Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

1:08

Data File : BG050894.D Acq On : 9 Nov 2021

Operator : CG/JU Sample : PB140516BS

Misc

ALS Vial : 20 Sample Multiplier: 1

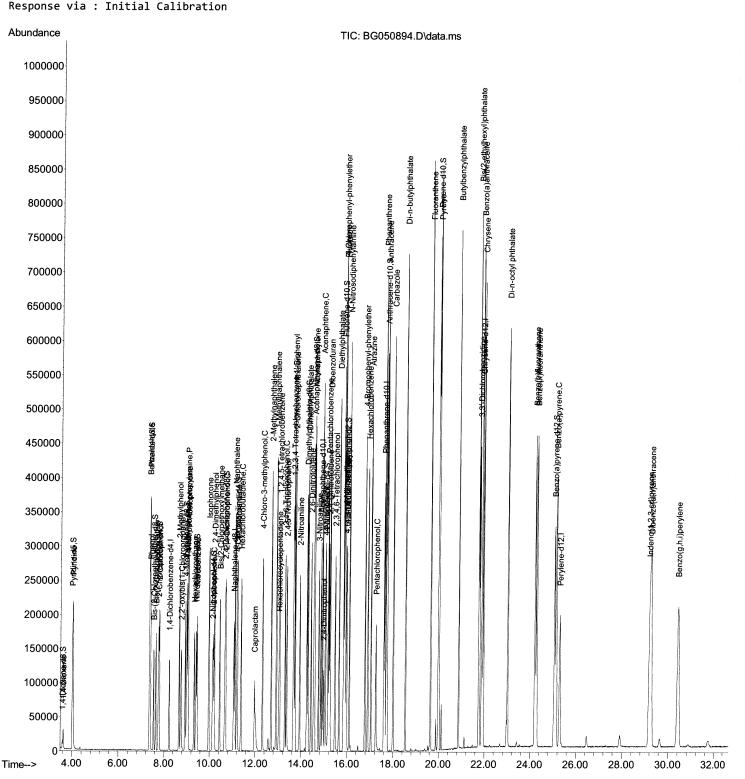
Quant Time: Nov 09 01:26:25 2021

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 14:49:05 2021

Instrument : BNA_G ClientSampleld :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/09/2021 Supervised By :mohammad ahmed 11/09/2021



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

Data File: BG050894.D

Acq On : 9 Nov 2021 1:08

Operator : CG/JU Sample : PB140516BS

Misc :

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Nov 09 01:26:25 2021

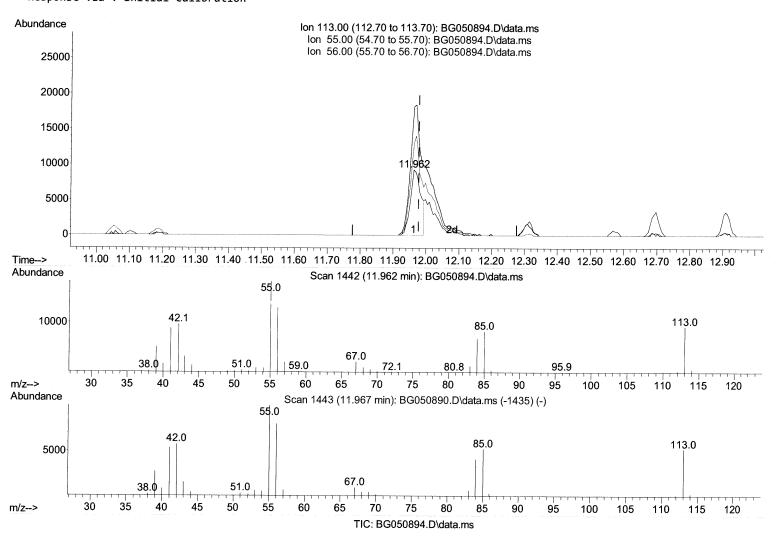
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 14:49:05 2021
Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/09/2021 Supervised By :mohammad ahmed 11/09/2021



(34) Caprolactam

11.962min (-0.015) 19.30 ng/ul

response	20499	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	183.80	196.42
56.00	136.50	140.41
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

Data File : BG050894.D

Acq On : 9 Nov 2021 1:08

Operator : CG/JU Sample : PB140516BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Nov 09 01:26:25 2021

 $\label{lem:quant_method} \mbox{Quant Methods} : \mbox{Z:\svoasrv} \mbox{HPCHEM1\BNA_G\Methods} : \mbox{Z:\svoasrv} \mbox{HPCHEM1\BNA_G\Methods} : \mbox{Methods} : \mbox{Z:\svoasrv} \mbox{HPCHEM1\BNA_G\Methods} : \mbox{Methods} : \mbox{Methods$

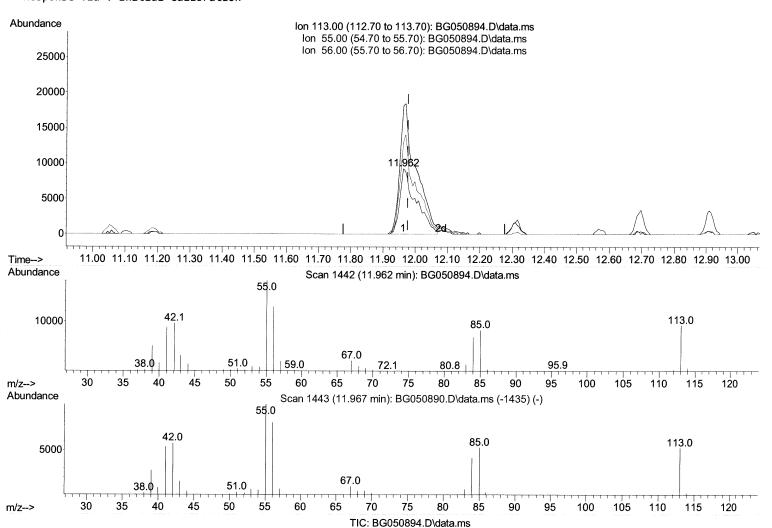
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 14:49:05 2021
Response via : Initial Calibration

Instrument : BNA_G ClientSampleId :

SLCS516

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/09/2021 Supervised By :mohammad ahmed 11/09/2021



(34) Caprolactam

11.962min (-0.015) 30.30 ng/ul m /////2しつい

response	32183	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	183.80	196.42
56.00	136.50	140.41
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

Data File : BG050894.D

Acq On : 9 Nov 2021 1:08

Operator : CG/JU Sample : PB140516BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Nov 09 01:26:25 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 14:49:05 2021
Response via : Initial Calibration

Instrument: BNA_G ClientSampleId: SLCS516

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/09/2021 Supervised By :mohammad ahmed 11/09/2021

Compound	R.T.	QIon	Response	Conc Ur	nits Dev	(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.226	152	36118	20.000	ng/ul	0.00
20) Naphthalene-d8	11.052		161118		ng/ul	0.00
38) Acenaphthene-d10	14.853	164	105294		ng/ul	0.00
64) Phenanthrene-d10	17.597		232089		ng/ul	0.00
79) Chrysene-d12	21.892		199202		ng/ul	-0.01
88) Perylene-d12	25.288	264	198490		ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.590	96	6577	5.877	ng/uL	0.00
4) Pyridine-d5	4.013	84	97833		ng/ul	0.00
7) Phenol-d5	7.368	99	112688		ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.544	67	74834		ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.756	132	79984	29.953	ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.931	113	85035		ng/ul	0.00
21) Nitrobenzene-d5	9.401	128	41069	29.995	ng/ul	0.00
24) 2-Nitrophenol-d4	10.129	143	45118		ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.670	165	77451		ng/ul	0.00
<pre>31) 4-Chloroaniline-d4</pre>	11.187	131	108483		ng/ul	0.00
46) Dimethylphthalate-d6	14.248	166	251350		ng/ul	0.00
49) Acenaphthylene-d8	14.548	160	316611		ng/ul	0.00
54) 4-Nitrophenol-d4	15.041	143	43878		ng/ul	0.00
60) Fluorene-d10	15.840	176	221782		ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.958	200	42923		ng/ul	0.00
73) Anthracene-d10	17.697	188	336663	30.681		0.00
81) Pyrene-d10	19.971	212	398785	30.995		0.00
92) Benzo(a)pyrene-d12	25.059	264	333329	30.378	•	-0.01
Farget Compounds					0va	lue
2) 1,4-Dioxane	3.625	88	15958	12.983	ng/uL	98
5) Pyridine	4.030	79	109266	31.531		98
6) Benzaldehyde	7.362	77	85769	35.292		93
8) Phenol	7.397	94	126847	31.824		99
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.638	93	95764	32.098		99
12) 2-Chlorophenol	7.785	128	87135	32.138		97
<pre>13) 2-Methylphenol</pre>	8.660	108	92176	31.287		98
14) 2,2'-oxybis(1-Chloropr	8.748	45	150695	32.067	ng/ul	99
16) Acetophenone	9.054	105	144995	30.770	ng/ul	99
17) N-Nitroso-di-n-propyla	9.031	70	86695	30.493	ng/ul	99
18) 4-Methylphenol	8.995	108	96731	30.836	ng/ul	93
19) Hexachloroethane	9.318	117	37198	32.814		99
22) Nitrobenzene	9.442	77	127411	33.365	ng/ul	98
23) Isophorone	9.965	82	236817	31.954	ng/ul	100
25) 2-Nitrophenol	10.159	139	51138	33.481	ng/ul	96
26) 2,4-Dimethylphenol	10.200	107	97668	29.055		98
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.441	93	132197	33.106	ng/ul	100
29) 2,4-Dichlorophenol	10.693	162	83130	33.232		97
30) Naphthalene	11.105	128	289692	32.881	_	99
22\ 4 Chlamandline		127	117988	30.596		99
32) 4-Chloroaniline	11.210	12/	11/500	30.330	iig/ ui	,,
33) Hexachlorobutadiene	11.210 11.375	225	53480	32.573		97
·				32.573		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

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Acq On : 9 Nov 2021 1:08

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Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Nov 09 01:26:25 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration

Instrument: BNA_G ClientSampleId: SLCS516

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/09/2021 Supervised By: mohammad ahmed 11/09/2021

Compound	R.T.	QIon	Response	Conc Un	its Dev(N	Min)
36) 2-Methylnaphthalene	12.691	142	194875	32.468	ng/ul	 97
37) 1-Methylnaphthalene	12.908	142	195176	32.092	_	97
39) 1,2,4,5-Tetrachloroben	13.055	216	103628	33.777		98
40) Hexachlorocyclopentadiene	13.026	237	38330	26.017	-	96
41) 2,4,6-Trichlorophenol	13.290	196	68240	33.992	_	99
42) 2,4,5-Trichlorophenol	13.367	196	73250	33.986		99
43) 1,1'-Biphenyl	13.690	154	258383	33.572	ng/ul	99
44) 2-Chloronaphthalene	13.737	162	204576	33.920	ng/ul	99
45) 2-Nitroaniline	13.937	65	79487	33.172	ng/ul	97
	14.295	163	269834	33.500	ng/ul	99
	14.424	165	58022	34.417	ng/ul	98
	14.577	152	336824	33.485	ng/ul	98
	14.753	138	60792	34.866		86
	14.918	1 53	225515	34.096	ng/ul	98
	14.965	184	26093	28.057	ng/ul	97
	15.053	109	44186	32.985	ng/ul	92
	15.247	168	320165	33.818	ng/ul	99
	15.206	165	82609	34.335	ng/ul#	99
	15.470	232	57606	34.010	ng/ul	98
	15.646	149	288634	33.477		99
•	15.899	166	250928	33.487	ng/ul	99
	15.881	204	132523	33.963	ng/ul	97
	15.917	138	60110	34.750		96
	15.969	198	45574	33.212	ng/ul	96
	16.093	169	222363	34.276		98
	16.774	248	79233	34.320	ng/ul	96
· ·	16.898	284	80217	33.798		95
	17.033	200	89138	32.401		98
	17.244	266	34074	31.271		95
	17.638	178	426285	34.409	_	99
	L7.732	178	409385	32.934		99
	L3.660	216	107772	34.104	-	99
	L5.164	250	95415	32.586	•	98
	L7.996	167	392786	35.255		98
	18.531	149	504367	34.452		100
	19.636	202	520073	33.681		98
	20.000	202	502585	33.312		99
	20.864	149	220179	33.937		98
	21.780	252	155222	31.996		98
	21.874	228	465225	33.735		100
	21.745	149	314954	33.819	_	100
	21.945	228	440148	33.411		100
	13.014	149	533362	33.006		100
	4.201	252	473703	33.500	_	100
	4.277	252	433082	32.639		99
	5.129	252	440709	32.724	•	98 07
	9.195 9.266	276 278	496748 421641	33.134		97 07
			415132	33.239 33.080		97 97
					_	97

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