Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\

Data File : BG051474.D

Acq On : 11 Dec 2021 3:52

Operator : CG/JU Sample : M5009-02

Misc

ALS Vial : 23 Sample Multiplier: 1

Quant Time: Dec 11 04:29:54 2021

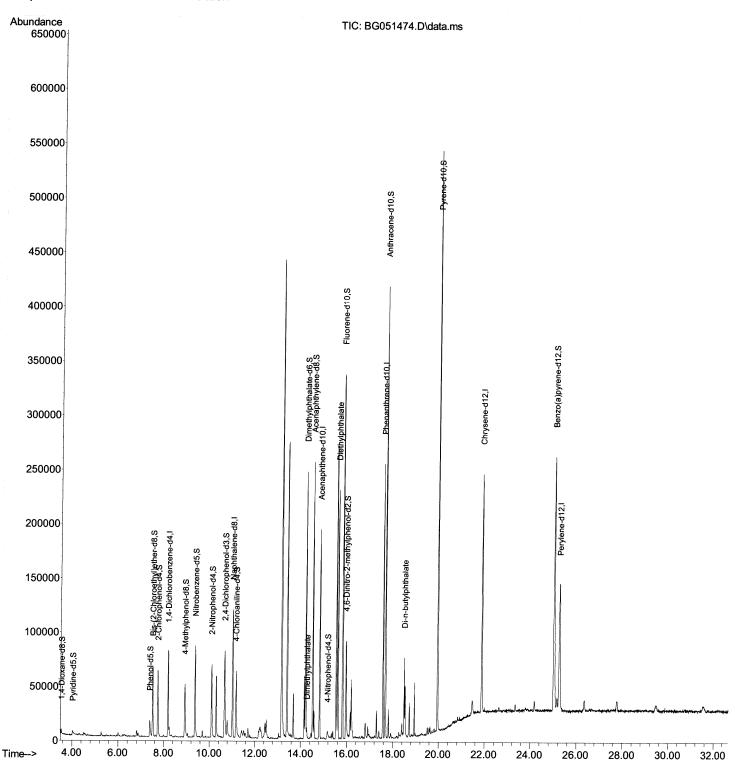
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/13/2021 Supervised By :Yogesh Patel 12/15/2021



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\

Data File : BG051474.D

Acq On : 11 Dec 2021 3:52

Operator : CG/JU Sample : M5009-02

Misc :

ALS Vial : 23 Sample Multiplier: 1

Quant Time: Dec 11 04:29:54 2021

 $\label{lem:quant_method} {\tt Quant\ Methods\SFAM-EPA-BG120821.M}$

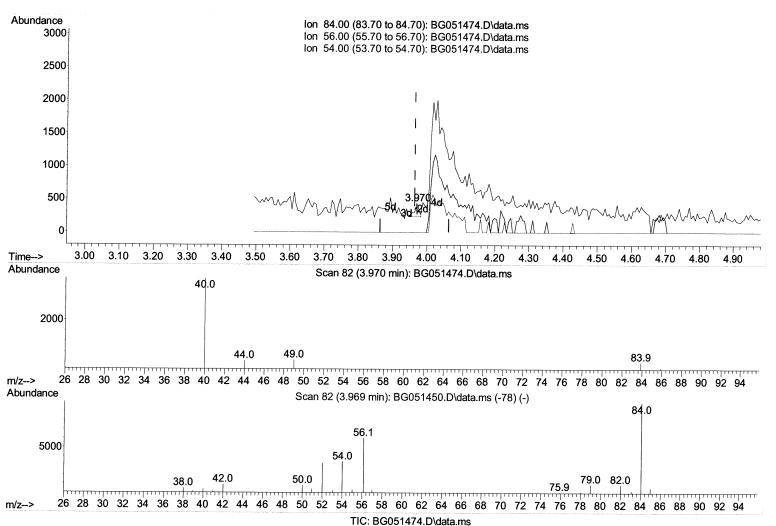
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration



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(4) Pyridine-d5 (S)

3.970min (+ 0.005) 0.09 ng/ul

response	185			
Ion	Ехр%	Act%		
84.00	100.00	100.00		
56.00	68.00	0.00#		
54.00	31.50	0.00#		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\

Data File : BG051474.D

Acq On : 11 Dec 2021 3:52

Operator : CG/JU Sample : M5009-02

Misc

ALS Vial : 23 Sample Multiplier: 1

Quant Time: Dec 11 04:29:54 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

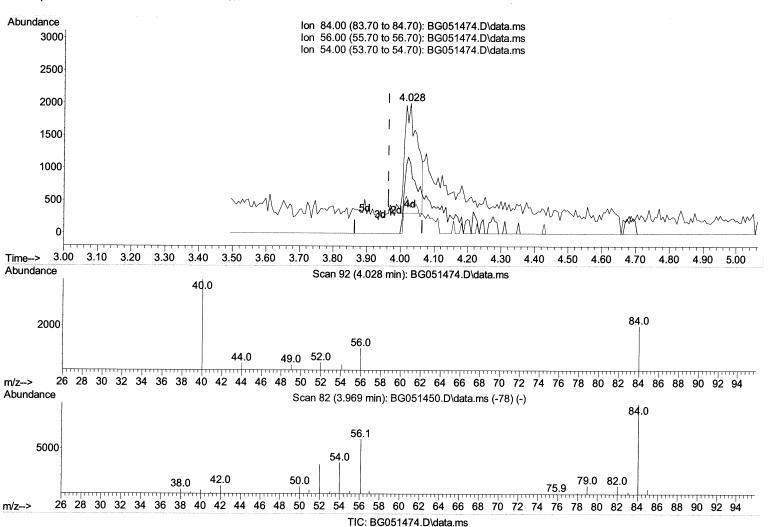
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

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(4) Pyridine-d5 (S)

4.028min (+ 0.064) 2.19 ng/ul m \2 //6 \3 (J)

response	4454	
Ion	Ежр%	Act%
84.00	100.00	100.00
56.00	68.00	53.76#
54.00	31.50	20.03#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\

Data File : BG051474.D

Acq On : 11 Dec 2021 3:52

: CG/JU **Operator** Sample : M5009-02

Misc

ALS Vial : 23 Sample Multiplier: 1

Quant Time: Dec 11 04:29:54 2021

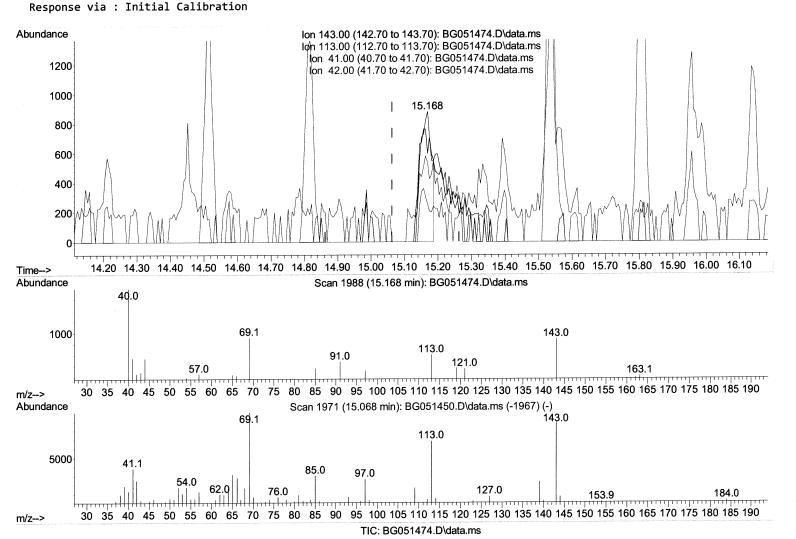
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 03:21:41 2021 Instrument: BNA_G ClientSampleId:

A0020

Manual Integrations APPROVED

Reviewed By: Jagrut Upadhyay 12/13/2021 Supervised By :Yogesh Patel 12/15/2021



(54) 4-Nitrophenol-d4 (S)

15.168min (+ 0.105) 2.53 ng/ul

response	2031			
Ion	Ехр%	Act*		
143.00	100.00	100.00		
113.00	80.30	67.80		
41.00	44.40	60.88#		
42.00	29.70	28.80		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\

Data File : BG051474.D

Acq On : 11 Dec 2021 3:52

Operator : CG/JU Sample : M5009-02

Misc :

ALS Vial : 23 Sample Multiplier: 1

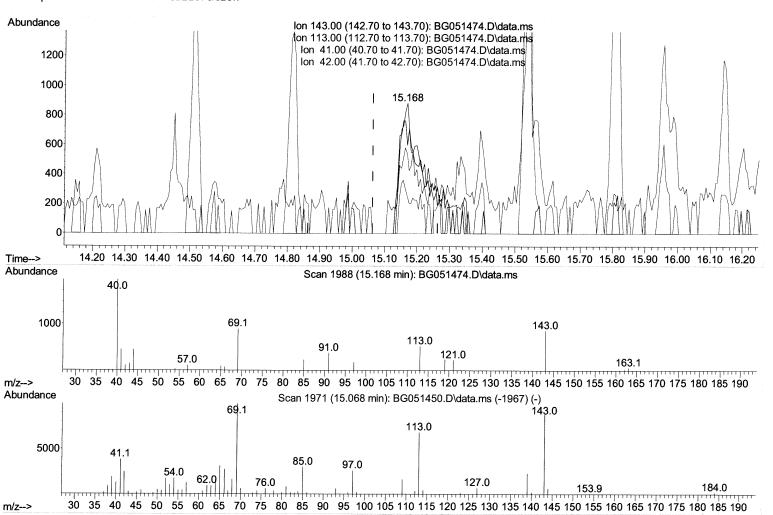
Quant Time: Dec 11 04:29:54 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId :

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/13/2021 Supervised By :Yogesh Patel 12/15/2021



TIC: BG051474.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.168min (+ 0.105) 4.60 ng/ul m \2//6/2/JU

response	3684		
Ion	Ежр%	Act%	
143.00	100.00	100.00	
113.00	80.30	67.80	
41.00	44.40	60.88#	
42.00	29.70	28.80	

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\

Data File : BG051474.D

Acq On : 11 Dec 2021 3:52

Operator : CG/JU Sample : M5009-02

Misc

ALS Vial : 23 Sample Multiplier: 1

Quant Time: Dec 11 04:29:54 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : A0020

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/13/2021 Supervised By :Yogesh Patel 12/15/2021

Compound	R.T.	QIon	Response C	onc Uni	ts Dev(M	lin)
Internal Standards				20.000	(]	0.00
 1,4-Dichlorobenzene-d4 	8.188	152	23261	20.000	-	0.00 0.00
20) Naphthalene-d8	11.008	136	101531	20.000	•	0.00
38) Acenaphthene-d10	14.816	164	68815	20.000	-	
64) Phenanthrene-d10	17.565		155750	20.000	•	0.00 0.00
79) Chrysene-d12	21.872		151880	20.000	•	
88) Perylene-d12	25.268	264	142009	20.000	ng/u1	0.00
System Monitoring Compounds					, ,	0.00
3) 1,4-Dioxane-d8	3.535		2963	4.183	ng/uL	0.00
4) Pyridine-d5	4.028					0.06 12/16/21/19
7) Phenol-d5	7.383		13166	5.560	_	0.03
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.501		35619	23.456	•	0.00
11) 2-Chlorophenol-d4	7.730		31916	18.944		0.00
15) 4-Methylphenol-d8	8.923		25912	13.929	•	0.01
21) Nitrobenzene-d5	9.369		22159	25.159	_	0.00
24) 2-Nitrophenol-d4	10.098		24046	24.127	•	0.00
28) 2,4-Dichlorophenol-d3	10.662		37786	23.305	_	0.01
31) 4-Chloroaniline-d4	11.167		44428	18.735	•	0.00
46) Dimethylphthalate-d6	14.210	166	162140	30.450	•	0.00
49) Acenaphthylene-d8	14.516	160	187425	27.792	•	0.00
54) 4-Nitrophenol-d4	15.168	143	3684m>		ng/ul>	0.11 11 (012)
60) Fluorene-d10	15.809	176	136943	28.890		0.00
65) 4,6-Dinitro-2-methylph	15.955	200	25230	27.260		0.00
73) Anthracene-d10	17.665	188		34.302		0.00
81) Pyrene-d10	19.945	212	309487	33.902		0.00
92) Benzo(a)pyrene-d12	25.033	264	261976	35.766	ng/ul	0.00
Target Compounds					Qva	
47) Dimethylphthalate	14.263	163	5827		ng/ul#	97
59) Diethylphthalate	15.609	149	127932		ng/ul	99
78) Di-n-butylphthalate	18.494	149	49023	4.890	ng/ul	100

^(#) = qualifier out of range (m) = manual integration (+) = signals summed