Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 42 Sample Multiplier: 1

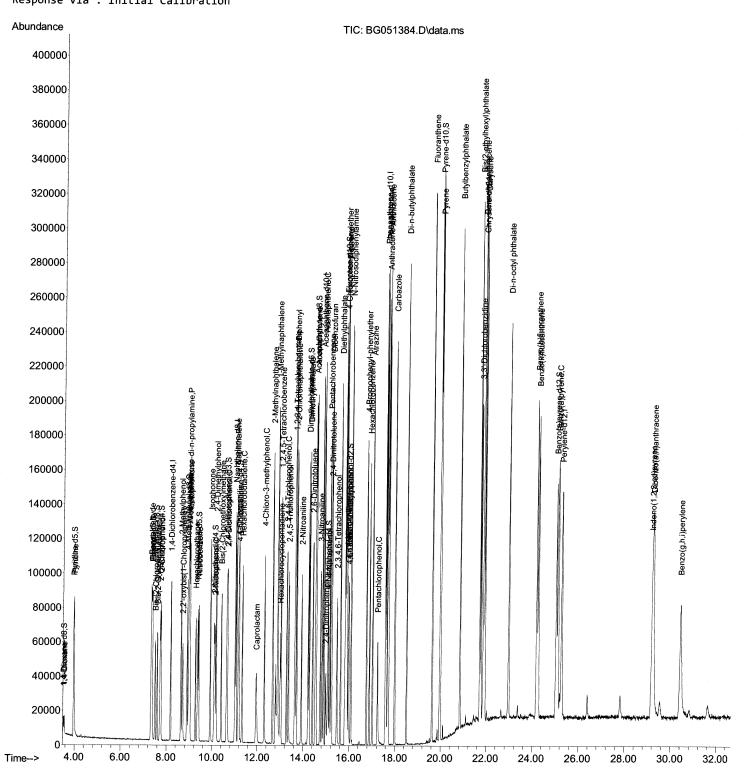
Quant Time: Dec 08 02:25:30 2021

 $\label{lem:quant_method} {\tt Quant_Methods\SFAM-EPA-BG112321.M}$

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument :
BNA_G
LabSampleId :
SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



SFAM-EPA-BG112321.M Wed Dec 08 03:39:03 2021

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

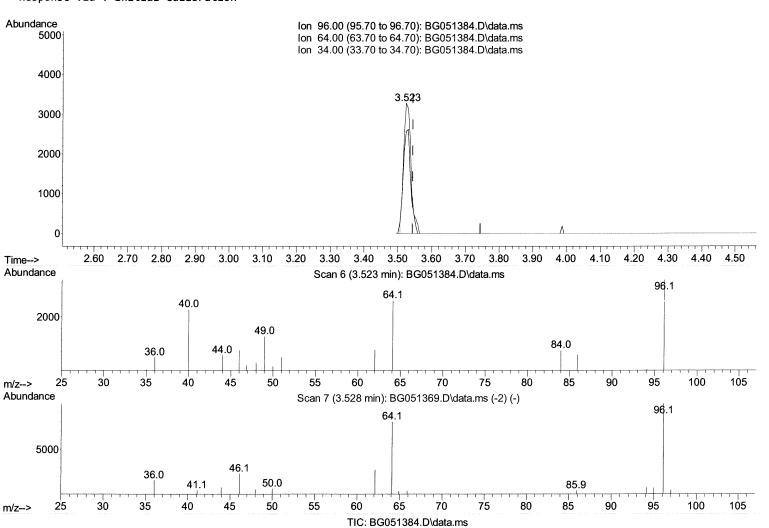
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/08/2021 Supervised By: mohammad ahmed 12/15/2021



(3) 1,4-Dioxane-d8 (S)

3.523min (-0.021) 6.89 ng/uL

response	5061			
Ion	Ехр%	Act*		
96.00	100.00	100.00		
64.00	77.60	78.38		
34.00	0.00	0.00		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

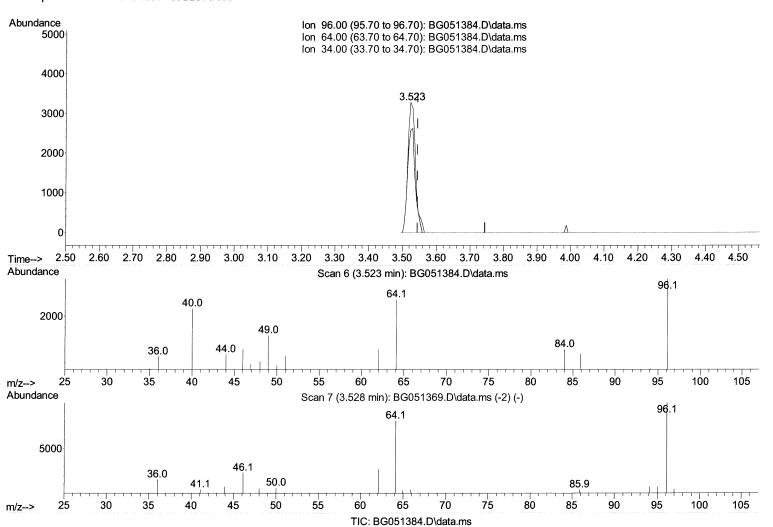
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



(3) 1,4-Dioxane-d8 (S)

3.523min (-0.021) 7.02 ng/uL m (2////2/J

response	5160	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	78.38
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

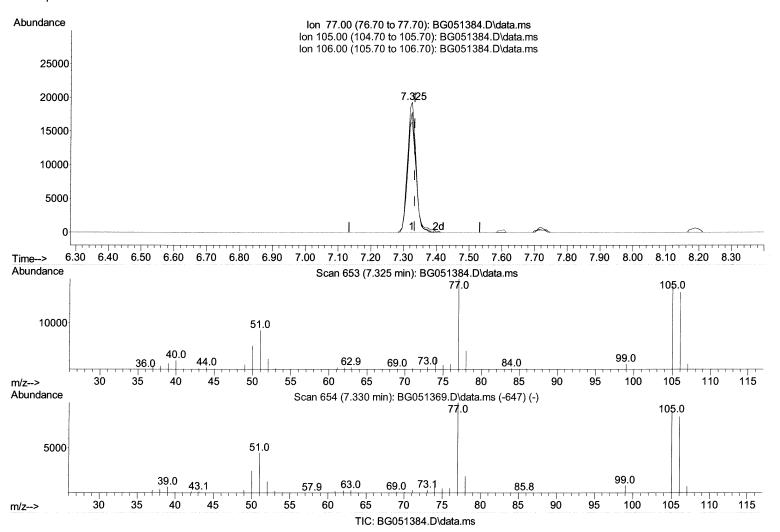
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/08/2021 Supervised By: mohammad ahmed 12/15/2021



(6) Benzaldehyde

7.325min (-0.009) 22.35 ng/ul

response	35910	
Ion	Ежр%	Act%
77.00	100.00	100.00
105.00	88.00	92.22
106.00	76.50	85.06
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

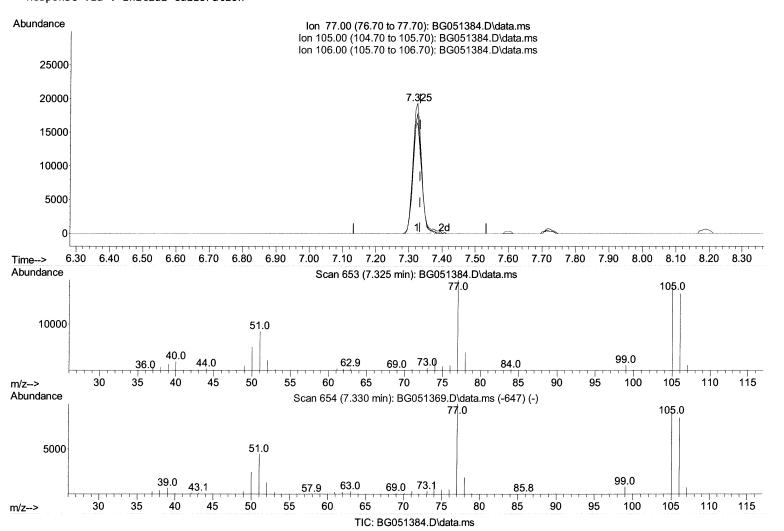
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



(6) Benzaldehyde

response	35128	
Ion	Ехр%	Act%
77.00	100.00	100.00
105.00	88.00	92.22
106.00	76.50	85.06
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

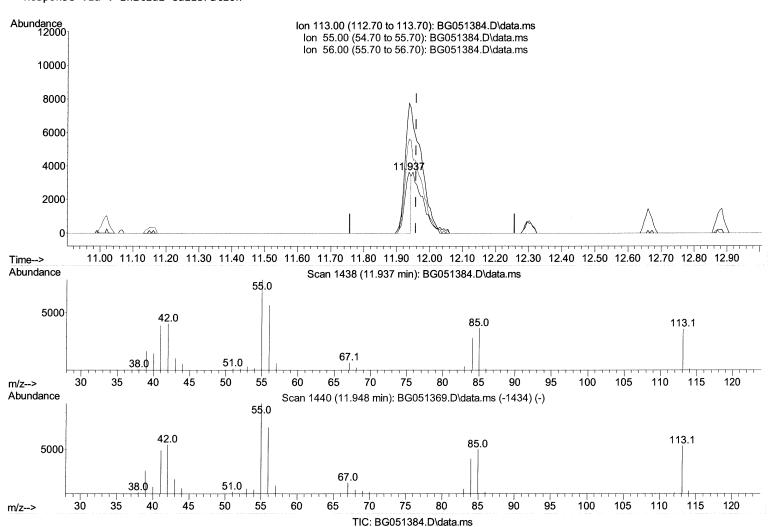
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/08/2021 Supervised By: mohammad ahmed 12/15/2021



(34) Caprolactam

11.937min (-0.021) 6.92 ng/ul

response	5000	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	183.80	214.03
56.00	136.50	154.86
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

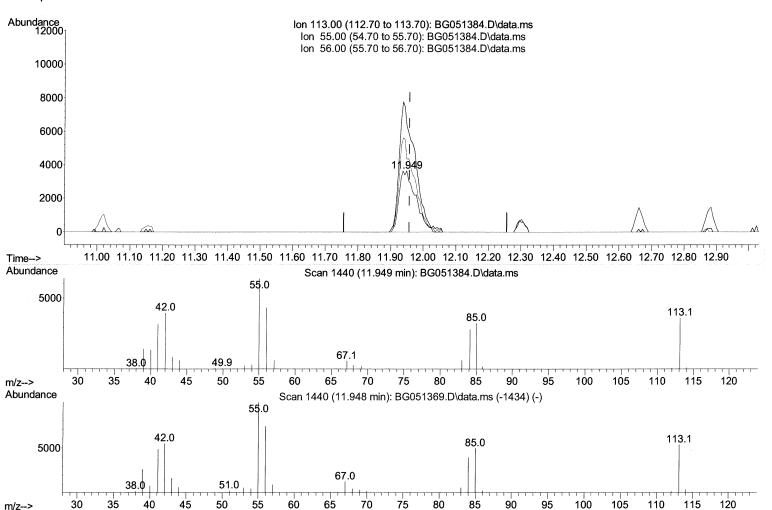
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



TIC: BG051384.D\data.ms

(34) Caprolactam

11.949min (-0.009) 17.87 ng/ul m \ 2 | | (2) U

response	12916	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	183.80	174.17
56.00	136.50	119.41
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

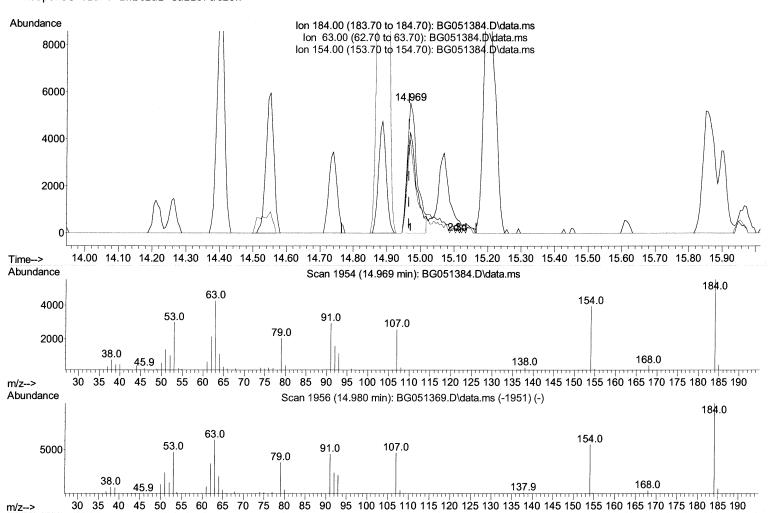
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



TIC: BG051384.D\data.ms

(53) 2,4-Dinitrophenol

14.969min (+ 0.003) 15.97 ng/ul

response	11195	
Ion	Exp%	Act%
184.00	100.00	100.00
63.00	82.70	77.38
154.00	67.00	71.57
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

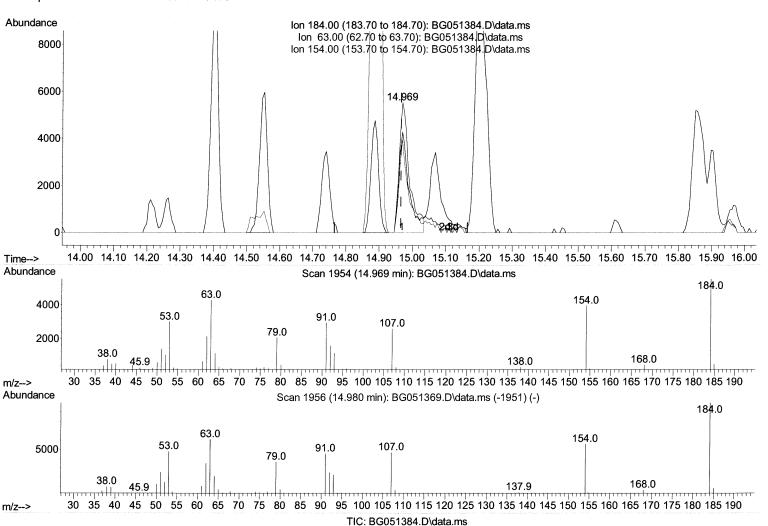
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



(53) 2,4-Dinitrophenol

14.969min (+ 0.003) 17.09 ng/ul m \J\/\()[J]

response	11977	
Ion	Ежр%	Act%
184.00	100.00	100.00
63.00	82.70	77.38
154.00	67.00	71.57
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

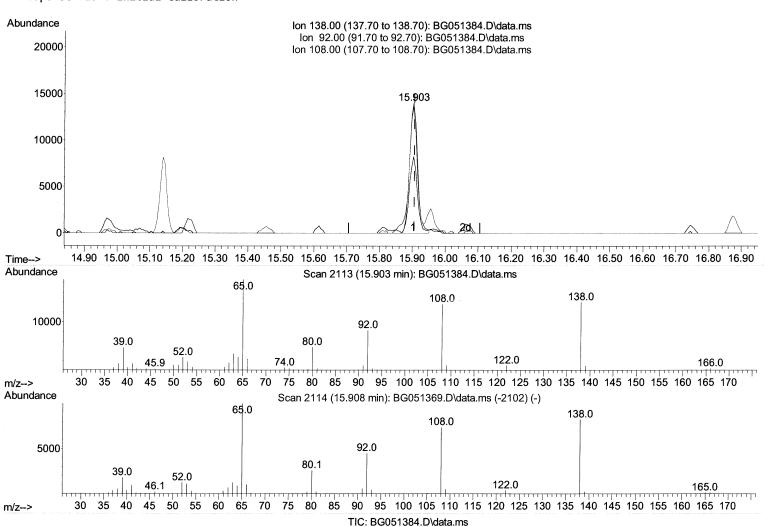
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



(63) 4-Nitroaniline

15.903min (-0.003) 21.08 ng/ul

response	25708	
Ion	Exp%	Act%
138.00	100.00	100.00
92.00	61.60	58.76
108.00	90.70	97.24
0.00	0 00	0 00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc :

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

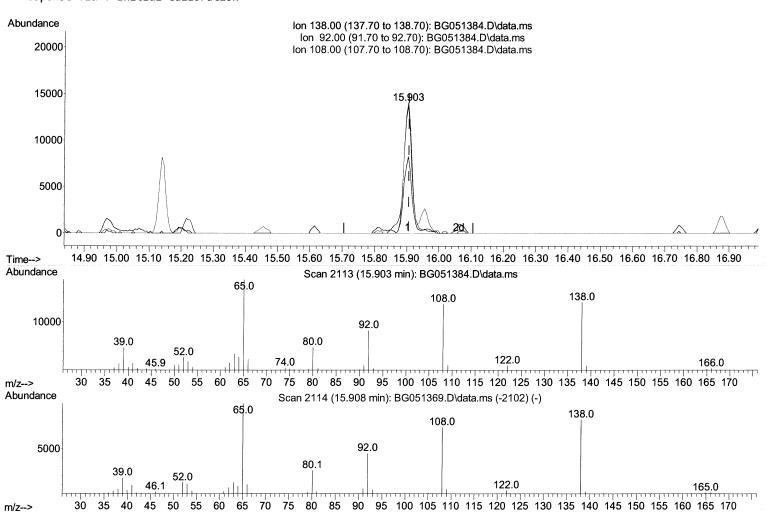
Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021



TIC: BG051384.D\data.ms

(63) 4-Nitroaniline

15.903min (-0.003) 21.48 ng/ul m (2////2/JU

response	26193	
Ion	Ехр%	Act%
138.00	100.00	100.00
92.00	61.60	58.76
108.00	90.70	97.24
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56 Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Instrument : BNA_G LabSampleId : SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021 Supervised By :mohammad ahmed 12/15/2021

	Compound	R.T.	QIon	Response	Conc Units Dev(Min)	
Inter	nal Standards					
1)	1,4-Dichlorobenzene-d4	8.188	152	25530	20.000 ng/ul -0.01	
	Naphthalene-d8	11.014	136	115593	20.000 ng/ul -0.01	
38)	Acenaphthene-d10	14.822	164	77501	20.000 ng/ul 0.00	
64)	Phenanthrene-d10	17.571	188	168699	20.000 ng/ul 0.00	
	Chrysene-d12	21.878	240	142995	20.000 ng/ul 0.00	
88)	Perylene-d12	25.280	264	143095	20.000 ng/ul 0.00	
Syste	m Monitoring Compounds				(01:	~ J
3)	1,4-Dioxane-d8	3.523	96	5160m _{>}	7.024 ng/uL> -0.02 12/16/21	٠, ۲
4)	Pyridine-d5	3.958	84	37313	17.308 ng/ul -0.02	
7)	Phenol-d5	7.354	99	47418	18.793 ng/ul 0.00	
9)	Bis-(2-Chloroethyl)eth	7.501	67	29763	18.781 ng/ul -0.01	
11)	2-Chlorophenol-d4	7.724	132	35013	19.270 ng/ul 0.00	
15) 4	4-Methylphenol-d8	8.905	113	37395	18.365 ng/ul 0.00	
21)	Nitrobenzene-d5	9.369	128	18731	19.196 ng/ul 0.00	
24)	2-Nitrophenol-d4	10.092	143	20531	18.652 ng/ul 0.00	
28)	2,4-Dichlorophenol-d3	10.644	165	35500	19.009 ng/ul 0.00	
31) 4	4-Chloroaniline-d4	11.155	131	50163	18.357 ng/ul 0.00	
46) [Dimethylphthalate-d6	14.217	166	108250	18.153 ng/ul 0.00	
49) /	Acenaphthylene-d8	14.522	160	143101	19.030 ng/ul 0.00	
54) 4	4-Nitrophenol-d4	15.051	143	14803	15.336 ng/ul 0.00	
	Fluorene-d10	15.815	176	100312	18.680 ng/ul 0.00	
65) 4	4,6-Dinitro-2-methylph	15.956	200	15468	14.859 ng/ul 0.00	
73) A	Anthracene-d10	17.671	188	154385	19.135 ng/ul 0.00	
81) F	Pyrene-d10	19.951	212	170247	19.677 ng/ul 0.00	
92) E	Benzo(a)pyrene-d12	25.039	264	141425	18.506 ng/ul 0.00	
Target	Compounds				Qvalue	
2) 1	l,4-Dioxane	3.558	88	5974	7.210 ng/uL 88	
5) F	Pyridine	3.976	79	40573	18.087 ng/ul 95	
6) E	Benzaldehyde	7.325	77	35128m >	21.861 ng/ul > (all 2 J	•
8) F	Phenol	7.378	94	48603	18.594 ng/ul 98	
10) E	Bis(2-Chloroethyl)ether	7.601	93	36779	18.598 ng/ul 97	
12) 2	2-Chlorophenol	7.754	128	35965	19.424 ng/ul 99	
	2-Methylphenol	8.635	108	36548	18.771 ng/ul 100	
14) 2	2,2'-oxybis(1-Chloropr	8.711	45	56432	19.775 ng/ul 96	
	Acetophenone	9.017	105	58743	18.652 ng/ul 100	
	I-Nitroso-di-n-propyla	8.987	70	34932	19.301 ng/ul 99	
18) 4	-Methylphenol	8.970	108	39341	18.896 ng/ul 97	
1 9) H	lexachloroethane	9.275	117	15372	19.656 ng/ul 95	
22) N	litrobenzene	9.410	77	50299	19.659 ng/ul 95	
	Sophorone	9.927	82	94931	19.097 ng/ul 99	
	-Nitrophenol	10.127	139	21322	18.702 ng/ul 95	
•	.,4-Dimethylphenol	10.174	107	44800	19.219 ng/ul 98	
	is(2-Chloroethoxy)met	10.403	93	50902	18.549 ng/ul 100	
	,4-Dichlorophenol	10.674	162	34996	19.036 ng/ul 96	
	laphthalene	11.067	128	120758	19.199 ng/ul 97	
	-Chloroaniline	11.185	127	49772	18.143 ng/ul 99	
	exachlorobutadiene	11.332	225	23598	18.610 ng/ul 98	
	aprolactam	11.949	113	12916m >	17.871 ng/ul > いより(は)しん	
35) 4	-Chloro-3-methylphenol	12.301	107	42691	19.331 ng/ul 96	

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051384.D

Acq On : 7 Dec 2021 15:56

Operator : CG/JU Sample : SSTDCCC020

Misc

ALS Vial : 42 Sample Multiplier: 1

Quant Time: Dec 08 02:25:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Instrument : BNA_G **LabSampleld :** SSTDCCC020

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/08/2021 Supervised By: mohammad ahmed 12/15/2021

Compound	R.T.	QIon	Response	Conc Ur	nits Dev(Min)
36) 2-Methylnaphthalene	12.660	142	80657	18.853	ng/ul	100
37) 1-Methylnaphthalene	12.877	142	83014	18.861	. ng/ul	100
39) 1,2,4,5-Tetrachloroben	13.024	216	46036	18.921	. ng/ul	97
40) Hexachlorocyclopentadiene	12.989	237	15772	16.038	ng/ul	97
41) 2,4,6-Trichlorophenol	13.271	196	28582	18.720	ng/ul	99
42) 2,4,5-Trichlorophenol	13.353	196	30306	18.954	ng/ul	97
43) 1,1'-Biphenyl	13.658	154	109311	18.884	ng/ul	96
44) 2-Chloronaphthalene	13.705	162	88289		ng/ul	99
45) 2-Nitroaniline	13.917	65	31708		ng/ul	94
47) Dimethylphthalate	14.264	163	108533		ng/ul	99
48) 2,6-Dinitrotoluene	14.405	165	23662	18.662	ng/ul	94
50) Acenaphthylene	14.551	152	139559	18.785	ng/ul	98
51) 3-Nitroaniline	14.739	138	25558		ng/ul	94
52) Acenaphthene	14.886	153	92027	18.783	ng/ul	94
53) 2,4-Dinitrophenol	14.969	184	11977m>		ng/ul>	12/1/10174
55) 4-Nitrophenol	15.068	109	15465	18.469		98
56) Dibenzofuran	15.221	168	132376	18.731		100
57) 2,4-Dinitrotoluene	15.198	165	34043		ng/ul#	94
58) 2,3,4,6-Tetrachlorophenol	15.456	232	20230	16.112		97
59) Diethylphthalate	15.615	149	116916	18.453	-	99
61) Fluorene	15.868	166	106286	18.776		99
62) 4-Chlorophenyl-phenyle	15.850	204	55150	18.078	-	1911/13127
63) 4-Nitroaniline	15.903	138	26193m>		ng/ul ➤	
66) 4,6-Dinitro-2-methylph	15.967	198	14461	14.404		99
67) N-Nitrosodiphenylamine	16.067	169	93332	19.325		99
68) 4-Bromophenyl-phenylether	16.749	248	34381	19.015	0.	93
69) Hexachlorobenzene	16.878	284	35037	19.004	_	98
70) Atrazine	17.007	200	38677	19.055		99
71) Pentachlorophenol72) Phenanthrene	17.237	266	15628	19.130		92
	17.618	178	180217	19.348		99
<pre>74) Anthracene 75) 1,2,3,4-Tetrachloroben</pre>	17.707	178 216	179360	19.389	-	98
76) Pentachlorobenzene	13.629 15.139	250	48807	19.835 19.314	•	96 100
77) Carbazole	17.983	167	44283 159526	19.646		100 100
78) Di-n-butylphthalate	18.500	149	205888	19.665	-	100
80) Fluoranthene	19.622	202	210842	19.840		97
82) Pyrene	19.980	202	207315	19.943		98
83) Butylbenzylphthalate	20.838	149	86723	20.067		100
84) 3,3'-Dichlorobenzidine	21.761	252	66729	20.043	-	98
85) Benzo(a)anthracene	21.855	228	186326	19.211	•	99
86) Bis(2-ethylhexyl)phtha	21.708	149	121565	19.548	-	98
87) Chrysene	21.925	228	176472	18.940		100
89) Di-n-octyl phthalate	22.971	149	204402	19.717		100
90) Benzo(b)fluoranthene	24.187	252	181346	18.779	_	99
91) Benzo(k)fluoranthene	24.258	252	167685	18.504		98
93) Benzo(a)pyrene	25.116	252	173446	18.826		99
94) Indeno(1,2,3-cd)pyrene	29.193	276	187624	18.199	•	97
95) Dibenzo(a,h)anthracene	29.246	278	159180	18.200		97
96) Benzo(g,h,i)perylene	30.427	276	152698	17.604		95

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