Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\

Data File : BG051188.D

Acq On : 23 Nov 2021 17:07

Operator : CG/JU Sample : SSTDICV020

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 24 06:08:17 2021

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

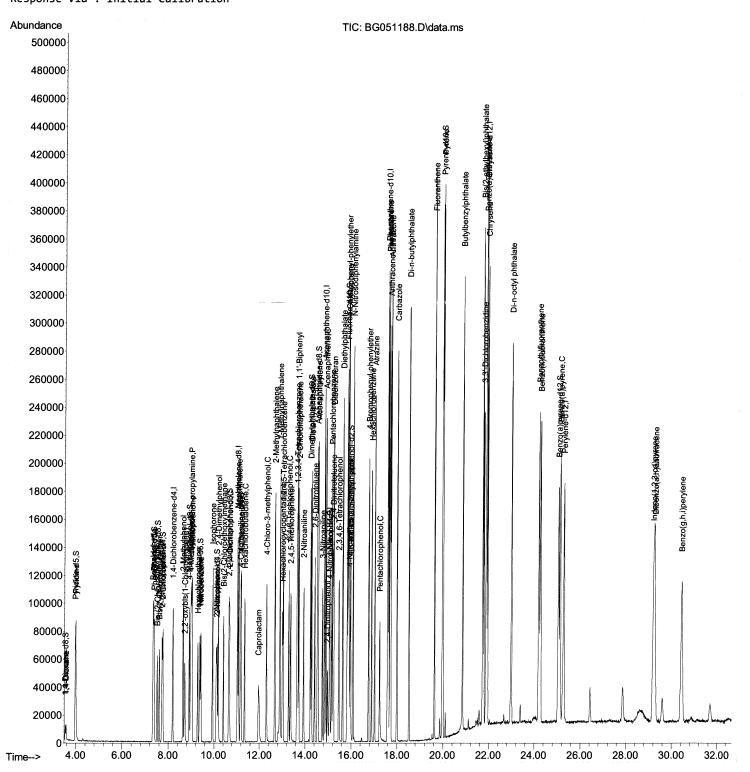
Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/30/2021



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\

Data File : BG051188.D

Acq On : 23 Nov 2021 17:07

Operator : CG/JU Sample : SSTDICV020

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 24 06:08:17 2021

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

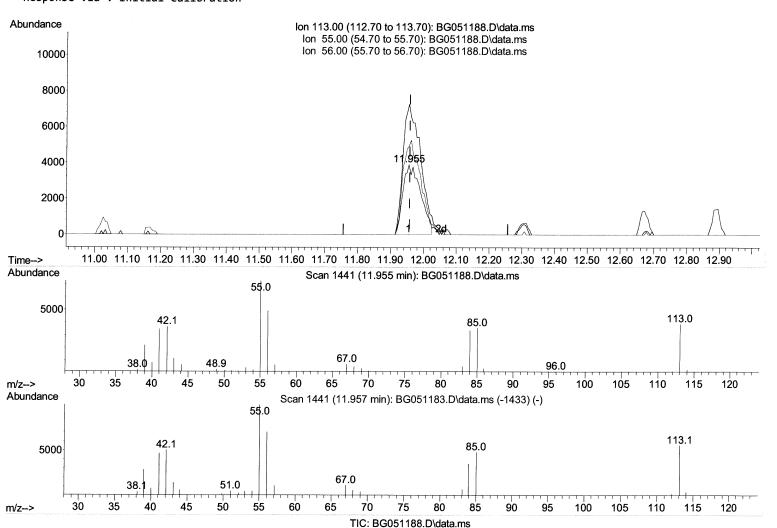
Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/30/2021



(34) Caprolactam

11.955min (-0.003) 18.75 ng/ul

response	14612	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	183.80	186.31
56.00	136.50	126.41
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\

Data File: BG051188.D

Acq On : 23 Nov 2021 17:07

Operator : CG/JU Sample : SSTDICV020

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 24 06:08:17 2021

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

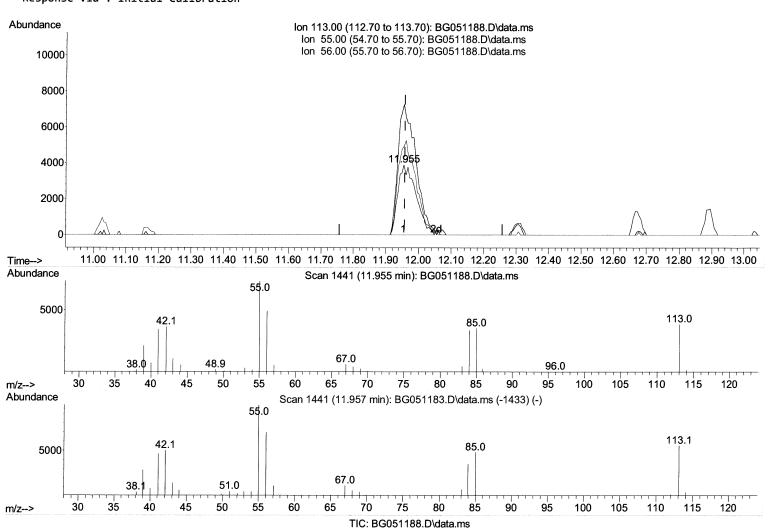
Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration



Manual Integrations APPROVED

Reviewed By: Jagrut Upadhyay 11/24/2021 Supervised By: mohammad ahmed 11/30/2021



(34) Caprolactam

11.955min (-0.003) 19.17 ng/ul m 1/29/21JU

response	14938	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	183.80	186.31
56.00	136.50	126.41
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG112321\

Data File: BG051188.D

Acq On : 23 Nov 2021 17:07

Operator : CG/JU Sample : SSTDICV020

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 24 06:08:17 2021

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

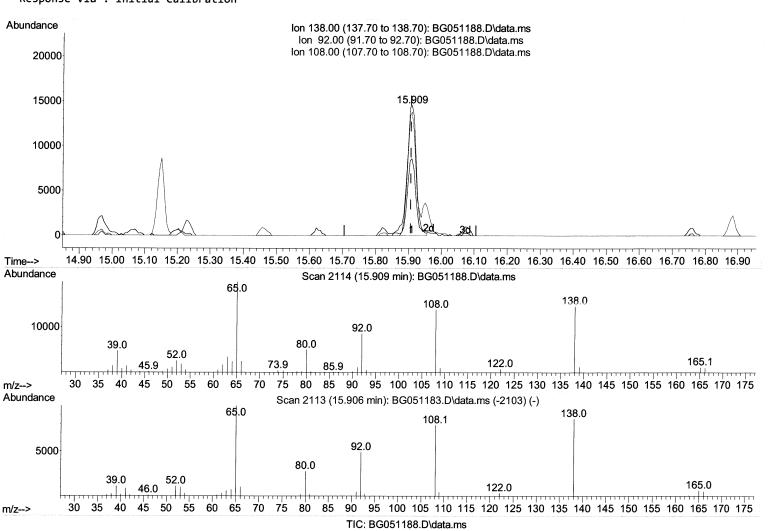
Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration



Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/30/2021



(63) 4-Nitroaniline

15.909min (+ 0.003) 20.64 ng/ul

response	28459				
Ion	Ежр%	Act%			
138.00	100.00	100.00			
92.00	61.60	59.15			
108.00	90.70	94.95			
0.00	0 00	0.00			

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG112321\

Data File : BG051188.D

Acq On : 23 Nov 2021 17:07

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 24 06:08:17 2021

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

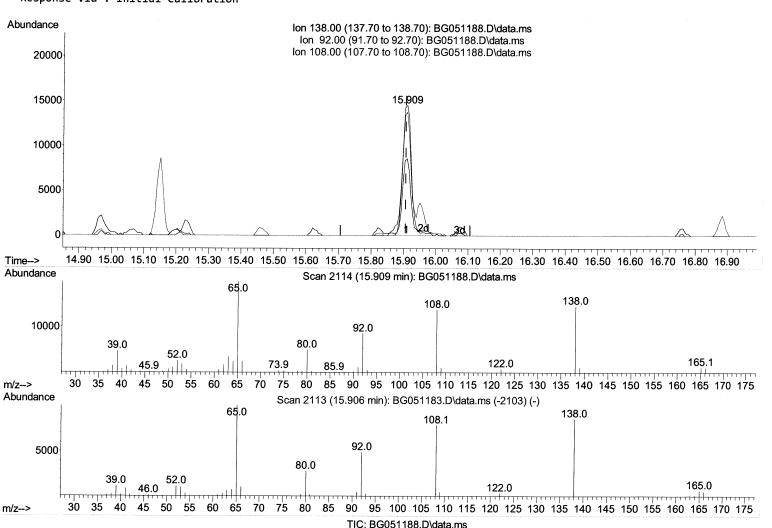
Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/24/2021 Supervised By: mohammad ahmed 11/30/2021



(63) 4-Nitroaniline

response	28917			
Ion	Ехр%	Act%		
138.00	100.00	100.00		
92.00	61.60	59.15		
108.00	90.70	94.95		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\

Data File : BG051188.D

Acq On : 23 Nov 2021 17:07

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 24 06:08:17 2021

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

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

Instrument : BNA_G ClientSampleId : SICV427

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/30/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Internal Standards					
 1,4-Dichlorobenzene-d4 	8.200		27208	20.000 ng/ul	0.00
20) Naphthalene-d8	11.026	136	124638	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.834	164	87608	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.583	188	206867	20.000 ng/ul	0.00
79) Chrysene-d12	21.884	240	188724	20.000 ng/ul	0.00
88) Perylene-d12	25.286	264	189891	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.535	96	5754	7.349 ng/uL	0.00
4) Pyridine-d5	3.970	84	41915	18.244 ng/ul	0.00
7) Phenol-d5	7.354	99	47152	17.535 ng/ul	0.00
Bis-(2-Chloroethyl)eth	7.513	67	30705	18.181 ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.730	132	34883	18.014 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.911	113	38708	17.838 ng/ul	0.00
21) Nitrobenzene-d5	9.375	128	19088	18.142 ng/ul	0.00
24) 2-Nitrophenol-d4	10.104	143	22091	18.613 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.650	165	36860	18.305 ng/ul	0.00
31) 4-Chloroaniline-d4	11.161	131	54448	18.479 ng/ul	0.00
46) Dimethylphthalate-d6	14.228	166	125482	18.615 ng/ul	0.00
49) Acenaphthylene-d8	14.528	160	156453	18.406 ng/ul	0.00
54) 4-Nitrophenol-d4	15.045	143	20190	18.504 ng/ul	0.00
60) Fluorene-d10	15.827	176	111019	18.289 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.950	200	22287	17.459 ng/ul	0.00
73) Anthracene-d10	17.683	188	181049	18.299 ng/ul	0.00
81) Pyrene-d10	19.963	212	204984	17.951 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.051	264	181332	17.880 ng/ul	0.00
Target Compounds				Qva	lue
2) 1,4-Dioxane	3.570	88	6570	7.440 ng/uL	86
5) Pyridine	3.987	79	43095	18.026 ng/ul	100
Benzaldehyde	7.331	77	36107	21.084 ng/ul	94
8) Phenol	7.383	94	49685	17.836 ng/ul	96
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.607	93	38195	18.123 ng/ul	97
12) 2-Chlorophenol	7.760	128	35476	17.978 ng/ul	95
<pre>13) 2-Methylphenol</pre>	8.641	108	36632	17.654 ng/ul	96
14) 2,2'-oxybis(1-Chloropr	8.717	45	55677	18.308 ng/ul	95
16) Acetophenone	9.029	105	61642	18.365 ng/ul	98
17) N-Nitroso-di-n-propyla	8.999	70	36499	18.923 ng/ul	96
18) 4-Methylphenol	8.976	108	39934	17.998 ng/ul	98
19) Hexachloroethane	9.287	117	14709	17.648 ng/ul	92
22) Nitrobenzene	9.416	77	51331	18.606 ng/ul	98
23) Isophorone	9.939	82	99003	18.471 ng/ul	98
25) 2-Nitrophenol	10.133	139	22229	18.082 ng/ul	98
26) 2,4-Dimethylphenol	10.186	107	46133	18.355 ng/ul	96
27) Bis(2-Chloroethoxy)met	10.415	93	54461	18.406 ng/ul	96
29) 2,4-Dichlorophenol	10.674	162	36072	18.198 ng/ul	97
30) Naphthalene	11.079	128	122243	18.025 ng/ul	99
32) 4-Chloroaniline	11.191	127	53481	18.080 ng/ul	99
33) Hexachlorobutadiene	11.344	225	24087	17.617 ng/ul	97
34) Caprolactam	11.955	113	14938m>	19.169 ng/ul>	11/29/21 J
35) 4-Chloro-3-methylphenol	12.307	107	44233	18.576 ng/ul	93

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\

Data File : BG051188.D

Acq On : 23 Nov 2021 17:07

Operator : CG/JU Sample : SSTDICV020

Misc

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 24 06:08:17 2021

 $\label{lem:quant_method} \textbf{Quant Method}: \textbf{Z:} \\ \textbf{SPAM-EPA-BG112321.M} \\ \\ \textbf{Quant Method}: \textbf{Z:} \\ \textbf{SPAM-EPA-BG112321.M} \\ \\ \textbf{Quant Method}: \textbf{Z:} \\ \textbf{Quant Method}: \textbf{Q:} \\ \textbf{Quant Method}: \textbf{Q:} \\ \textbf{$

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

Instrument : BNA_G ClientSampleId :

SICV427

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/24/2021 Supervised By :mohammad ahmed 11/30/2021

Compound	R.T.	QIon	Response	Conc Un	its De	v(Min)
36) 2-Methylnaphthalene	12.671	142	85761	18.592	ng/ul	97
37) 1-Methylnaphthalene	12.889	142	86304	18.185		97
39) 1,2,4,5-Tetrachloroben	13.036	216	48647	17.687		96
40) Hexachlorocyclopentadiene	13.000	237	26137	23.511	ng/ula	# 98
41) 2,4,6-Trichlorophenol	13.277	196	31205	18.080	ng/ul	97
42) 2,4,5-Trichlorophenol	13.359	196	31878	17.637	ng/ul	99
43) 1,1'-Biphenyl	13.664	154	119504	18.263	ng/ul	97
44) 2-Chloronaphthalene	13.717	162	93295	17.924	ng/ul	98
45) 2-Nitroaniline	13.923	65	34022	18.886	ng/ul	94
47) Dimethylphthalate	14.275	163	127086	18.626	ng/ul	99
48) 2,6-Dinitrotoluene	14.411	165	26264	18.325	ng/ul	95
50) Acenaphthylene	14.563	152	153221	18.245	ng/ul	98
51) 3-Nitroaniline	14.745	138	28346	20.008		92
52) Acenaphthene	14.898	153	100087	18.071	ng/ul	97
53) 2,4-Dinitrophenol	14.969	184	14400	18.177		‡ 87
55) 4-Nitrophenol	15.063	109	20482	21.639	ng/ul	93
56) Dibenzofuran	15.233	168	145026	18.154	ng/ul	98
57) 2,4-Dinitrotoluene	15.198	165	39208	19.153	٠.	[‡] 91
58) 2,3,4,6-Tetrachlorophenol	15.462	232	26181	18.446	ng/ul	93
59) Diethylphthalate	15.627	149	134867	18.831		100
61) Fluorene	15.879	166	116998	18.284	ng/ul	98
62) 4-Chlorophenyl-phenyle	15.862	204	62675	18.175		93
63) 4-Nitroaniline	15.909	138	28917m⊅	20.975	-	> ([Balai Ju
66) 4,6-Dinitro-2-methylph	15.968	198	21959	17.837	_	96
67) N-Nitrosodiphenylamine	16.079	169	108586	18.335	•	95
68) 4-Bromophenyl-phenylether	16.755	248	39655	17.886	-	94
69) Hexachlorobenzene	16.884	284	40850	18.069	_	94
70) Atrazine	17.019	200	45762	18.386	_	98
71) Pentachlorophenol	17.237	266	19501	19.467	_	96
72) Phenanthrene	17.624	178	205697	18.009	_	99
74) Anthracene	17.718	178	208283	18.361		99
75) 1,2,3,4-Tetrachloroben	13.641	216	51777	17.160		96
76) Pentachlorobenzene	15.151	250	49725	17.686	-	99
77) Carbazole	17.989	167	185821	18.662		99
78) Di-n-butylphthalate	18.517	149	237220	18.477		98
80) Fluoranthene	19.628	202	252052	17.971		97
82) Pyrene	19.992	202	247668	18.052		97
83) Butylbenzylphthalate	20.850	149	103100	18.076		95
84) 3,3'-Dichlorobenzidine	21.772	252	81844	18.626	-	96
85) Benzo(a)anthracene	21.866	228	229549	17.933	_	100
86) Bis(2-ethylhexyl)phtha	21.725	149	149473	18.212	•	99
87) Chrysene	21.931	228	220420	17.925	_	99
<pre>89) Di-n-octyl phthalate 90) Benzo(b)fluoranthene</pre>	22.995	149	253594	18.434		100
· · · · · · · · · · · · · · · · · · ·	24.199	252	232207	18.120	_	99
91) Benzo(k)fluoranthene 93) Benzo(a)pyrene	24.270	252	211582	17.594	<u> </u>	99
94) Indeno(1,2,3-cd)pyrene	25.127	252 276	221051	18.081		98
95) Dibenzo(a,h)anthracene	29.199	276 278	243706	17.813		97
96) Benzo(g,h,i)perylene	29.264 30.427	276	203832 204898	17.562 17.801	-	98 97
ytene		_,U 	204030	T1.00T	g/ u ±	<i>JI</i>

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