ND	
53 2,4-Dichlorophenol	162		5.798					ND	
54 1,2,4-Trichlorobenzene	180		5.887					ND	
55 2,6-Dichlorophenol	162		6.022					ND	
56 Hexachloropropene	213		6.045					ND	
57 Naphthalene	128		5.975					ND	
58 4-Chloroaniline	127		6.016					ND	
59 Hexachlorobutadiene	225		6.069					ND	
60 N-Nitrosodi-n-butylamine	84		6.310					ND	
61 p-Phenylene diamine	108		6.363					ND	
62 Caprolactam	55		6.369					ND	
63 Safrole, Total	162		6.540					ND	
64 4-Chloro-3-methylphenol	107		6.481					ND	
65 2-Methylnaphthalene	142		6.651					ND	
66 Isosafrole Peak 1	162		6.834					ND	
67 1-Methylnaphthalene	142		6.757					ND	
68 Hexachlorocyclopentadiene	237		6.798					ND	
69 1,2,4,5-Tetrachlorobenzene	216		6.822					ND	
70 2,4,6-Trichlorophenol	196		6.934					ND	
71 Isosafrole Peak 2	104		7.063					ND	
72 2,4,5-Trichlorophenol	196		6.969					ND	
73 1-Chloronaphthalene	162		7.169					ND	
74 1,1'-Biphenyl	154		7.110					ND	
75 2-Chloronaphthalene	162		7.151					ND	
76 1,4-Naphthoquinone	158		7.328					ND	
77 2-Nitroaniline	65		7.251					ND	
78 1,4-Dinitrobenzene	168		7.381					ND	
79 Dimethyl phthalate	163		7.398					ND	
80 1,3-Dinitrobenzene	168		7.463					ND	
81 2,6-Dinitrotoluene	165		7.481					ND	
82 Acenaphthylene	152		7.569					ND	
83 3-Nitroaniline	138		7.663					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
84 Acenaphthene	153		7.739					ND	
85 Pentachlorobenzene	250		7.863					ND	
86 2,4-Dinitrophenol	184		7.763					ND	
87 4-Nitrophenol	109		7.816					ND	
88 1-Naphthylamine	143		7.998					ND	
89 2,4-Dinitrotoluene	165		7.887					ND	
90 Dibenzofuran	168		7.910					ND	
91 2-Naphthylamine	143		8.075					ND	
92 2,3,4,6-Tetrachlorophenol	232		8.028					ND	
93 Thionazin	97		8.169					ND	
94 Diethyl phthalate	149		8.092					ND	
95 N-Nitro-o-toluidine	152		8.263					ND	
96 4-Chlorophenyl phenyl ethe	204		8.234					ND	
97 Diphenylamine	169		8.351					ND	
98 Fluorene	166		8.257					ND	
99 4-Nitroaniline	138		8.286					ND	
100 4,6-Dinitro-2-methylphenol	198		8.298					ND	
101 Sulfotepp	97		8.451					ND	
102 N-Nitrosodiphenylamine	169		8.357					ND	
103 Azobenzene	77		8.392					ND	
104 1,2-Diphenylhydrazine	77		8.392					ND	
105 Diallate Peak 1	86		8.610					ND	
106 1,3,5-Trinitrobenzene	213		8.622					ND	
107 Phorate	121		8.628					ND	
108 Phenacetin	108		8.645					ND	
109 Diallate Peak 2	86		8.704					ND	
110 Dimethoate	87		8.822					ND	
111 4-Bromophenyl phenyl ether	248		8.728					ND	
112 Hexachlorobenzene	284		8.810					ND	
113 4-Aminobiphenyl	169		8.998					ND	
114 Pentachloronitrobenzene	237		9.010					ND	
115 Pronamide	173		9.016					ND	
116 Pentachlorophenol	266		9.010					ND	
117 Disulfoton	88		9.145					ND	
118 Dinoseb	211		9.157					ND	
119 Phenanthrene	178		9.233					ND	
120 Anthracene	178		9.286					ND	
121 Methyl parathion	109		9.551					ND	
122 Carbazole	167		9.445					ND	
123 Di-n-butyl phthalate	149		9.739					ND	
124 Ethyl Parathion	109		9.980					ND	
125 4-Nitroquinoline-1-oxide	190		10.104					ND	
126 Methapyrilene	97		10.128					ND	
127 Isodrin	193		10.433					ND	
128 Fluoranthene	202		10.675					ND	
129 Benzidine	184		10.466					ND	
130 Aramite Peak 1	185		11.139					ND	
131 Pyrene	202		11.039					ND	
132 Aramite Peak 2	185		11.274					ND	
133 p-Dimethylamino azobenzene	120		11.492					ND	
134 Chlorobenzilate	251		11.545					ND	
135 3,3'-Dimethylbenzidine	212		12.180					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
136 Famphur	218		12.045					ND	
137 Butyl benzyl phthalate	149		12.151					ND	
138 2-Acetylaminofluorene	181		12.745					ND	
139 4,4'-Methylene bis(2-chlor	231		13.463					ND	
140 3,3'-Dichlorobenzidine	252		13.468					ND	
141 Benzo[a]anthracene	228		13.504					ND	
142 Bis(2-ethylhexyl) phthalat	149		13.492					ND	
143 Chrysene	228		13.598					ND	
144 Di-n-octyl phthalate	149		15.362					ND	
145 7,12-Dimethylbenz(a)anthra	256		16.480					ND	
146 Benzo[b]fluoranthene	252		16.509					ND	
147 Benzo[k]fluoranthene	252		16.592					ND	
148 Benzo[a]pyrene	252		17.492					ND	
149 3-Methylcholanthrene	268		18.550					ND	
150 Dibenz[a,j]acridine	279		20.350					ND	
151 Indeno[1,2,3-cd]pyrene	276		20.862					ND	
152 Dibenz(a,h)anthracene	278		20.939					ND	
153 Benzo[g,h,i]perylene	276		21.633					ND	
S 160 Aramite, Total	185		15.047					ND	
S 161 Isosafrole	162		15.047					ND	
S 162 Diallate	86		15.047					ND	
154 Total Cresols	1		0.000					ND	
155 Tetraethyl Pyrophosphate (	1		0.000					ND	
157 4,4'-DDE	246		5.063					ND	
158 4,4'-DDD	235		5.375					ND	
159 4,4'-DDT	235		5.604					ND	
S 163 Methyl Phenols,Total	1		0.000					ND	

Reagents:

MS-IS_00007 Amount Added: 20.00 Units: uL Run Reagent

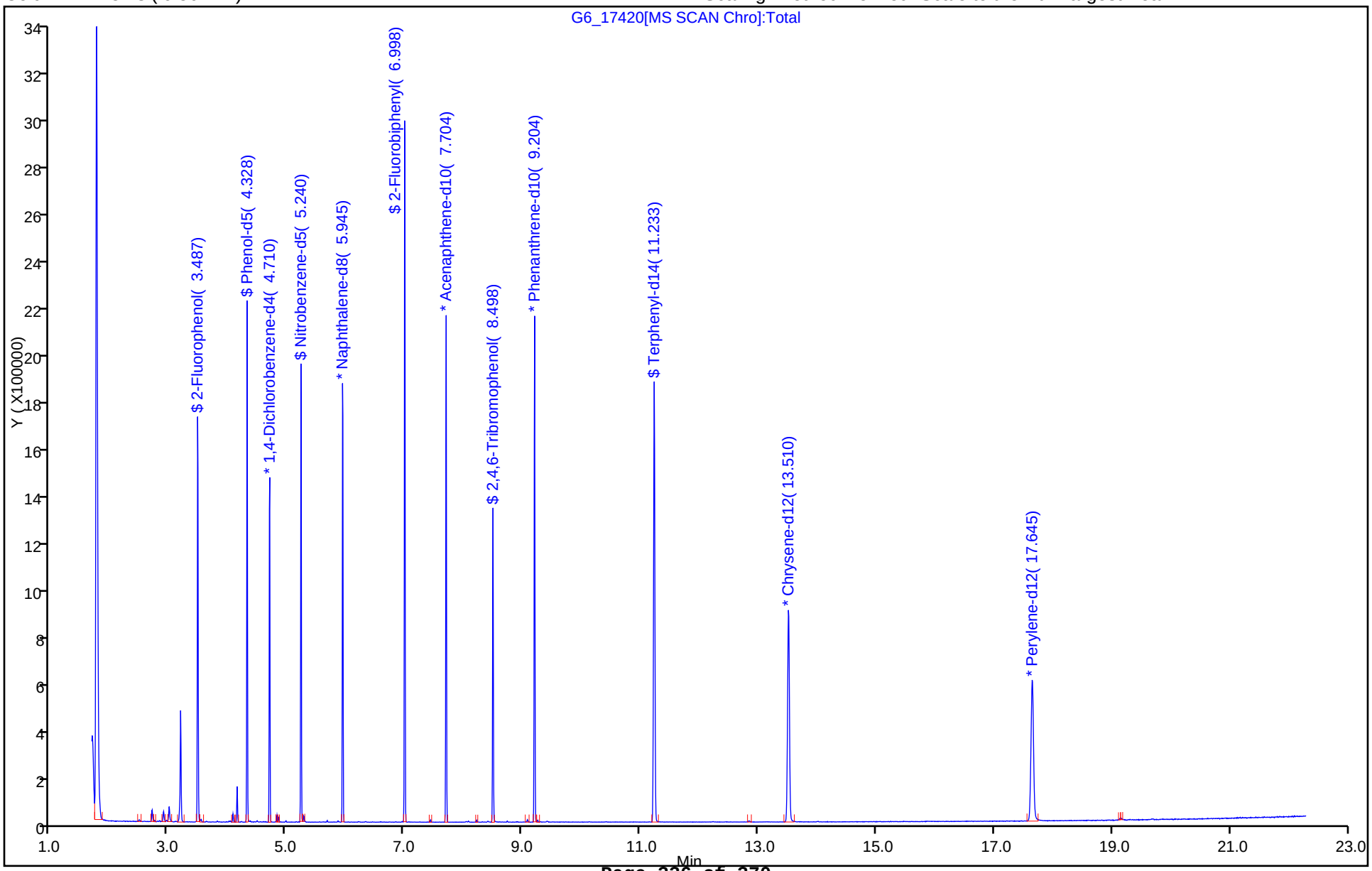
Report Date: 20-Apr-2015 07:09:14 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17420.D

Injection Date:	18-Apr-2015 17:49:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	MB 280-272870/1-A			Worklist Smp#:	27
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	8
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Denver | Job No.: 280-67711-2

SDG No.:

Client Sample ID: | Lab Sample ID: LCS 280-272870/2-A

Matrix: Water | Lab File ID: G6_17421.D

Analysis Method: 8270D | Date Collected:

Extract. Method: 3520C | Date Extracted: 04/15/2015 18:50

Sample wt/vol: 1000(mL) | Date Analyzed: 04/18/2015 18:16

Con. Extract Vol.: 1(mL) | Dilution Factor: 1

Injection Volume: 0.5(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 273380 | Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
105-60-2	Caprolactam	52.3		5.0	2.5

CAS NO.	SURROGATE	%REC	Q	LIMITS
118-79-6	2,4,6-Tribromophenol (Surr)	68		48-135
321-60-8	2-Fluorobiphenyl	61		48-135
367-12-4	2-Fluorophenol (Surr)	60		41-135
4165-60-0	Nitrobenzene-d5 (Surr)	61		42-135
4165-62-2	Phenol-d5 (Surr)	60		46-135
1718-51-0	Terphenyl-d14 (Surr)	62		20-135

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17421.D
Lims ID: LCS 280-272870/2-A
Client ID:
Sample Type: LCS
Inject. Date: 18-Apr-2015 18:16:30 ALS Bottle#: 9 Worklist Smp#: 28
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: LCS280-273368_2-A
Operator ID: HOEFLERA Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:13 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: RT Order ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: kiekeld Date: 20-Apr-2015 07:05:12

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	95	220573	40.0	40.0	
* 2 Naphthalene-d8	136	5.951	5.951	0.000	99	836102	40.0	40.0	
* 3 Acenaphthene-d10	164	7.704	7.704	0.000	92	460509	40.0	40.0	
* 4 Phenanthrene-d10	188	9.204	9.210	-0.006	97	780849	40.0	40.0	
* 5 Chrysene-d12	240	13.515	13.516	-0.001	96	716077	40.0	40.0	
* 6 Perylene-d12	264	17.650	17.651	-0.001	95	613809	40.0	40.0	
$ 7 2-Fluorophenol	112	3.487	3.493	-0.006	93	477204	100.0	59.5	
$ 8 Phenol-d5	99	4.328	4.334	-0.006	99	607080	100.0	60.4	
$ 9 Nitrobenzene-d5	82	5.245	5.245	0.000	89	520529	100.0	60.7	
$ 10 2-Fluorobiphenyl	172	7.004	7.004	0.000	99	952495	100.0	61.2	
$ 11 2,4,6-Tribromophenol	330	8.504	8.504	0.000	93	117984	100.0	68.1	
$ 12 Terphenyl-d14	244	11.233	11.239	-0.006	98	882084	100.0	62.0	
13 1,4-Dioxane	88	2.034	2.040	-0.006	97	152404	80.0	45.0	
14 N-Nitrosodimethylamine	74	2.304	2.305	0.000	92	265571	80.0	51.7	
15 Pyridine	79	2.346	2.352	-0.006	92	371789	80.0	42.4	
23 Phenol	94	4.340	4.346	-0.006	99	499291	80.0	49.5	
24 Aniline	93	4.398	4.404	-0.006	98	503575	80.0	40.9	
25 Bis(2-chloroethyl)ether	93	4.434	4.434	0.000	95	394820	80.0	49.9	
26 2-Chlorophenol	128	4.510	4.516	-0.006	97	414835	80.0	50.4	
27 1,3-Dichlorobenzene	146	4.657	4.657	0.000	97	287607	80.0	34.4	
28 1,4-Dichlorobenzene	146	4.728	4.728	0.000	94	302049	80.0	35.8	
29 Benzyl alcohol	108	4.822	4.828	-0.006	92	273532	80.0	52.8	
30 1,2-Dichlorobenzene	146	4.875	4.875	0.000	95	300275	80.0	37.0	
31 2-Methylphenol	108	4.916	4.916	0.000	94	377968	80.0	50.6	
32 2,2'-oxybis[1-chloropropan	45	4.934	4.934	0.000	94	522666	80.0	44.9	
37 Benzaldehyde	106	5.081	5.015	0.066	70	44007	NC	NC	
38 3 & 4 Methylphenol	108	5.057	5.063	-0.006	93	399732	80.0	51.6	
39 3-Methylphenol	108	5.057	5.063	-0.006	93	399732	80.0	51.6	
40 4-Methylphenol	108	5.057	5.063	-0.006	93	399732	80.0	51.6	
41 N-Nitrosodi-n-propylamine	70	5.063	5.069	-0.006	85	276909	80.0	49.2	
42 Acetophenone	105	5.081	5.087	-0.006	97	552998	80.0	50.8	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
43 Hexachloroethane	117	5.198	5.198	0.000	92	114452	80.0	31.6	
44 Nitrobenzene	77	5.263	5.263	0.000	88	412244	80.0	49.9	
46 Isophorone	82	5.475	5.475	0.000	99	757833	80.0	54.0	
48 2-Nitrophenol	139	5.569	5.569	0.000	93	218443	80.0	53.1	
49 2,4-Dimethylphenol	107	5.575	5.581	-0.006	92	356522	80.0	47.3	
50 Bis(2-chloroethoxy)methane	93	5.669	5.669	0.000	99	480367	80.0	52.1	
52 Benzoic acid	105	5.651	5.687	-0.036	88	299733	80.0	50.6	
53 2,4-Dichlorophenol	162	5.798	5.798	0.000	95	331628	80.0	53.7	
54 1,2,4-Trichlorobenzene	180	5.881	5.887	-0.006	93	245435	80.0	39.1	
55 2,6-Dichlorophenol	162	6.022	6.022	0.000	94	318072	80.0	54.7	
57 Naphthalene	128	5.969	5.975	-0.006	97	981081	80.0	44.1	
58 4-Chloroaniline	127	6.016	6.016	0.000	82	412003	80.0	41.0	
59 Hexachlorobutadiene	225	6.063	6.069	-0.006	96	116037	80.0	33.9	
62 Caprolactam	55	6.351	6.369	-0.018	77	206963	80.0	52.3	
64 4-Chloro-3-methylphenol	107	6.475	6.481	-0.006	96	347066	80.0	52.5	
65 2-Methylnaphthalene	142	6.651	6.651	0.000	91	676260	80.0	47.3	
67 1-Methylnaphthalene	142	6.751	6.757	-0.006	91	595310	80.0	44.3	
68 Hexachlorocyclopentadiene	237	6.792	6.798	-0.006	93	37627	80.0	10.4	
69 1,2,4,5-Tetrachlorobenzene	216	6.816	6.822	-0.006	97	282238	80.0	47.4	
70 2,4,6-Trichlorophenol	196	6.928	6.934	-0.006	93	221926	80.0	53.7	
72 2,4,5-Trichlorophenol	196	6.969	6.969	0.000	94	241904	80.0	54.9	
74 1,1'-Biphenyl	154	7.110	7.110	0.000	95	850377	80.0	49.1	
75 2-Chloronaphthalene	162	7.145	7.151	-0.006	97	636545	80.0	48.7	
77 2-Nitroaniline	65	7.245	7.251	-0.006	82	249401	80.0	52.0	
79 Dimethyl phthalate	163	7.392	7.398	-0.006	98	793195	80.0	51.7	
80 1,3-Dinitrobenzene	168	7.457	7.463	-0.006	87	154953	80.0	57.0	
81 2,6-Dinitrotoluene	165	7.475	7.481	-0.006	94	195566	80.0	55.2	
82 Acenaphthylene	152	7.569	7.569	0.000	94	1106711	80.0	52.2	
83 3-Nitroaniline	138	7.657	7.663	-0.006	94	198128	80.0	43.2	
84 Acenaphthene	153	7.739	7.739	0.000	94	687561	80.0	50.2	
86 2,4-Dinitrophenol	184	7.757	7.763	-0.006	81	267133	160.0	105.3	
87 4-Nitrophenol	109	7.810	7.816	-0.006	95	259488	160.0	105.0	
89 2,4-Dinitrotoluene	165	7.881	7.887	-0.006	90	266797	80.0	57.4	
90 Dibenzofuran	168	7.910	7.910	0.000	90	1001525	80.0	51.9	
92 2,3,4,6-Tetrachlorophenol	232	8.028	8.028	0.000	74	192418	80.0	53.9	
94 Diethyl phthalate	149	8.086	8.092	-0.006	97	851413	80.0	54.6	
96 4-Chlorophenyl phenyl ethe	204	8.228	8.234	-0.006	95	354210	80.0	50.7	
97 Diphenylamine	169	8.351	8.351	0.000	95	1182526	137.1	89.8	
98 Fluorene	166	8.251	8.257	-0.006	85	826096	80.0	52.6	
99 4-Nitroaniline	138	8.280	8.286	-0.006	83	244786	80.0	54.6	
100 4,6-Dinitro-2-methylphenol	198	8.292	8.298	-0.006	79	317037	160.0	114.8	
102 N-Nitrosodiphenylamine	169	8.351	8.357	-0.006	62	1182526	160.0	102.2	
103 Azobenzene	77	8.392	8.392	0.000	100	843818	80.0	48.9	
104 1,2-Diphenylhydrazine	77	8.392	8.392	0.000	100	843818	80.9	49.5	
111 4-Bromophenyl phenyl ether	248	8.722	8.728	-0.006	70	190813	80.0	52.2	
112 Hexachlorobenzene	284	8.804	8.810	-0.006	87	202556	80.0	53.3	
116 Pentachlorophenol	266	9.004	9.010	-0.006	87	241454	160.0	96.7	
119 Phenanthrene	178	9.227	9.233	-0.006	98	1152717	80.0	53.2	
120 Anthracene	178	9.280	9.286	-0.006	98	1150419	80.0	51.8	
122 Carbazole	167	9.445	9.445	0.000	82	1210183	80.0	53.5	
123 Di-n-butyl phthalate	149	9.733	9.739	-0.006	100	1408848	80.0	52.9	
128 Fluoranthene	202	10.663	10.675	-0.012	98	1230444	80.0	51.9	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
129 Benzidine	184		10.466				ND	ND	
131 Pyrene	202	11.027	11.039	-0.012	97	1288589	80.0	53.9	
137 Butyl benzyl phthalate	149	12.145	12.151	-0.006	96	626461	80.0	55.5	
140 3,3'-Dichlorobenzidine	252	13.451	13.468	-0.017	73	287004	80.0	40.6	
141 Benzo[a]anthracene	228	13.492	13.504	-0.012	99	1074591	80.0	52.2	
142 Bis(2-ethylhexyl) phthalat	149	13.480	13.492	-0.012	93	834928	80.0	56.6	
143 Chrysene	228	13.586	13.598	-0.012	96	1020493	80.0	53.1	
144 Di-n-octyl phthalate	149	15.351	15.362	-0.011	99	1428540	80.0	54.9	
146 Benzo[b]fluoranthene	252	16.492	16.509	-0.017	94	940541	80.0	52.5	
147 Benzo[k]fluoranthene	252	16.574	16.592	-0.018	98	967526	80.0	52.5	
148 Benzo[a]pyrene	252	17.474	17.492	-0.018	79	910137	80.0	52.8	
151 Indeno[1,2,3-cd]pyrene	276	20.838	20.862	-0.024	96	765436	80.0	52.4	
152 Dibenz(a,h)anthracene	278	20.921	20.939	-0.018	90	759647	80.0	53.3	
153 Benzo[g,h,i]perylene	276	21.609	21.633	-0.024	90	817592	80.0	52.7	

QC Flag Legend

Processing Flags

NC - Not Calibrated

ND - Not Detected or Marked ND

Reagents:

MS-IS_00007 Amount Added: 20.00 Units: uL Run Reagent

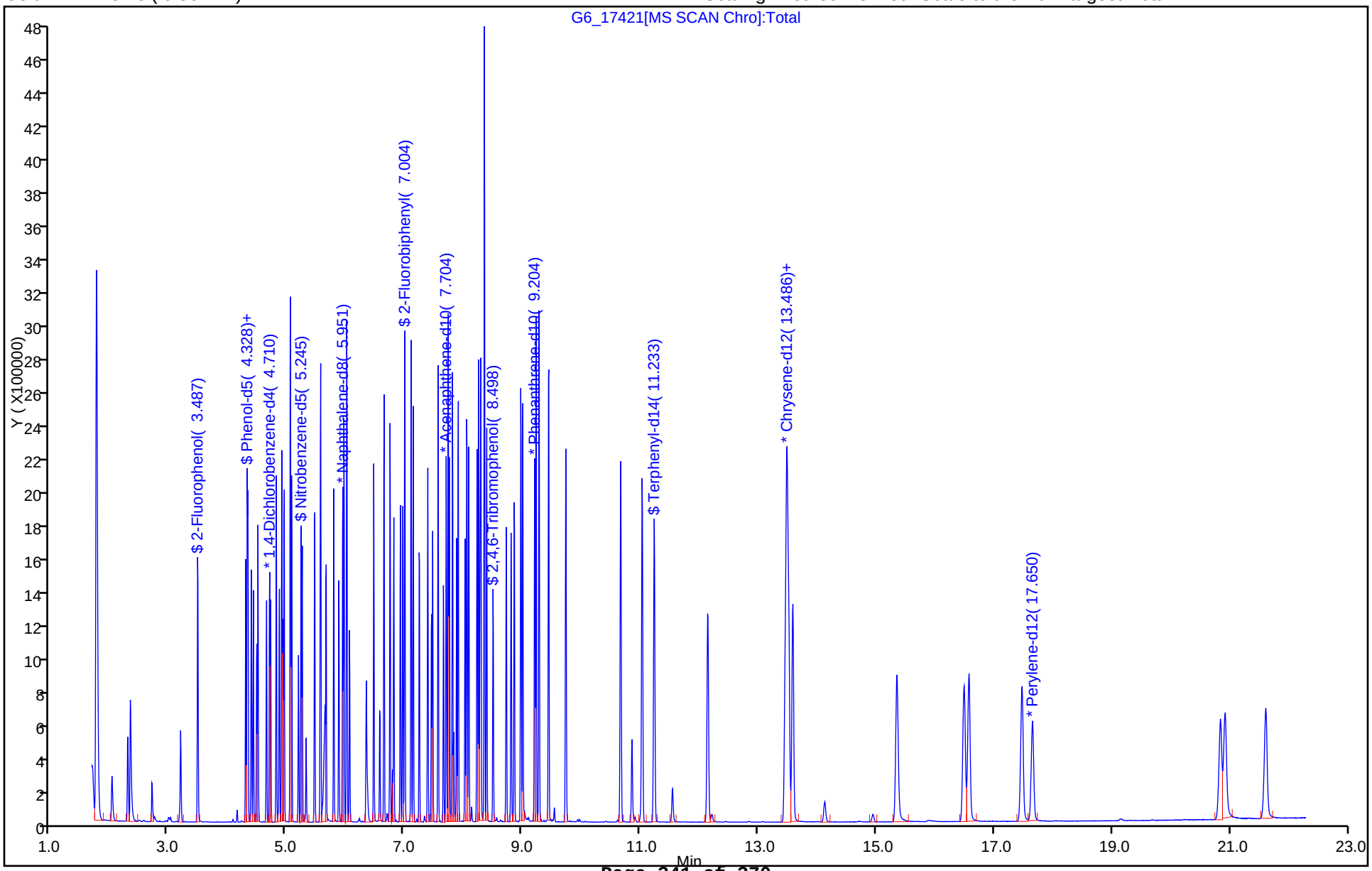
Report Date: 20-Apr-2015 07:09:15 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17421.D

Injection Date:	18-Apr-2015 18:16:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	LCS 280-272870/2-A			Worklist Smp#:	28
Client ID:					
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	9
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Denver | Job No.: 280-67711-2

SDG No.:

Client Sample ID: TMW43042015 MS | Lab Sample ID: 280-67711-1 MS

Matrix: Water | Lab File ID: G6_17429.D

Analysis Method: 8270D | Date Collected: 04/10/2015 08:25

Extract. Method: 3520C | Date Extracted: 04/15/2015 18:50

Sample wt/vol: 885(mL) | Date Analyzed: 04/18/2015 21:39

Con. Extract Vol.: 1(mL) | Dilution Factor: 1

Injection Volume: 0.5(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 273380 | Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
105-60-2	Caprolactam	62.1		5.6	2.8

CAS NO.	SURROGATE	%REC	Q	LIMITS
118-79-6	2,4,6-Tribromophenol (Surr)	72		48-135
321-60-8	2-Fluorobiphenyl	61		48-135
367-12-4	2-Fluorophenol (Surr)	58		41-135
4165-60-0	Nitrobenzene-d5 (Surr)	58		42-135
4165-62-2	Phenol-d5 (Surr)	60		46-135
1718-51-0	Terphenyl-d14 (Surr)	65		20-135

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17429.D
Lims ID: 280-67711-B-1-B MS
Client ID: TMW43042015
Sample Type: MS
Inject. Date: 18-Apr-2015 21:39:30 ALS Bottle#: 17 Worklist Smp#: 30
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: 280-67711-B-1-CMS
Operator ID: HOEFLERA Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:13 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: RT Order ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: kiekeld Date: 20-Apr-2015 07:07:04

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	95	207648	40.0	40.0	
* 2 Naphthalene-d8	136	5.951	5.951	0.000	99	815083	40.0	40.0	
* 3 Acenaphthene-d10	164	7.704	7.704	0.000	93	448804	40.0	40.0	
* 4 Phenanthrene-d10	188	9.204	9.210	-0.006	97	777208	40.0	40.0	
* 5 Chrysene-d12	240	13.521	13.516	0.005	96	709191	40.0	40.0	
* 6 Perylene-d12	264	17.656	17.651	0.005	96	608383	40.0	40.0	
$ 7 2-Fluorophenol	112	3.487	3.493	-0.006	92	441230	100.0	58.4	
$ 8 Phenol-d5	99	4.328	4.334	-0.006	99	563129	100.0	59.5	
$ 9 Nitrobenzene-d5	82	5.240	5.245	-0.005	91	487702	100.0	58.4	
$ 10 2-Fluorobiphenyl	172	7.004	7.004	0.000	99	924919	100.0	61.0	
$ 11 2,4,6-Tribromophenol	330	8.498	8.504	-0.006	85	121838	100.0	72.2	
$ 12 Terphenyl-d14	244	11.233	11.239	-0.006	99	915343	100.0	65.0	
13 1,4-Dioxane	88	2.028	2.040	-0.012	98	142154	80.0	44.6	
14 N-Nitrosodimethylamine	74	2.299	2.305	-0.005	92	245493	80.0	50.7	
15 Pyridine	79	2.340	2.352	-0.012	91	340798	80.0	41.3	
16 2-Picoline	93		2.987					ND	
17 N-Nitrosomethylethylamine	88		3.081					ND	
18 Methyl methanesulfonate	80		3.351					ND	
19 N-Nitrosodiethylamine	102		3.704					ND	
20 Pentachlorophenol_T	266		3.822					ND	
21 Ethyl methanesulfonate	79		3.969					ND	
22 Pentachloroethane	117		4.446					ND	
23 Phenol	94	4.340	4.346	-0.006	99	465500	80.0	49.1	
24 Aniline	93	4.399	4.404	-0.006	98	452438	80.0	39.0	
25 Bis(2-chloroethyl)ether	93	4.434	4.434	0.000	96	370074	80.0	49.7	
26 2-Chlorophenol	128	4.510	4.516	-0.006	96	385730	80.0	49.8	
27 1,3-Dichlorobenzene	146	4.657	4.657	0.000	97	325204	80.0	41.4	
28 1,4-Dichlorobenzene	146	4.722	4.728	-0.006	92	337569	80.0	42.5	
29 Benzyl alcohol	108	4.822	4.828	-0.006	93	256476	80.0	52.5	
30 1,2-Dichlorobenzene	146	4.875	4.875	0.000	96	328260	80.0	43.0	
31 2-Methylphenol	108	4.916	4.916	0.000	94	346447	80.0	49.2	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
32 2,2'-oxybis[1-chloropropan	45	4.934	4.934	0.000	93	482803	80.0	44.0	
33 N-Nitrosopyrrolidine	100		5.075					ND	
34 N-Nitrosomorpholine	116		5.104					ND	
35 Benzidine_T	184		4.952					ND	
36 2-Toluidine	106		5.122					ND	
37 Benzaldehyde	106		5.015				ND	ND	
38 3 & 4 Methylphenol	108	5.057	5.063	-0.006	92	363840	80.0	49.9	
39 3-Methylphenol	108	5.057	5.063	-0.006	92	363840	80.0	49.9	
40 4-Methylphenol	108	5.057	5.063	-0.006	93	363840	80.0	49.9	
41 N-Nitrosodi-n-propylamine	70	5.063	5.069	-0.006	86	257220	80.0	48.6	
42 Acetophenone	105	5.081	5.087	-0.006	97	512292	80.0	50.0	
43 Hexachloroethane	117	5.198	5.198	0.000	92	136979	80.0	40.2	
44 Nitrobenzene	77	5.263	5.263	0.000	88	445661	80.0	55.4	
45 N-Nitrosopiperidine	114		5.410					ND	
46 Isophorone	82	5.475	5.475	0.000	99	710631	80.0	52.0	
47 o,o',o''-Triethylphosphoro	198		5.622					ND	
48 2-Nitrophenol	139	5.569	5.569	0.000	93	204550	80.0	51.0	
49 2,4-Dimethylphenol	107	5.575	5.581	-0.006	92	292006	80.0	39.7	
50 Bis(2-chloroethoxy)methane	93	5.663	5.669	-0.006	97	446476	80.0	49.7	
51 alpha,alpha-Dimethyl phene	58		5.792					ND	
52 Benzoic acid	105	5.651	5.687	-0.036	88	303041	80.0	52.5	
53 2,4-Dichlorophenol	162	5.798	5.798	0.000	94	307727	80.0	51.1	
54 1,2,4-Trichlorobenzene	180	5.881	5.887	-0.006	92	269867	80.0	44.1	
55 2,6-Dichlorophenol	162	6.022	6.022	0.000	94	300039	80.0	52.9	
56 Hexachloropropene	213		6.045					ND	
57 Naphthalene	128	5.969	5.975	-0.006	97	997817	80.0	46.0	
58 4-Chloroaniline	127	6.016	6.016	0.000	81	376360	80.0	38.4	
59 Hexachlorobutadiene	225	6.063	6.069	-0.006	95	136337	80.0	40.9	
60 N-Nitrosodi-n-butylamine	84		6.310					ND	
61 p-Phenylene diamine	108		6.363					ND	
62 Caprolactam	55	6.351	6.369	-0.018	76	212024	80.0	55.0	
63 Safrole, Total	162		6.540					ND	
64 4-Chloro-3-methylphenol	107	6.475	6.481	-0.006	96	338782	80.0	52.6	
65 2-Methylnaphthalene	142	6.651	6.651	0.000	91	681236	80.0	48.9	
66 Isosafrole Peak 1	162		6.834					ND	
67 1-Methylnaphthalene	142	6.751	6.757	-0.006	93	609893	80.0	46.5	
68 Hexachlorocyclopentadiene	237	6.792	6.798	-0.006	88	37392	80.0	10.6	
69 1,2,4,5-Tetrachlorobenzene	216	6.816	6.822	-0.006	97	284191	80.0	49.0	
70 2,4,6-Trichlorophenol	196	6.928	6.934	-0.006	93	221297	80.0	54.9	
71 Isosafrole Peak 2	104		7.063					ND	
72 2,4,5-Trichlorophenol	196	6.969	6.969	0.000	94	240330	80.0	56.0	
73 1-Chloronaphthalene	162		7.169					ND	
74 1,1'-Biphenyl	154	7.110	7.110	0.000	95	841751	80.0	49.8	
75 2-Chloronaphthalene	162	7.145	7.151	-0.006	97	641188	80.0	50.3	
76 1,4-Naphthoquinone	158		7.328					ND	
77 2-Nitroaniline	65	7.245	7.251	-0.006	83	244042	80.0	52.2	
78 1,4-Dinitrobenzene	168		7.381					ND	
79 Dimethyl phthalate	163	7.392	7.398	-0.006	98	806565	80.0	54.0	
80 1,3-Dinitrobenzene	168	7.457	7.463	-0.006	87	157949	80.0	59.6	
81 2,6-Dinitrotoluene	165	7.475	7.481	-0.006	95	196719	80.0	57.0	
82 Acenaphthylene	152	7.569	7.569	0.000	92	1109350	80.0	53.6	
83 3-Nitroaniline	138	7.657	7.663	-0.006	94	184282	80.0	41.3	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
84 Acenaphthene	153	7.739	7.739	0.000	94	685401	80.0	51.3	
85 Pentachlorobenzene	250		7.863					ND	
86 2,4-Dinitrophenol	184	7.757	7.763	-0.006	81	286265	160.0	115.5	
87 4-Nitrophenol	109	7.810	7.816	-0.006	94	278351	160.0	115.6	
88 1-Naphthylamine	143		7.998					ND	
89 2,4-Dinitrotoluene	165	7.881	7.887	-0.005	89	273025	80.0	60.3	
90 Dibenzofuran	168	7.910	7.910	0.000	91	994810	80.0	52.9	
91 2-Naphthylamine	143		8.075					ND	
92 2,3,4,6-Tetrachlorophenol	232	8.028	8.028	0.000	73	200079	80.0	57.5	
93 Thionazin	97		8.169					ND	
94 Diethyl phthalate	149	8.087	8.092	-0.006	98	880091	80.0	57.9	
95 N-Nitro-o-toluidine	152		8.263					ND	
96 4-Chlorophenyl phenyl ethe	204	8.228	8.234	-0.006	94	359224	80.0	52.8	
97 Diphenylamine	169	8.351	8.351	0.000	95	1181676	137.1	92.1	
98 Fluorene	166	8.251	8.257	-0.006	84	833578	80.0	54.5	
99 4-Nitroaniline	138	8.281	8.286	-0.005	81	209512	80.0	48.0	
100 4,6-Dinitro-2-methylphenol	198	8.292	8.298	-0.006	81	331008	160.0	120.4	
101 Sulfotepp	97		8.451					ND	
102 N-Nitrosodiphenylamine	169	8.351	8.357	-0.006	61	1181676	160.0	102.6	
103 Azobenzene	77	8.392	8.392	0.000	99	850210	80.0	50.6	
104 1,2-Diphenylhydrazine	77	8.392	8.392	0.000	100	850210	80.9	51.1	
105 Diallate Peak 1	86		8.610					ND	
106 1,3,5-Trinitrobenzene	213		8.622					ND	
107 Phorate	121		8.628					ND	
108 Phenacetin	108		8.645					ND	
109 Diallate Peak 2	86		8.704					ND	
110 Dimethoate	87		8.822					ND	
111 4-Bromophenyl phenyl ether	248	8.722	8.728	-0.006	69	198713	80.0	54.6	
112 Hexachlorobenzene	284	8.804	8.810	-0.006	87	208252	80.0	55.1	
113 4-Aminobiphenyl	169		8.998					ND	
114 Pentachloronitrobenzene	237		9.010					ND	
115 Pronamide	173		9.016					ND	
116 Pentachlorophenol	266	9.004	9.010	-0.006	87	258819	160.0	104.1	
117 Disulfoton	88		9.145					ND	
118 Dinoseb	211		9.157					ND	
119 Phenanthrene	178	9.228	9.233	-0.005	98	1220612	80.0	56.6	
120 Anthracene	178	9.281	9.286	-0.005	98	1209310	80.0	54.7	
121 Methyl parathion	109		9.551					ND	
122 Carbazole	167	9.439	9.445	-0.006	82	1274869	80.0	56.6	
123 Di-n-butyl phthalate	149	9.733	9.739	-0.006	100	1508815	80.0	57.0	
124 Ethyl Parathion	109		9.980					ND	
125 4-Nitroquinoline-1-oxide	190		10.104					ND	
126 Methapyrilene	97		10.128					ND	
127 Isodrin	193		10.433					ND	
128 Fluoranthene	202	10.663	10.675	-0.012	98	1306451	80.0	55.3	
129 Benzidine	184		10.466				ND	ND	
130 Aramite Peak 1	185		11.139					ND	
131 Pyrene	202	11.027	11.039	-0.012	97	1376232	80.0	58.1	
132 Aramite Peak 2	185		11.274					ND	
133 p-Dimethylamino azobenzene	120		11.492					ND	
134 Chlorobenzilate	251		11.545					ND	
135 3,3'-Dimethylbenzidine	212		12.180					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
136 Famphur	218		12.045					ND	
137 Butyl benzyl phthalate	149	12.139	12.151	-0.012	97	676371	80.0	60.5	
138 2-Acetylaminofluorene	181		12.745					ND	
139 4,4'-Methylene bis(2-chlor	231		13.463					ND	
140 3,3'-Dichlorobenzidine	252	13.451	13.468	-0.017	70	138395	80.0	19.7	
141 Benzo[a]anthracene	228	13.492	13.504	-0.012	99	1140467	80.0	56.0	
142 Bis(2-ethylhexyl) phthalat	149	13.480	13.492	-0.012	92	893451	80.0	61.2	
143 Chrysene	228	13.586	13.598	-0.012	96	1090182	80.0	57.3	
144 Di-n-octyl phthalate	149	15.351	15.362	-0.011	99	1530362	80.0	59.2	
145 7,12-Dimethylbenz(a)anthra	256		16.480					ND	
146 Benzo[b]fluoranthene	252	16.492	16.509	-0.017	96	994304	80.0	55.9	
147 Benzo[k]fluoranthene	252	16.574	16.592	-0.018	98	1043489	80.0	57.2	
148 Benzo[a]pyrene	252	17.474	17.492	-0.018	77	950803	80.0	55.6	
149 3-Methylcholanthrene	268		18.550					ND	
150 Dibenz[a,j]acridine	279		20.350					ND	
151 Indeno[1,2,3-cd]pyrene	276	20.839	20.862	-0.023	97	779069	80.0	53.8	M
152 Dibenz(a,h)anthracene	278	20.921	20.939	-0.018	83	792435	80.0	56.1	
153 Benzo[g,h,i]perylene	276	21.615	21.633	-0.018	90	845314	80.0	55.0	
S 160 Aramite, Total	185		15.047					ND	
S 161 Isosafrole	162		15.047					ND	
S 162 Diallate	86		15.047					ND	
154 Total Cresols	1		0.000					ND	
155 Tetraethyl Pyrophosphate (	1		0.000					ND	
157 4,4'-DDE	246		5.063					ND	
158 4,4'-DDD	235		5.375					ND	
159 4,4'-DDT	235		5.604					ND	
S 163 Methyl Phenols,Total	1		0.000					ND	

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

Review Flags

M - Manually Integrated

Reagents:

MS-IS_00007 Amount Added: 20.00 Units: uL Run Reagent

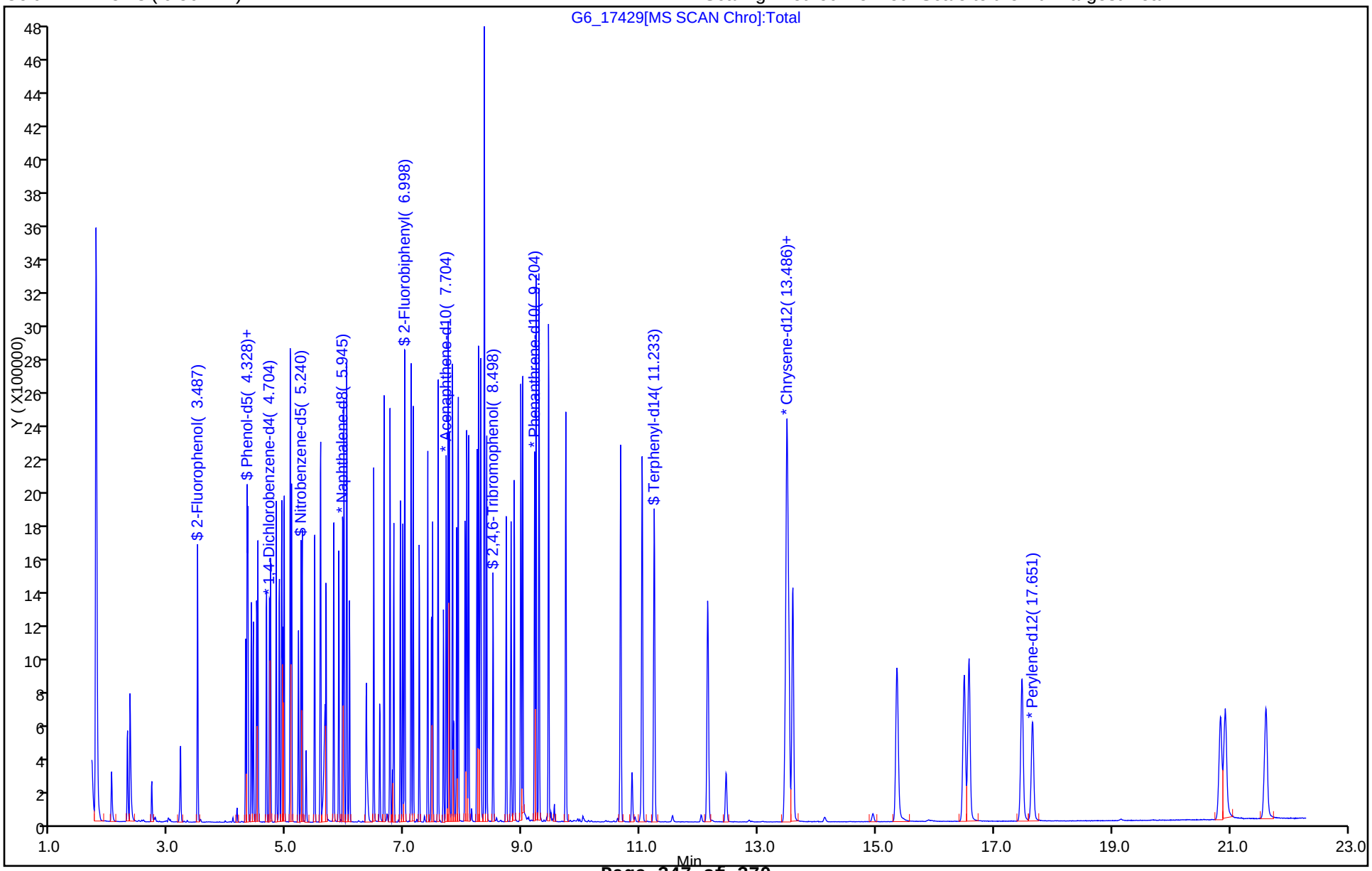
Report Date: 20-Apr-2015 07:09:16 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17429.D

Injection Date:	18-Apr-2015 21:39:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	280-67711-B-1-B MS			Worklist Smp#:	30
Client ID:	TMW43042015				
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	17
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



FORM I
GC/MS SEMI VOA ORGANICS ANALYSIS DATA SHEET

Lab Name: TestAmerica Denver | Job No.: 280-67711-2

SDG No.:

Client Sample ID: TMW43042015 MSD | Lab Sample ID: 280-67711-1 MSD

Matrix: Water | Lab File ID: G6_17430.D

Analysis Method: 8270D | Date Collected: 04/10/2015 08:25

Extract. Method: 3520C | Date Extracted: 04/15/2015 18:50

Sample wt/vol: 903.3(mL) | Date Analyzed: 04/18/2015 22:06

Con. Extract Vol.: 1(mL) | Dilution Factor: 1

Injection Volume: 0.5(uL) | Level: (low/med) Low

% Moisture: | GPC Cleanup:(Y/N) N

Analysis Batch No.: 273380 | Units: ug/L

CAS NO.	COMPOUND NAME	RESULT	Q	RL	MDL
105-60-2	Caprolactam	58.4		5.5	2.8

CAS NO.	SURROGATE	%REC	Q	LIMITS
118-79-6	2,4,6-Tribromophenol (Surr)	66		48-135
321-60-8	2-Fluorobiphenyl	61		48-135
367-12-4	2-Fluorophenol (Surr)	62		41-135
4165-60-0	Nitrobenzene-d5 (Surr)	63		42-135
4165-62-2	Phenol-d5 (Surr)	62		46-135
1718-51-0	Terphenyl-d14 (Surr)	62		20-135

TestAmerica Denver
Target Compound Quantitation Report

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17430.D
Lims ID: 280-67711-L-1-B MSD
Client ID: TMW43042015
Sample Type: MSD
Inject. Date: 18-Apr-2015 22:06:30 ALS Bottle#: 18 Worklist Smp#: 31
Injection Vol: 0.5 ul Dil. Factor: 1.0000
Sample Info: 280-67711-L-1-CMSD
Operator ID: HOEFLERA Instrument ID: SMS_G6

Method: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\SMS_G6_8270D.m
Limit Group: MSSV - 8270D
Method Label: 8270D
Last Update: 20-Apr-2015 07:09:13 Calib Date: 26-Feb-2015 10:02:30
Integrator: RTE ID Type: RT Order ID
Quant Method: Internal Standard Quant By: Initial Calibration
Last ICal File: \\Denchrom\ChromData\SMS_G6\20150416-33972.b\G6_16673.D

Column 1 : VF-5ms (0.50 mm) Det: MS SCAN
Process Host: XAWRK031

First Level Reviewer: kiekeld Date: 20-Apr-2015 07:08:15

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
* 1 1,4-Dichlorobenzene-d4	152	4.710	4.710	0.000	95	222633	40.0	40.0	
* 2 Naphthalene-d8	136	5.951	5.951	0.000	99	860621	40.0	40.0	
* 3 Acenaphthene-d10	164	7.704	7.704	0.000	92	488251	40.0	40.0	
* 4 Phenanthrene-d10	188	9.204	9.210	-0.006	97	823179	40.0	40.0	
* 5 Chrysene-d12	240	13.516	13.516	0.000	96	756154	40.0	40.0	
* 6 Perylene-d12	264	17.651	17.651	0.000	95	637043	40.0	40.0	
$ 7 2-Fluorophenol	112	3.487	3.493	-0.006	92	502901	100.0	62.1	
$ 8 Phenol-d5	99	4.328	4.334	-0.006	99	628452	100.0	61.9	
$ 9 Nitrobenzene-d5	82	5.240	5.245	-0.005	91	554318	100.0	62.8	
$ 10 2-Fluorobiphenyl	172	7.004	7.004	0.000	100	1003669	100.0	60.8	
$ 11 2,4,6-Tribromophenol	330	8.498	8.504	-0.006	84	121473	100.0	66.2	
$ 12 Terphenyl-d14	244	11.233	11.239	-0.006	99	930081	100.0	61.9	
13 1,4-Dioxane	88	2.028	2.040	-0.012	97	160580	80.0	47.0	
14 N-Nitrosodimethylamine	74	2.299	2.305	-0.005	91	279265	80.0	53.8	
15 Pyridine	79	2.340	2.352	-0.012	92	388641	80.0	43.9	
16 2-Picoline	93		2.987					ND	
17 N-Nitrosomethylethylamine	88		3.081					ND	
18 Methyl methanesulfonate	80		3.351					ND	
19 N-Nitrosodiethylamine	102		3.704					ND	
20 Pentachlorophenol_T	266		3.822					ND	
21 Ethyl methanesulfonate	79		3.969					ND	
22 Pentachloroethane	117		4.446					ND	
23 Phenol	94	4.340	4.346	-0.006	99	524550	80.0	51.6	
24 Aniline	93	4.399	4.404	-0.006	98	506394	80.0	40.7	
25 Bis(2-chloroethyl)ether	93	4.434	4.434	0.000	95	423777	80.0	53.1	
26 2-Chlorophenol	128	4.510	4.516	-0.006	97	438500	80.0	52.8	
27 1,3-Dichlorobenzene	146	4.657	4.657	0.000	97	377084	80.0	44.7	
28 1,4-Dichlorobenzene	146	4.722	4.728	-0.006	92	385662	80.0	45.3	
29 Benzyl alcohol	108	4.822	4.828	-0.006	93	285266	80.0	54.5	
30 1,2-Dichlorobenzene	146	4.875	4.875	0.000	96	380905	80.0	46.5	
31 2-Methylphenol	108	4.916	4.916	0.000	95	386083	80.0	51.2	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
32 2,2'-oxybis[1-chloropropan	45	4.934	4.934	0.000	93	555549	80.0	47.2	
33 N-Nitrosopyrrolidine	100		5.075					ND	
34 N-Nitrosomorpholine	116		5.104					ND	
35 Benzidine_T	184		4.952					ND	
36 2-Toluidine	106		5.122					ND	
37 Benzaldehyde	106		5.015				ND	ND	
38 3 & 4 Methylphenol	108	5.063	5.063	0.000	87	410524	80.0	52.5	
39 3-Methylphenol	108	5.063	5.063	0.000	87	410524	80.0	52.5	
40 4-Methylphenol	108	5.063	5.063	0.000	89	410524	80.0	52.5	
41 N-Nitrosodi-n-propylamine	70	5.063	5.069	-0.006	84	290693	80.0	51.2	
42 Acetophenone	105	5.081	5.087	-0.006	96	581887	80.0	53.0	
43 Hexachloroethane	117	5.198	5.198	0.000	93	157274	80.0	43.0	
44 Nitrobenzene	77	5.263	5.263	0.000	89	516712	80.0	60.8	
45 N-Nitrosopiperidine	114		5.410					ND	
46 Isophorone	82	5.475	5.475	0.000	99	792437	80.0	54.9	
47 o,o',o''-Triethylphosphoro	198		5.622					ND	
48 2-Nitrophenol	139	5.569	5.569	0.000	94	231023	80.0	54.5	
49 2,4-Dimethylphenol	107	5.575	5.581	-0.006	92	302059	80.0	38.9	
50 Bis(2-chloroethoxy)methane	93	5.669	5.669	0.000	99	500669	80.0	52.8	
51 alpha,alpha-Dimethyl phene	58		5.792					ND	
52 Benzoic acid	105	5.651	5.687	-0.036	88	329552	80.0	54.1	
53 2,4-Dichlorophenol	162	5.798	5.798	0.000	95	347259	80.0	54.6	
54 1,2,4-Trichlorobenzene	180	5.881	5.887	-0.006	93	314578	80.0	48.7	
55 2,6-Dichlorophenol	162	6.022	6.022	0.000	94	335720	80.0	56.1	
56 Hexachloropropene	213		6.045					ND	
57 Naphthalene	128	5.969	5.975	-0.006	97	1128991	80.0	49.3	
58 4-Chloroaniline	127	6.016	6.016	0.000	81	392918	80.0	37.9	
59 Hexachlorobutadiene	225	6.063	6.069	-0.006	95	160139	80.0	45.5	
60 N-Nitrosodi-n-butylamine	84		6.310					ND	
61 p-Phenylene diamine	108		6.363					ND	
62 Caprolactam	55	6.351	6.369	-0.018	78	214704	80.0	52.7	
63 Safrole, Total	162		6.540					ND	
64 4-Chloro-3-methylphenol	107	6.475	6.481	-0.006	94	365815	80.0	53.7	
65 2-Methylnaphthalene	142	6.651	6.651	0.000	92	770374	80.0	52.4	
66 Isosafrole Peak 1	162		6.834					ND	
67 1-Methylnaphthalene	142	6.751	6.757	-0.006	92	676939	80.0	48.9	
68 Hexachlorocyclopentadiene	237	6.792	6.798	-0.006	92	21667	80.0	5.65	7
69 1,2,4,5-Tetrachlorobenzene	216	6.816	6.822	-0.006	96	315008	80.0	51.4	
70 2,4,6-Trichlorophenol	196	6.928	6.934	-0.006	92	235449	80.0	53.7	
71 Isosafrole Peak 2	104		7.063					ND	
72 2,4,5-Trichlorophenol	196	6.969	6.969	0.000	94	254686	80.0	54.5	
73 1-Chloronaphthalene	162		7.169					ND	
74 1,1'-Biphenyl	154	7.110	7.110	0.000	94	932585	80.0	50.7	
75 2-Chloronaphthalene	162	7.145	7.151	-0.006	97	707726	80.0	51.1	
76 1,4-Naphthoquinone	158		7.328					ND	
77 2-Nitroaniline	65	7.245	7.251	-0.006	82	253710	80.0	49.9	
78 1,4-Dinitrobenzene	168		7.381					ND	
79 Dimethyl phthalate	163	7.392	7.398	-0.006	98	833062	80.0	51.3	
80 1,3-Dinitrobenzene	168	7.457	7.463	-0.006	89	164060	80.0	56.9	
81 2,6-Dinitrotoluene	165	7.475	7.481	-0.006	95	207634	80.0	55.3	
82 Acenaphthylene	152	7.569	7.569	0.000	94	1198003	80.0	53.2	
83 3-Nitroaniline	138	7.657	7.663	-0.006	93	175374	80.0	36.1	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
84 Acenaphthene	153	7.739	7.739	0.000	94	737152	80.0	50.8	
85 Pentachlorobenzene	250		7.863					ND	
86 2,4-Dinitrophenol	184	7.757	7.763	-0.006	81	292006	160.0	108.5	
87 4-Nitrophenol	109	7.810	7.816	-0.006	94	277267	160.0	105.9	
88 1-Naphthylamine	143		7.998					ND	
89 2,4-Dinitrotoluene	165	7.881	7.887	-0.005	84	271353	80.0	55.1	
90 Dibenzofuran	168	7.910	7.910	0.000	90	1061122	80.0	51.9	
91 2-Naphthylamine	143		8.075					ND	
92 2,3,4,6-Tetrachlorophenol	232	8.028	8.028	0.000	72	207035	80.0	54.7	
93 Thionazin	97		8.169					ND	
94 Diethyl phthalate	149	8.086	8.092	-0.006	98	888032	80.0	53.7	
95 N-Nitro-o-toluidine	152		8.263					ND	
96 4-Chlorophenyl phenyl ethe	204	8.228	8.234	-0.006	95	382804	80.0	51.7	
97 Diphenylamine	169	8.351	8.351	0.000	95	1187951	137.1	85.1	
98 Fluorene	166	8.251	8.257	-0.006	85	881278	80.0	52.9	
99 4-Nitroaniline	138	8.281	8.286	-0.005	81	206704	80.0	43.5	
100 4,6-Dinitro-2-methylphenol	198	8.292	8.298	-0.006	82	338852	160.0	116.3	
101 Sulfotepp	97		8.451					ND	
102 N-Nitrosodiphenylamine	169	8.351	8.357	-0.006	61	1187951	160.0	97.4	
103 Azobenzene	77	8.392	8.392	0.000	99	885645	80.0	48.4	
104 1,2-Diphenylhydrazine	77	8.392	8.392	0.000	99	885645	80.9	49.0	
105 Diallate Peak 1	86		8.610					ND	
106 1,3,5-Trinitrobenzene	213		8.622					ND	
107 Phorate	121		8.628					ND	
108 Phenacetin	108		8.645					ND	
109 Diallate Peak 2	86		8.704					ND	
110 Dimethoate	87		8.822					ND	
111 4-Bromophenyl phenyl ether	248	8.722	8.728	-0.006	69	206339	80.0	53.5	
112 Hexachlorobenzene	284	8.804	8.810	-0.006	87	212748	80.0	53.1	
113 4-Aminobiphenyl	169		8.998					ND	
114 Pentachloronitrobenzene	237		9.010					ND	
115 Pronamide	173		9.016					ND	
116 Pentachlorophenol	266	9.004	9.010	-0.006	88	263968	160.0	100.2	
117 Disulfoton	88		9.145					ND	
118 Dinoseb	211		9.157					ND	
119 Phenanthrene	178	9.228	9.233	-0.005	98	1223368	80.0	53.6	
120 Anthracene	178	9.281	9.286	-0.005	98	1213940	80.0	51.9	
121 Methyl parathion	109		9.551					ND	
122 Carbazole	167	9.439	9.445	-0.006	82	1261491	80.0	52.9	
123 Di-n-butyl phthalate	149	9.733	9.739	-0.006	100	1502398	80.0	53.6	
124 Ethyl Parathion	109		9.980					ND	
125 4-Nitroquinoline-1-oxide	190		10.104					ND	
126 Methapyrilene	97		10.128					ND	
127 Isodrin	193		10.433					ND	
128 Fluoranthene	202	10.663	10.675	-0.012	98	1296626	80.0	51.9	
129 Benzidine	184		10.466				ND	ND	
130 Aramite Peak 1	185		11.139					ND	
131 Pyrene	202	11.027	11.039	-0.012	97	1362039	80.0	53.9	
132 Aramite Peak 2	185		11.274					ND	
133 p-Dimethylamino azobenzene	120		11.492					ND	
134 Chlorobenzilate	251		11.545					ND	
135 3,3'-Dimethylbenzidine	212		12.180					ND	

Compound	Sig	RT (min.)	Exp RT (min.)	Dlt RT (min.)	Q	Response	Cal Amt ug/ml	OnCol Amt ug/ml	Flags
136 Famphur	218		12.045					ND	
137 Butyl benzyl phthalate	149	12.139	12.151	-0.012	97	663246	80.0	55.6	
138 2-Acetylaminofluorene	181		12.745					ND	
139 4,4'-Methylene bis(2-chlor	231		13.463					ND	
140 3,3'-Dichlorobenzidine	252	13.451	13.468	-0.017	68	121422	80.0	16.2	
141 Benzo[a]anthracene	228	13.492	13.504	-0.012	99	1121851	80.0	51.6	
142 Bis(2-ethylhexyl) phthalat	149	13.480	13.492	-0.012	92	877691	80.0	56.4	
143 Chrysene	228	13.586	13.598	-0.012	96	1068504	80.0	52.7	
144 Di-n-octyl phthalate	149	15.351	15.362	-0.011	99	1505903	80.0	54.8	
145 7,12-Dimethylbenz(a)anthra	256		16.480					ND	
146 Benzo[b]fluoranthene	252	16.492	16.509	-0.017	94	977550	80.0	52.5	
147 Benzo[k]fluoranthene	252	16.574	16.592	-0.018	98	1015707	80.0	53.1	
148 Benzo[a]pyrene	252	17.474	17.492	-0.018	78	906481	80.0	50.7	
149 3-Methylcholanthrene	268		18.550					ND	
150 Dibenz[a,j]acridine	279		20.350					ND	
151 Indeno[1,2,3-cd]pyrene	276	20.844	20.862	-0.018	97	766686	80.0	49.7	M
152 Dibenz(a,h)anthracene	278	20.921	20.939	-0.018	87	753533	80.0	50.9	
153 Benzo[g,h,i]perylene	276	21.609	21.633	-0.024	89	818821	80.0	50.9	
S 160 Aramite, Total	185		15.047					ND	
S 161 Isosafrole	162		15.047					ND	
S 162 Diallate	86		15.047					ND	
154 Total Cresols	1		0.000					ND	
155 Tetraethyl Pyrophosphate (	1		0.000					ND	
157 4,4'-DDE	246		5.063					ND	
158 4,4'-DDD	235		5.375					ND	
159 4,4'-DDT	235		5.604					ND	
S 163 Methyl Phenols,Total	1		0.000					ND	

QC Flag Legend

Processing Flags

ND - Not Detected or Marked ND

7 - Failed Limit of Detection

Review Flags

M - Manually Integrated

Reagents:

MS-IS_00007 Amount Added: 20.00 Units: uL Run Reagent

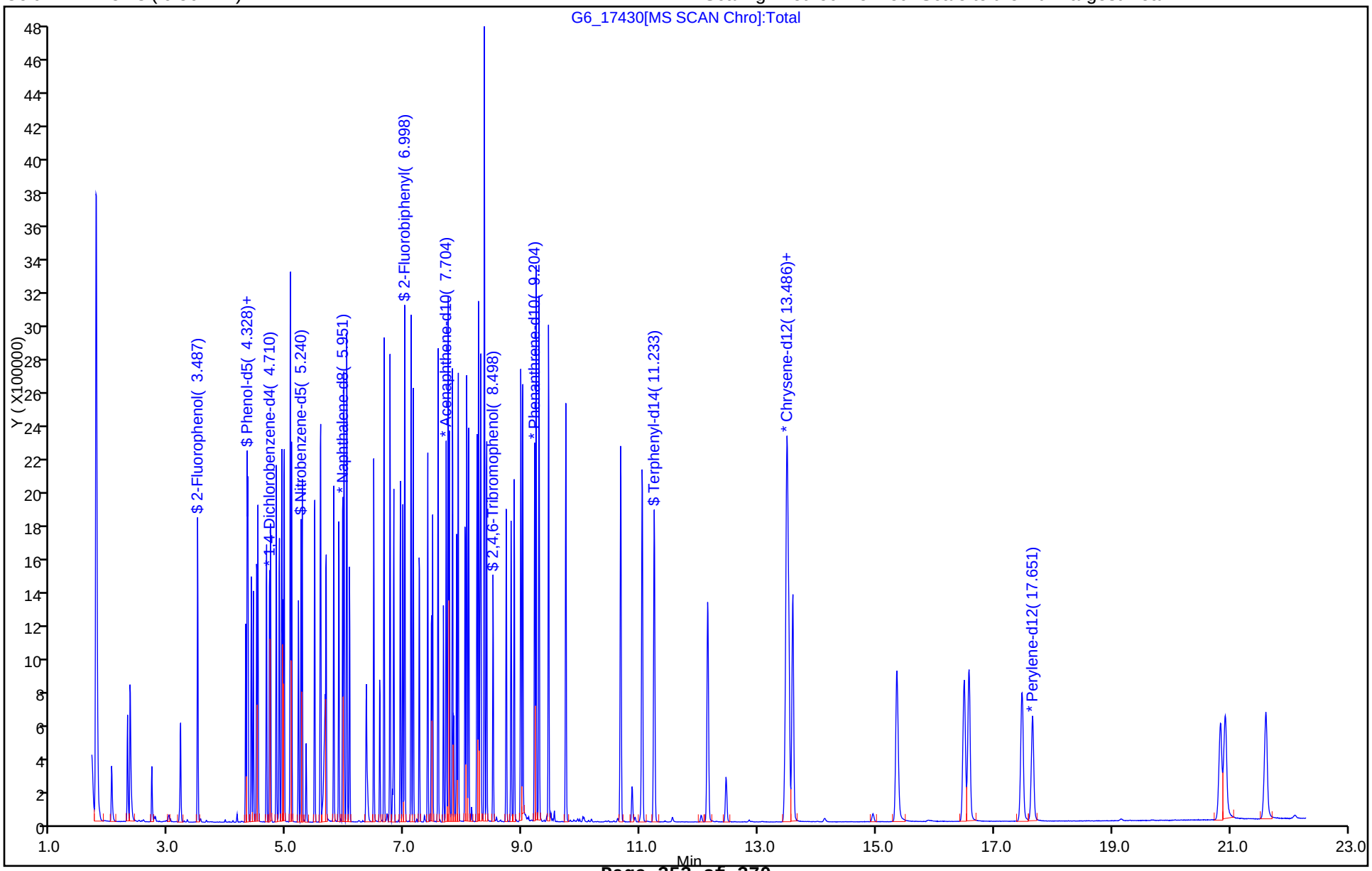
Report Date: 20-Apr-2015 07:09:17 Chrom Revision: 2.2 09-Apr-2015 10:05:40

TestAmerica Denver

Data File: \\Denchrom\ChromData\SMS_G6\20150418-34074.b\G6_17430.D

Injection Date:	18-Apr-2015 22:06:30	Instrument ID:	SMS_G6	Operator ID:	HOEFLERA
Lims ID:	280-67711-L-1-B MSD			Worklist Smp#:	31
Client ID:	TMW43042015				
Injection Vol:	0.5 ul	Dil. Factor:	1.0000	ALS Bottle#:	18
Method:	SMS_G6_8270D	Limit Group:	MSSV - 8270D		

Column: VF-5ms (0.50 mm) Y Scaling: Method Defined: Scale to the Nth Largest Peak: 1



GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Instrument ID: SMS_G6 Start Date: 02/25/2015 11:44

Analysis Batch Number: 272059 End Date: 02/25/2015 16:19

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
DFTPP 280-272059/2		02/25/2015 11:44	1	G6_16636.D	Vf-5MS (30.25) 0.25(mm)
ICIS 280-272059/3		02/25/2015 11:53	1	G6_16637.D	Vf-5MS (30.25) 0.25(mm)
STD004 280-272059/4 IC		02/25/2015 12:19	1	G6_16638.D	Vf-5MS (30.25) 0.25(mm)
STD010 280-272059/5 IC		02/25/2015 12:46	1	G6_16639.D	Vf-5MS (30.25) 0.25(mm)
STD020 280-272059/6 IC		02/25/2015 13:12	1	G6_16640.D	Vf-5MS (30.25) 0.25(mm)
STD050 280-272059/7 IC		02/25/2015 13:39	1	G6_16641.D	Vf-5MS (30.25) 0.25(mm)
STD120 280-272059/8 IC		02/25/2015 14:06	1	G6_16642.D	Vf-5MS (30.25) 0.25(mm)
STD160 280-272059/9 IC		02/25/2015 14:32	1	G6_16643.D	Vf-5MS (30.25) 0.25(mm)
STD200 280-272059/10 IC		02/25/2015 14:59	1	G6_16644.D	Vf-5MS (30.25) 0.25(mm)
ICV 280-272059/11		02/25/2015 15:26	1	G6_16645.D	Vf-5MS (30.25) 0.25(mm)
ICV 280-272059/12		02/25/2015 15:52	1	G6_16646.D	Vf-5MS (30.25) 0.25(mm)
ICV 280-272059/13		02/25/2015 16:19	1	G6_16647.D	Vf-5MS (30.25) 0.25(mm)

GC/MS SEMI VOA ANALYSIS RUN LOG

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Instrument ID: SMS_G6 Start Date: 04/18/2015 15:09

Analysis Batch Number: 273380 End Date: 04/18/2015 23:26

LAB SAMPLE ID	CLIENT SAMPLE ID	DATE ANALYZED	DILUTION FACTOR	LAB FILE ID	COLUMN ID
DFTPP 280-273380/2		04/18/2015 15:09	1	G6_17413.D	Vf-5MS (30.25) 0.25(mm)
CCV 280-273380/3		04/18/2015 15:20	1	G6_17414.D	Vf-5MS (30.25) 0.25(mm)
CCV 280-273380/4		04/18/2015 16:09	1	G6_17416.D	Vf-5MS (30.25) 0.25(mm)
MB 280-272870/1-A		04/18/2015 17:49	1	G6_17420.D	Vf-5MS (30.25) 0.25(mm)
LCS 280-272870/2-A		04/18/2015 18:16	1	G6_17421.D	Vf-5MS (30.25) 0.25(mm)
280-67711-1	TMW43042015	04/18/2015 21:13	1	G6_17428.D	Vf-5MS (30.25) 0.25(mm)
280-67711-1 MS	TMW43042015 MS	04/18/2015 21:39	1	G6_17429.D	Vf-5MS (30.25) 0.25(mm)
280-67711-1 MSD	TMW43042015 MSD	04/18/2015 22:06	1	G6_17430.D	Vf-5MS (30.25) 0.25(mm)
280-67711-2	DTW43042015	04/18/2015 22:32	1	G6_17431.D	Vf-5MS (30.25) 0.25(mm)
280-67711-4	TMW47042015	04/18/2015 22:59	1	G6_17432.D	Vf-5MS (30.25) 0.25(mm)
ZZZZZ		04/18/2015 23:26	1		Vf-5MS (30.25) 0.25(mm)

GC/MS SVOAContinuing Calibration Review Checklist *TestAmerica*
Denver

Instrument ID and Date: **G6_HSL/AFC/AP9/BZHD/CUST/PHE_8270D_041815** Work List **34074/34075**

Check Method Used: Analysis ☐ 625 ☒ 8270 ☐ 8270_SIM ☐ Other SV _____

Review Items	Level 1 Yes	Level 1 No	Level 1 N/A	Level 2	Comments
Continuing Calibration					
1. DFTPP meets criteria?	X			✓	
2. ICAL date and instrument ID verified?	X			✓	
3. Do SPCC RRFs and CCC %Ds meet method criteria?	X			✓	
4. Does %D meet criteria for non-CCC compounds?	X			✓	Circle one: DOD / non-DOD
5. Isomeric pairs checked for correct peak assignment? Aniline and bis(2-chloroethyl) ether 1,3-Dichlorobenzene, 1,4-dichlorobenzene and 1,2-dichlorobenzene 2-Methylphenol, 3/4-Methylphenol 2,4-Dimethylphenol and 3,5-dimethylphenol 2,4,6-Trichlorophenol and 2,4,5-Trichlorophenol Phenanthrene and anthracene Fluoranthene and pyrene Benzo(a)anthracene and chrysene Benzo(e)pyrene and benzo(a)pyrene Benzo(b) and (k)fluoranthene	X			✓	
6. Standards traceability properly documented?	X			✓	
7. Manual integrations documented and checked?	X			✓	
8. Do the Internal Standards meet criteria for %D against ICAL?	X			✓	

1st Level Reviewer: Alex Hoffler Date: 042015

2nd Level Reviewer: Dawn Dyer Date: 042015

Alex Hoffer 04/20/15

Injection Log

Directory: C:\HPCHEM\1\DATA\G6\041815.B

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	100	G6_17412.D	1.	RINSE		18 Apr 2015 14:43
2	1	G6_17413.D	1.	DFTPP		18 Apr 2015 15:09
3	2	G6_17414.D	1.	CCV HSL		18 Apr 2015 15:20
4	3	G6_17415.D	1.	CCV AFC		18 Apr 2015 15:47
5	4	G6_17416.D	1.	CCV AP9		18 Apr 2015 16:09
6	5	G6_17417.D	1.	CCV BZHD		18 Apr 2015 16:34
7	6	G6_17418.D	1.	CCV CUST		18 Apr 2015 16:56
8	7	G6_17419.D	1.	CCV PHE		18 Apr 2015 17:23
9	8	G6_17420.D	1.	MB280-273368_1-A		18 Apr 2015 17:49
10	9	G6_17421.D	1.	LCS280-273368_2-A		18 Apr 2015 18:16
11	10	G6_17422.D	1.	MB280-273142_1-A		18 Apr 2015 18:43
12	11	G6_17423.D	1.	LCS280-273142_2-A		18 Apr 2015 19:05
13	12	G6_17424.D	1.	LCSD280-273142_3-A		18 Apr 2015 19:31
14	13	G6_17425.D	1.	LB3280-270930_1-D		18 Apr 2015 19:53
15	14	G6_17426.D	1.	LCS280-270930_2-D		18 Apr 2015 20:19
16	15	G6_17427.D	1.	160-11229-C-3-B		18 Apr 2015 20:46
17	16	G6_17428.D	1.	280-67711-C-1-B		18 Apr 2015 21:13
18	17	G6_17429.D	1.	280-67711-B-1-CMS		18 Apr 2015 21:39
19	18	G6_17430.D	1.	280-67711-L-1-CMSD		18 Apr 2015 22:06
20	19	G6_17431.D	1.	280-67711-B-2-B		18 Apr 2015 22:32
21	20	G6_17432.D	1.	280-67711-C-4-B		18 Apr 2015 22:59
22	21	G6_17433.D	1.	280-67625-A-9-B		18 Apr 2015 23:26
23	22	G6_17434.D	1.	280-67852-C-1-C 5X - NC		18 Apr 2015 23:52
24	23	G6_17435.D	1.	280-67852-C-2-C 5X - NC		19 Apr 2015 00:19
25	24	G6_17436.D	1.	280-67261-L-1-H		19 Apr 2015 00:45
26	25	G6_17437.D	1.	280-67261-L-1-IMS		19 Apr 2015 01:12
27	26	G6_17438.D	1.	280-67261-L-1-JMSD		19 Apr 2015 01:39
28	98	G6_17439.D	1.	RINSE		19 Apr 2015 02:05
29	99	G6_17440.D	1.	RINSE		19 Apr 2015 02:32
30	100	G6_17441.D	1.	RINSE		19 Apr 2015 02:59
31	98	G6_17442.D	1.	RINSE		19 Apr 2015 03:25
32	99	G6_17443.D	1.	RINSE		19 Apr 2015 03:52
33	100	G6_17444.D	1.	RINSE		19 Apr 2015 04:19
34		G6_17445.D	1.			

Dilution Solvent Lot #: 87975 **Pipette ID:** SV-18/SV-23 **Method(s) Performed:** 8270 DoD/8270C-UTS/8270C

Daily Maintenance: Check box if maintenance was performed. No mark indicates the item was not changed.

Changed Septum ☑ **Changed Liner** ☑ **Changed Seal** ☑ **Changed Ferrule** ☐ **Clipped Column** ☑

0.1 m 77

TestAmerica Denver
GC/MS SVOA Initial Calibration Review Checklist

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

Instrument ID and Date: G6 HSL-8270D 022515
Calibration Event 21947 Batch 272059

Check Method Used: Analysis ☐ 625 ☒ 8270 ☐ Other SV _____

Review Items	----- Level 1 ----- Yes	No	N/A	Level 2	Comments
Initial Calibration					
1. DFTPP meets criteria?	X			✓	
2. ICAL date and instrument ID verified?	X			✓	
3. Does the Form VI match the data in the CHROM source method?	X			✓	
4. Sufficient number of calibration points used?	X			✓	
5. Reasons for removal of points documented?	X			✓	Some points < RL removed
6. %RSD or correlation coefficient within method limits?	X			✓	
7. If RRF used for ICAL, were all compounds within 20% RSD?			X	✓ NA	List all exceptions below (cpd & RSD)
8. Response factors meet criteria?	X			✓	
9. Isomeric pairs checked for correct peak assignment?	X			✓	
10. Data checked for detector saturation?	X			✓	
11. Standards traceability properly documented?	X			✓	
12. Manual integrations documented and checked?	X			✓	
13. 2nd source ICV recovery 75-125% (±25%drift)for DoD projects, 65-135% (±35%,or ±55% of expected for poor performers) for non-DoD? Exceptions noted in comment section.	**X**			✓	

1st Level Reviewer: _____ Date: 04/10/15

2nd Level Reviewer: _____ Date: 04/10/15

Injection Log

Directory: C:\HPCHEM\1\DATA\G6\022515.B

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	100	G6_16635.D	1.	RINSE		25 Feb 2015 11:18
2	1	G6_16636.D	1.	DFTPP		25 Feb 2015 11:44
3	2	G6_16637.D	1.	ICIS HSL		25 Feb 2015 11:53
4	3	G6_16638.D	1.	STD004 HSL		25 Feb 2015 12:19
5	4	G6_16639.D	1.	STD010 HSL		25 Feb 2015 12:46
6	5	G6_16640.D	1.	STD020 HSL		25 Feb 2015 13:12
7	6	G6_16641.D	1.	STD050 HSL		25 Feb 2015 13:39
8	7	G6_16642.D	1.	STD120 HSL		25 Feb 2015 14:06
9	8	G6_16643.D	1.	STD160 HSL		25 Feb 2015 14:32
10	9	G6_16644.D	1.	STD200 HSL		25 Feb 2015 14:59
11	10	G6_16645.D	1.	ICV HSL 1		25 Feb 2015 15:26
12	11	G6_16646.D	1.	ICV HSL 2		25 Feb 2015 15:52
13	12	G6_16647.D	1.	ICV FAM		25 Feb 2015 16:19
14	13	G6_16648.D	1.	8270Surrogate_00077		25 Feb 2015 16:45
15	14	G6_16649.D	1.	LB3280-265126_1-B		25 Feb 2015 17:07
16	15	G6_16650.D	1.	LCS280-265126_2-B		25 Feb 2015 17:30
17	16	G6_16651.D	1.	LB280-265170_1-D		25 Feb 2015 17:52
18	17	G6_16652.D	1.	LCS280-265170_2-D		25 Feb 2015 18:14
19	18	G6_16653.D	1.	250-24545-B-2-B		25 Feb 2015 18:36
20	19	G6_16654.D	1.	250-24545-B-2-CMS		25 Feb 2015 18:58
21	20	G6_16655.D	1.	250-24545-B-2-DMSD		25 Feb 2015 19:20
22	21	G6_16656.D	1.	280-65511-D-1-H		25 Feb 2015 19:43
23	22	G6_16657.D	1.	280-65511-D-1-IMS		25 Feb 2015 20:05
24	23	G6_16658.D	1.	280-65511-D-1-JMSD		25 Feb 2015 20:27
25	98	G6_16659.D	1.	RINSE		25 Feb 2015 20:49
26	99	G6_16660.D	1.	RINSE		25 Feb 2015 21:16
27	100	G6_16661.D	1.	RINSE		25 Feb 2015 21:42
28	98	G6_16662.D	1.	RINSE		25 Feb 2015 22:09
29	99	G6_16663.D	1.	RINSE		25 Feb 2015 22:36
30	100	G6 16664.D	1.	RINSE		25 Feb 2015 23:02

Good

02.26.15

Dilution Solvent Lot #: MeOH Pipette ID: SV-12/SV-23 Method(s) Performed: 8270C

Daily Maintenance: Check box if maintenance was performed. No mark indicates the item was not changed.

Changed Septum [2] Changed Liner [2] Changed Seal [2] Changed Ferrule [2] Clipped Column [2]

New column

GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Batch Number: 272870 Batch Start Date: 04/15/15 18:50 Batch Analyst: Knauf, James R

Batch Method: 3520C Batch End Date: 04/17/15 19:05

Lab Sample ID	Client Sample ID	Method Chain	Basis	Initial pH	GrossWeight	TareWeight	InitialAmount	FinalAmount	ReceivedpH
MB 280-272870/1		3520C, 8270D		7			1000 mL	1 mL	7
LCS 280-272870/2		3520C, 8270D		7			1000 mL	1 mL	7
280-67711-C-1	TMW43042015	3520C, 8270D	T	7	1425.4 g	500.4 g	925 mL	1 mL	7
280-67711-B-1 MS	TMW43042015	3520C, 8270D	T	7	1386.9 g	501.9 g	885 mL	1 mL	7
280-67711-L-1 MSD	TMW43042015	3520C, 8270D	T	7	1403.2 g	499.9 g	903.3 mL	1 mL	7
280-67711-B-2	DTW43042015	3520C, 8270D	T	7	1530.7 g	501.7 g	1029 mL	1 mL	7
280-67711-C-4	TMW47042015	3520C, 8270D	T	7	1463.1 g	499.9 g	963.2 mL	1 mL	7

Lab Sample ID	Client Sample ID	Method Chain	Basis	FirstAdjustpH	SecondAdjustpH	8270_LCS_Main 00020	8270_LCS_Supp 00103	8270Surrogate 00078	AnalysisComment
MB 280-272870/1		3520C, 8270D		2	12			1 mL	same as batch 273368
LCS 280-272870/2		3520C, 8270D		2	12	1 mL	1 mL	1 mL	same as batch 273368
280-67711-C-1	TMW43042015	3520C, 8270D	T	2	12			1 mL	
280-67711-B-1 MS	TMW43042015	3520C, 8270D	T	2	12	1 mL	1 mL	1 mL	
280-67711-L-1 MSD	TMW43042015	3520C, 8270D	T	2	12	1 mL	1 mL	1 mL	
280-67711-B-2	DTW43042015	3520C, 8270D	T	2	12			1 mL	
280-67711-C-4	TMW47042015	3520C, 8270D	T	2	12			1 mL	

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

GC/MS SEMI VOA BATCH WORKSHEET

Lab Name: TestAmerica Denver Job No.: 280-67711-2

SDG No.:

Batch Number: 272870 Batch Start Date: 04/15/15 18:50 Batch Analyst: Knauf, James R

Batch Method: 3520C Batch End Date: 04/17/15 19:05

Batch Notes	
Acid used for pH adjustment	1:1 H2S04
Acid used for pH adjust Lot #	1:1 H2S04_00040
Balance ID	24350888
Base used for pH adjustment	10N_NaOH
Base used for pH adjust Lot #	10N_NaOH_00061
Batch Comment	DV OP 0008/0007 N. Elga Water NaCl: 145475
Person's name who did the concentration	EJ/ME
Time the first extraction ended 24hr	4.16.15@1315
Time the first extraction started 24 hr	4.15.15@1910
Na2SO4 Lot Number	145224_00001
Oven, Bath or Block Temperature 1	see above Celsius
Prep Solvent Lot #	MeCl2_Cycl_00211
Prep Solvent Name	MeCl2
Prep Solvent Volume Used	300 mL
Person's name who did the prep	James Knauf Pipette: K
Person's name who witnessed reagent drop	Reviewer:AH
Time the second extraction ended 24hr	4.17.15@0805
Time the second extraction started 24hr	4.16.15@1400
Sufficient volume for MS/MSD?	yes
Uncorrected Temperature	84 Celsius
Water Bath ID	A/B
Water Bath Temperature	84 Celsius

Basis	Basis Description
T	Total/NA

The pound sign (#) in the amount added field denotes that the reagent was used undiluted. All calculations are performed using the stated concentration for this reagent.

Shipping and Receiving Documents

TestAmerica Denver
4955 Yarrow Street
Arvada, CO 80002
phone 303-736-0100 fax 303-431-7171

280-67711 Chain of Custody

Chain of Custody Record

0.8, 1.2, 5.4, 0.7
1.0 + 0.2 IR-5
4/11/15 CO

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica Laboratories, Inc.

Client Contact	Project Manager: John Nance	Site Contact: John Nance 505 321 7260	Date: 4/10/15	COC No: FWDAAPR15-071
John Nance	Tel: 505 835 7660	Lab Contact: Michelle Johnston	Carrier: Federal Express	1 of 1 COCs
6700 Jefferson Street NE Suite C3	Analysis Turnaround Time			Job No.
Albuquerque, NM 87109	Calendar (C) or Work Days (W) W			
505 835 7660 PHONE	TAT if different from Below 15			SDG No.
	☐ 2 weeks			
Project Name: Fort Wingate	☐ 1 week			
Site: Fort Wingate, New Mexico	☐ 2 days			
PO#	☐ 1 day			

Sample Identification	Sample Date	Sample Time	Sample Type	Matrix	# of Cont.	Filtered Sample	Explosives 8330B	Nitrate Nitrite 9056	Perchlorate 6860	TAL Metal6010C/6020A/7470A Total	TAL Metal6010C/6020A/7470A Diss.	Organochlorine Pesticides 8081A	Volatile Organic Compounds 8260C	Semivolatile Organics 8270D	ORO 8015C C24-C36	TPH-DRO 8015C C10-C28	TPH-GRO 8015C C6-C10	TCLP VOCs 1311/8260B	TCLP SVOCs 1311/8270D	Metals 6010C/6020A dissolved	TCLP Pesticides 1311/8081A	TCLP Herbicides 1311/8151A	pH 9045C	Ignitibility 1010	Sample Specific Notes:
TMW43042015	4/10/2015	0825	grab	GW	6	Y		1		1	1		3												
TMW43042015MS	4/10/2015	0825	grab	GW	6	Y		1		1	1		3												
TMW43042015MSD	4/10/2015	0825	grab	GW	6	Y		1		1	1		3												
DTW43042015	4/10/2015	0825	grab	GW	6	Y		1		1	1		3												
TB -08-042015	4/10/2015	0800	grab	GW	2	N							2												
TMW47042015	4/10/2015	0930	grab	GW	3	N							3												

Preservation Used: 1= Ice, 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other _____

Possible Hazard Identification
☐ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

Sample Disposal
☐ Return To Client ☑ Disposal By Lab ☑ Archive For 60 days after invoice

Special Instructions/QC Requirements & Comments:

Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Todd Arrowood	Sundance	4/10/15 1500		TAD	4/11/15 930
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:

TestAmerica Denver
4955 Yarrow Street
Arvada, CO 80002
phone 303-736-0100 fax 303-431-7171

Chain of Custody Record

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica Laboratories, Inc.

Client Contact	Project Manager: John Nance	Site Contact: John Nance 505 321 7260	Date: 4/10/15	COC No: FWDAAPR15-069
John Nance	Tel: 505 835 7660	Lab Contact: Michelle Johnston	Carrier: Federal Express	1 of 1 COCs
6700 Jefferson Street NE Suite C3	Analysis Turnaround Time			Job No.
Albuquerque, NM 87109	Calendar (C) or Work Days (W) W			
505 835 7660 PHONE	TAT if different from Below 15			SDG No.
	☐ 2 weeks			
Project Name: Fort Wingate	☐ 1 week			
Site: Fort Wingate, New Mexico	☐ 2 days			
PO#	☐ 1 day			

Sample Identification	Sample Date	Sample Time	Sample Type	Matrix	# of Cont.	Filtered Sample	Explosives 8330B	Nitrate 9056	Perchlorate 6860	TAL Metal6010C/6020A/7470A Total	TAL Metal6010C/6020A/7470A Diss.	Organochlorine Pesticides 8081A	Volatile Organic Compounds 8260C	Semivolatile Organics 8270D	ORO 8015C C24-C36	TPH-DRO 8015C C10-C28	TPH-GRO 8015C C6-C10	TCLP VOCs 1311/8260B	TCLP SVOCs 1311/8270D	Metals 6010C/6020A dissolved	TCLP Pesticides 1311/8081A	TCLP Herbicides 1311/8151A	pH 9045C	Ignitibility 1010	Sample Specific Notes:
TMW43042015	4/10/2015	0825	grab	GW	2	N						2													
TMW43042015MS	4/10/2015	0825	grab	GW	2	N						2													
TMW43042015MSD	4/10/2015	0825	grab	GW	2	N						2													
DMW43042015	4/10/2015	0825	grab	GW	2	N						2													

Page 364 of 370

Preservation Used: 1= Ice, 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other _____

Possible Hazard Identification
☐ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

Sample Disposal
☐ Return To Client ☑ Disposal By Lab ☑ Archive For 60 days after invoice

Special Instructions/QC Requirements & Comments:

Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Todd Arrowood	Sundance	4/10/15 1500		TAD	4/11/15 930
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:

TestAmerica Denver
4955 Yarrow Street
Arvada, CO 80002
phone 303-736-0100 fax 303-431-7171

Chain of Custody Record

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
TestAmerica Laboratories, Inc.

Client Contact	Project Manager: John Nance	Site Contact: John Nance 505 321 7260	Date: 4/10/15	COC No: FWDAAPR15-070
John Nance	Tel: 505 835 7660	Lab Contact: Michelle Johnston	Carrier: Federal Express	1 of 1 COCs
6700 Jefferson Street NE Suite C3	Analysis Turnaround Time			Job No.
Albuquerque, NM 87109	Calendar (C) or Work Days (W) W			
505 835 7660 PHONE	TAT if different from Below 15			SDG No.
	☐ 2 weeks			
Project Name: Fort Wingate	☐ 1 week			
Site: Fort Wingate, New Mexico	☐ 2 days			
PO #	☐ 1 day			

Sample Identification	Sample Date	Sample Time	Sample Type	Matrix	# of Cont.	Filtered Sample	Explosives 8330B	Nitrate Nitrite 9056	Perchlorate 6860	TAL Metal6010C/6020A/7470A Total	TAL Metal6010C/6020A/7470A Diss.	Organochlorine Pesticides 8081A	Volatile Organic Compounds 8260C	Semivolatile Organics 8270D	ORO 8015C C24-C36	TPH-DRO 8015C C10-C28	TPH-GRO 8015C C6-C10	TCLP VOCs 1311/8260B	TCLP SVOCs 1311/8270D	Metals 6010C/6020A dissolved	TCLP Pesticides 1311/8081A	TCLP Herbicides 1311/8151A	pH 9045C	Ignitibility 1010	Sample Specific Notes:
TMW43042015	4/10/2015	0825	grab	GW	3	Y	2		1																
TMW43042015MS	4/10/2015	0825	grab	GW	3	Y	2		1																
TMW43042015MSD	4/10/2015	0825	grab	GW	3	Y	2		1																
DUP43042015	4/10/2015	0825	grab	GW	3	Y	2		1																

Preservation Used: 1= Ice, 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other _____

Possible Hazard Identification
☐ *Non-Hazard* ☐ *Flammable* ☐ *Skin Irritant* ☐ *Poison B* ☐ *Unknown*

Sample Disposal
☐ *Return To Client* ☑ *Disposal By Lab* ☑ *Archive For 60 days after invoice*

Special Instructions/QC Requirements & Comments:

Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Todd Arrowood	Sundance	4/10/15 1500		TAD	4/11/15 930
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:

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THE LEADER IN ENVIRONMENTAL TESTING
TestAmerica Laboratories, Inc.

Client Contact	Project Manager: John Nance	Site Contact: John Nance 505 321 7260	Date: 4/10/15	COC No: FWDAAPR15-068
John Nance	Tel: 505 835 7660	Lab Contact: Michelle Johnston	Carrier: Federal Express	1 of 1 COCs
6700 Jefferson Street NE Suite C3	Analysis Turnaround Time			Job No.
Albuquerque, NM 87109	Calendar (C) or Work Days (W) W			
505 835 7660 PHONE	TAT if different from Below 15			SDG No.
	☐ 2 weeks			
Project Name: Fort Wingate	☐ 1 week			
Site: Fort Wingate, New Mexico	☐ 2 days			
PO#	☐ 1 day			

Sample Identification	Sample Date	Sample Time	Sample Type	Matrix	# of Cont.	Filtered Sample	Explosives 8330B	Nitrate Nitrite 9056	Perchlorate 6860	TAL Metal6010C/6020A/7470A Total	TAL Metal6010C/6020A/7470A Diss.	Organochlorine Pesticides 8081A	Volatile Organic Compounds 8260C	Semivolatile Organics 8270D	ORO 8015C C24-C36	TPH-DRO 8015C C10-C28	TPH-GRO 8015C C6-C10	TCLP VOCs 1311/8260B	TCLP SVOCs 1311/8270D	Metals 6010C/6020A dissolved	TCLP Pesticides 1311/8081A	TCLP Herbicides 1311/8151A	pH 9045C	Ignitibility 1010	Sample Specific Notes:
TMW43042015	4/10/2015	0825	grab	GW	2	N								2											
TMW43042015MS	4/10/2015	0825	grab	GW	2	N								2											
TMW43042015MSD	4/10/2015	0825	grab	GW	2	N								2											
DUP43042015	4/10/2015	0825	grab	GW	2	N								2											

Page 366 of 370

Preservation Used: 1= Ice, 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other _____

Possible Hazard Identification
☐ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

Sample Disposal
☐ Return To Client ☑ Disposal By Lab ☑ Archive For 60 days after invoice

Special Instructions/QC Requirements & Comments:

Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Todd Arrowood	Sundance	4/10/15 1500		FHD	4/11/15 930
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:

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Chain of Custody Record

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TestAmerica Laboratories, Inc.

Client Contact	Project Manager: John Nance	Site Contact: John Nance 505 321 7260	Date: 4/10/15	COC No: FWDAAPR15-072
John Nance	Tel: 505 835 7660	Lab Contact: Michelle Johnston	Carrier: Federal Express	1 of 1 COCs
6700 Jefferson Street NE Suite C3	Analysis Turnaround Time			Job No.
Albuquerque, NM 87109	Calendar (C) or Work Days (W) W			
505 835 7660 PHONE	TAT if different from Below 15			SDG No.
	☐ 2 weeks			
Project Name: Fort Wingate	☐ 1 week			
Site: Fort Wingate, New Mexico	☐ 2 days			
P O #	☐ 1 day			

Sample Identification	Sample Date	Sample Time	Sample Type	Matrix	# of Cont.	Filtered Sample	Explosives 8330B	Nitrate 9056	Perchlorate 6860	TAL Metal6010C/6020A/7470A Total	TAL Metal6010C/6020A/7470A Diss.	Organochlorine Pesticides 8081A	Volatile Organic Compounds 8260C	Semivolatile Organics 8270D	ORO 8015C C24-C36	TPH-DRO 8015C C10-C28	TPH-GRO 8015C C6-C10	TCLP VOCs 1311/8260B	TCLP SVOCs 1311/8270D	Metals 6010C/6020A dissolved	TCLP Pesticides 1311/8081A	TCLP Herbicides 1311/8151A	pH 9045C	Ignitibility 1010	Sample Specific Notes:
TMW47042015	4/10/2015	0930	grab	GW	10	Y	2	1	1	1	1	2		2											

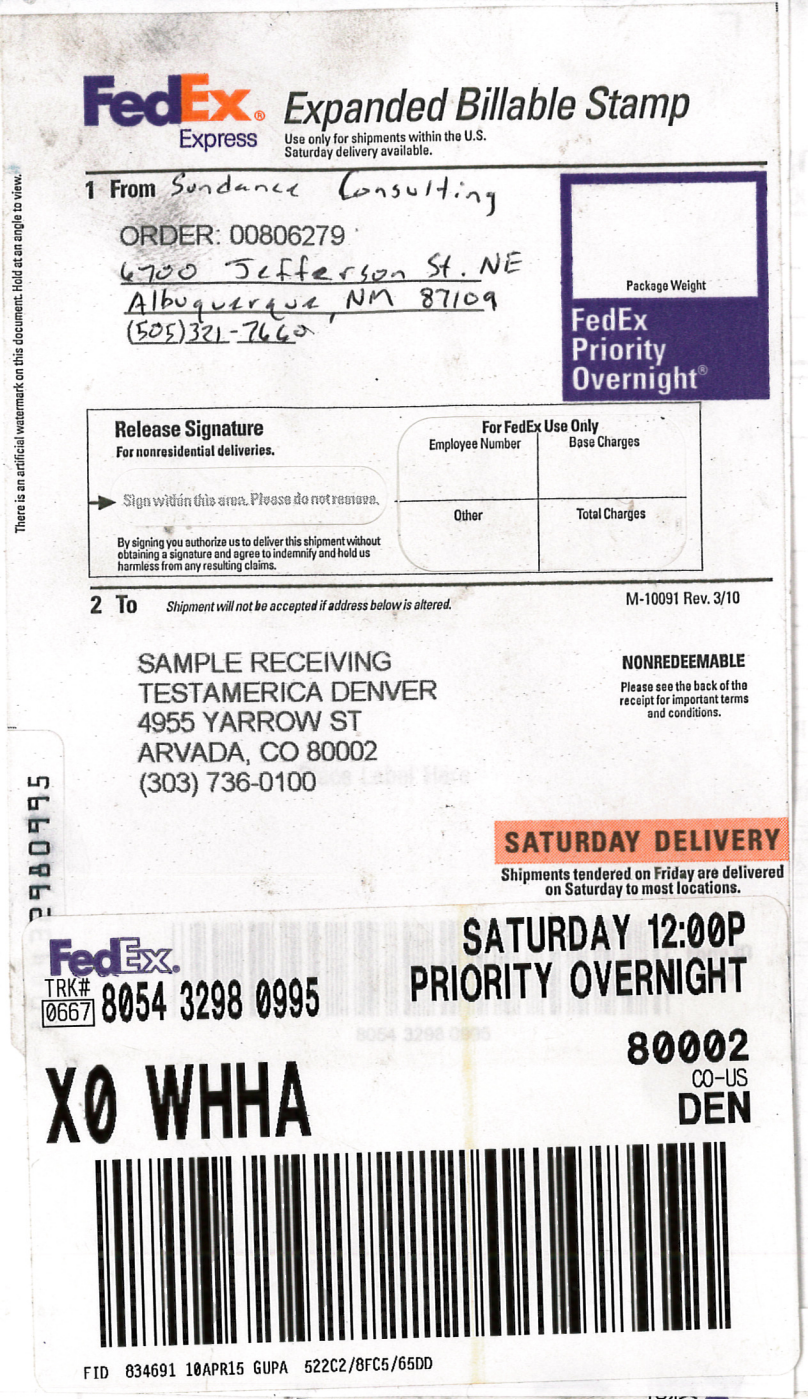
Preservation Used: 1= Ice, 2= HCl; 3= H2SO4; 4=HNO3; 5=NaOH; 6= Other _____

Possible Hazard Identification
☐ Non-Hazard ☐ Flammable ☐ Skin Irritant ☐ Poison B ☐ Unknown

Sample Disposal
☐ Return To Client ☑ Disposal By Lab ☑ Archive For 60 days after invoice

Special Instructions/QC Requirements & Comments:

Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Todd Arrowood	Sundance	4/10/15 1500		TAD	4/11/15 930
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:
Relinquished by:	Company:	Date/Time:	Received by:	Company:	Date/Time:



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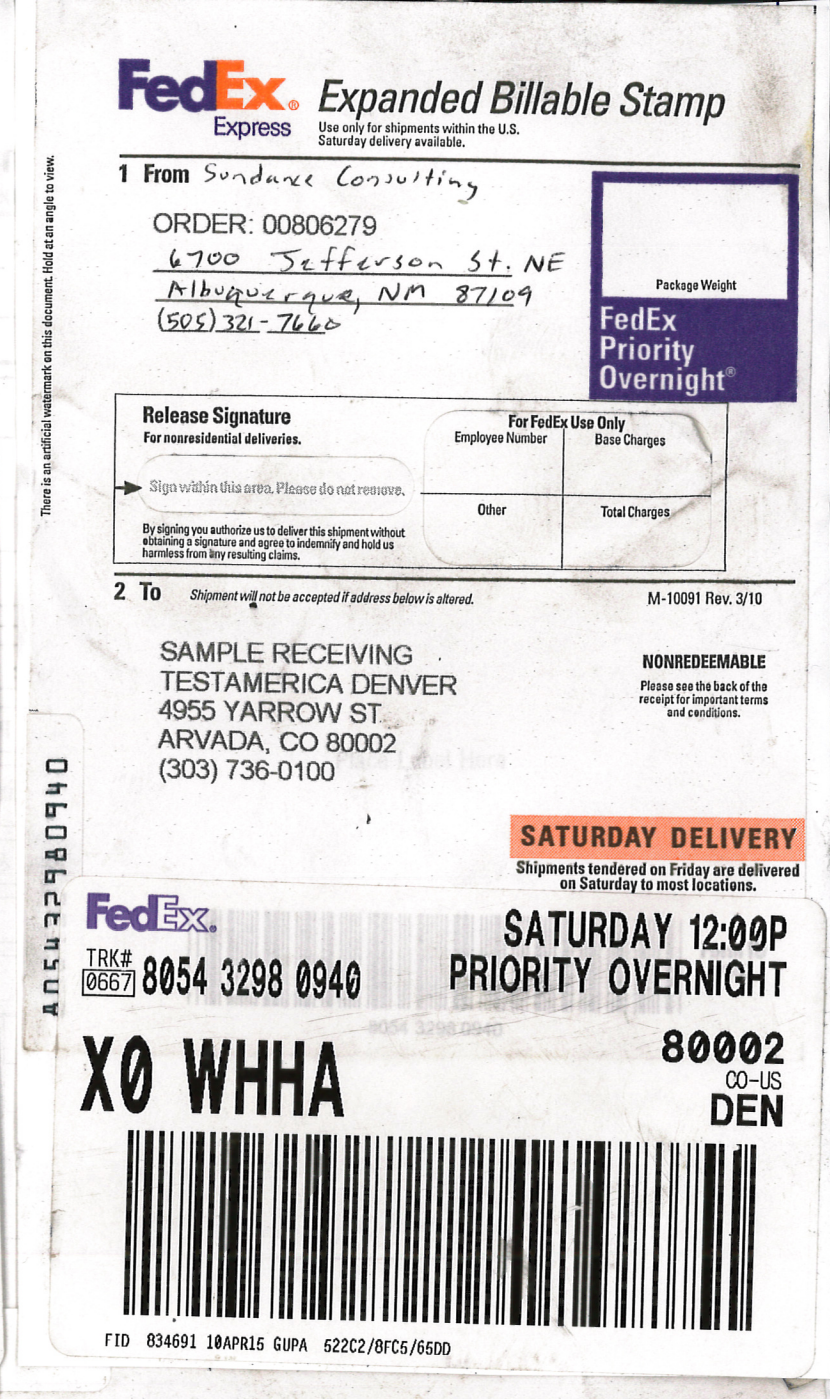

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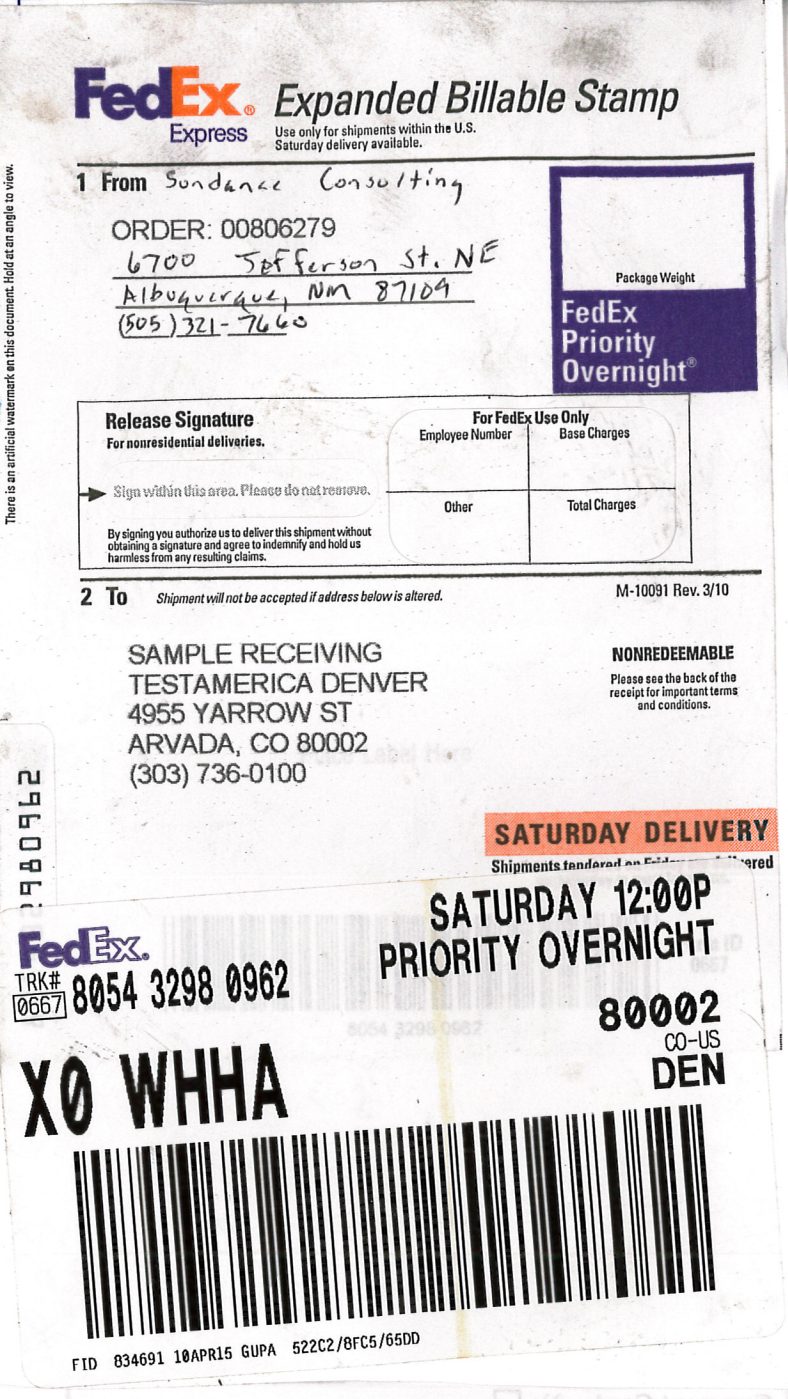
834691 10APR15 GUPA 522C2/8FC5/65DD



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Other | Total Charges

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SAMPLE RECEIVING
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ARVADA, CO 80002
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NONREDEEMABLE

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SATURDAY DELIVERY

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SATURDAY 12:00P
PRIORITY OVERNIGHT

80002
CO-US
DEN

X0 WHHA

FID 834691 10APR15 GUPA 522C2/8FC5/65DD

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1 From Sundance Consulting

ORDER: 00806279

6700 Jefferson St. NE
Albuquerque, NM 87109
(505) 321-7660

Package Weight

FedEx Priority Overnight®

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For nonresidential deliveries.

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Other | Total Charges

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CO-US
DEN

X0 WHHA

FID 834691 10APR15 GUPA 522C2/8FC5/65DD

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Login Sample Receipt Checklist

Client: Sundance Consulting, Inc

Job Number: 280-67711-2

Login Number: 67711

List Source: TestAmerica Denver

List Number: 1

Creator: Orfield, Tayler C

Question	Answer	Comment
Radioactivity wasn't checked or is </= background as measured by a survey meter.	N/A	
The cooler's custody seal, if present, is intact.	True	
Sample custody seals, if present, are intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
Is the Field Sampler's name present on COC?	False	Not requested on COC.
There are no discrepancies between the containers received and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
Sample Preservation Verified.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	
Residual Chlorine Checked.	N/A	