Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050814.D

Acq On : 2 Nov 2021 12:11

Operator : CG/JU Sample : SSTD08016

Misc

ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 02 14:33:17 2021

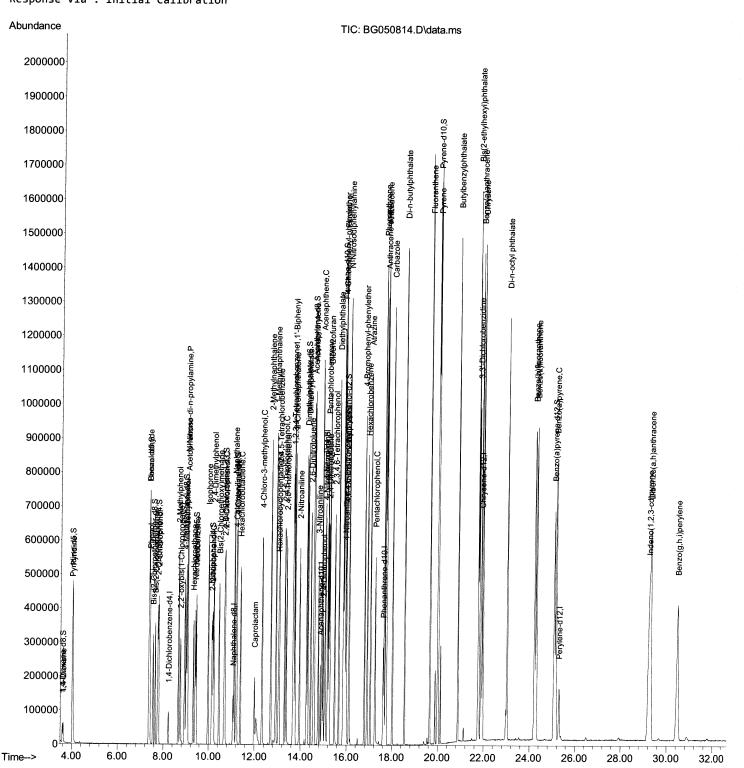
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : SSTD080416

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050814.D

Acq On : 2 Nov 2021 12:11

Operator : CG/JU Sample : SSTD08016

Misc

ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 02 14:33:17 2021

 $\label{lem:quant_def} Quant \ \ Method: Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M$

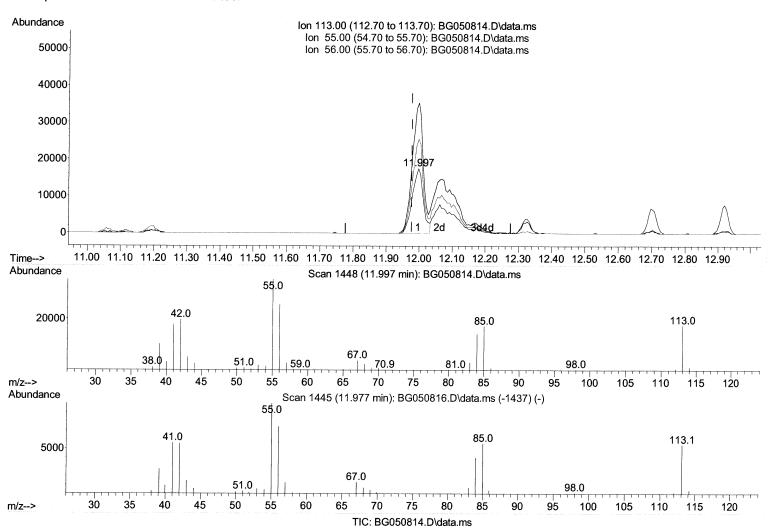
Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



(34) Caprolactam

11.997min (+ 0.020) 50.84 ng/ul

response	40926	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	183.80	201.28
56.00	136.50	145.31
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050814.D

Acq On : 2 Nov 2021 12:11

Operator : CG/JU Sample : SSTD08016

Misc

ALS Vial : 6 Sample Multiplier: 1

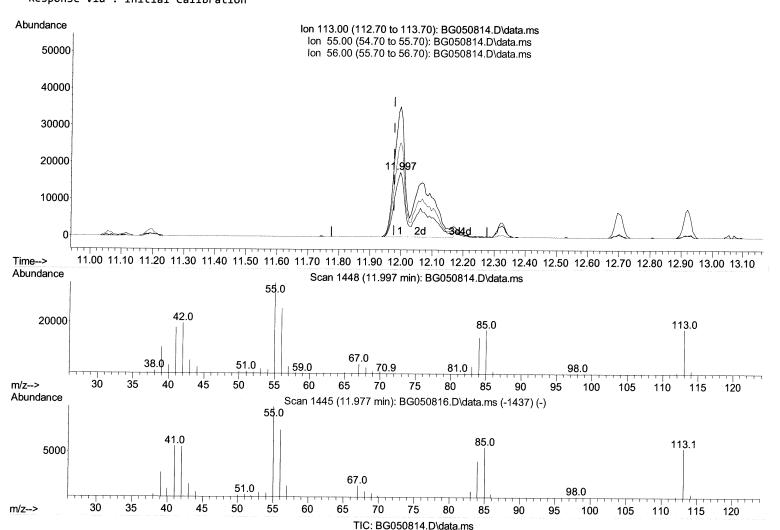
Quant Time: Nov 02 14:33:17 2021

 $\label{lem:quant_def} Quant \ \mbox{Methods\SFAM-EPA-BG110321.M}$

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration Instrument :
BNA_G
ClientSampleId :
SSTD080416

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



(34) Caprolactam

11.997min (+ 0.020) 92.23 ng/ul m \\04/2\JU

response	74239	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	183.80	201.28
56.00	136.50	145.31
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050814.D

Acq On : 2 Nov 2021 12:11

Operator : CG/JU Sample : SSTD08016

Misc :

ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 02 14:33:17 2021

 $\label{lem:quant_method} \mbox{Quant Methods} : \mbox{Z:\svoasrv\hPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M}$

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : SSTD080416

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/02/2021 Supervised By: mohammad ahmed 11/08/2021

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.237	152	25929	20.000 ng/ul	0.00
20) Naphthalene-d8	11.063		122054	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.858		81312	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.602	188	186323	20.000 ng/ul	0.00
79) Chrysene-d12	21.903	240	149539	20.000 ng/ul	0.00
88) Perylene-d12	25.305	264	149023	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.589	96	30473	37.932 ng/uL	0.00
4) Pyridine-d5	4.012	84	225160	93.660 ng/ul	0.00
7) Phenol-d5	7.379	99	260617	94.216 ng/ul	0.00
Bis-(2-Chloroethyl)eth	7.549	67	162586	90.990 ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.761	132	181449	94.651 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.942	113	201738	92.640 ng/ul	0.00
21) Nitrobenzene-d5	9.412	128	98696	95. 1 54 ng/ul	0.00
24) 2-Nitrophenol-d4	10.134	143	112043	97. 1 50 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.675	165	184412	94.924 ng/ul	0.00
31) 4-Chloroaniline-d4	11.192	131	266239	90. 49 4 ng/ul	0.00
46) Dimethylphthalate-d6	14.253	166	572966	92.104 ng/ul	0.00
49) Acenaphthylene-d8	14.559	160	711462	91.796 ng/ul	0.00
54) 4-Nitrophenol-d4	15.058	143	110158	97.663 ng/ul	0.01
60) Fluorene-d10	15.851	176	503328	91.334 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.969	200	111661	98.847 ng/ul	0.00
73) Anthracene-d10	17.708	188	772062	87.643 ng/ul	0.00
81) Pyrene-d1092) Benzo(a)pyrene-d12	19.982 25.087	212 264	877647 757345	90.868 ng/ul 91.930 ng/ul	0.00 0.01
Target Compounds				_	
2) 1,4-Dioxane	3.624	00	22127	Qva]	
5) Pyridine	4.036	88 79	33137	37.554 ng/uL	95
6) Benzaldehyde	7.367	79 77	233555 163804	93.540 ng/ul	98 05
8) Phenol	7.408	94	264586	93.887 ng/ul 92.465 ng/ul	95 99
10) Bis(2-Chloroethyl)ether	7.643	93	198009	92.448 ng/ul	99
12) 2-Chlorophenol	7.796	128	180382	92.674 ng/ul	97
13) 2-Methylphenol	8.671	108	198974	94.076 ng/ul	99
14) 2,2'-oxybis(1-Chloropr	8.754	45	305448	90.539 ng/ul	99
16) Acetophenone	9.065	105	305559	90.324 ng/ul	96
17) N-Nitroso-di-n-propyla	9.053	70	185104	90.689 ng/ul	96
18) 4-Methylphenol	9.006	108	207498	92.140 ng/ul	98
19) Hexachloroethane	9.324	117	76844	94.426 ng/ul	98
22) Nitrobenzene	9.453	77	263503	91.089 ng/ul	98
23) Isophorone	9.976	82	515609	91.839 ng/ul	99
25) 2-Nitrophenol	10.164	139	112552	97.275 ng/ul	96
26) 2,4-Dimethylphenol	10.211	107	234948	92.265 ng/ul	99
27) Bis(2-Chloroethoxy)met	10.452	93	277795	91.833 ng/ul	98
29) 2,4-Dichlorophenol	10.704	162	179218	94.574 ng/ul	98
30) Naphthalene	11.116	128	604059	90.507 ng/ul	98
32) 4-Chloroaniline	11.221	127	267355	91.518 ng/ul	100
33) Hexachlorobutadiene	11.380	225	114851	92.341 ng/ul	97
34) Caprolactam	11.997	113	74239m >	<u>.</u>	utic (Poil)
35) 4-Chloro-3-methylphenol	12.326	107	227529	93.990 ng/ul	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050814.D

Acq On : 2 Nov 2021 12:11

Operator : CG/JU Sample : SSTD08016

Misc :

ALS Vial : 6 Sample Multiplier: 1

Quant Time: Nov 02 14:33:17 2021

Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 14:25:39 2021
Response via : Initial Calibration

Instrument: BNA_G ClientSampleId: SSTD080416

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
36) 2-Methylnaphthalene	12.702	142	410220	90.220 ng/ul	100
37) 1-Methylnaphthalene	12.919	142	416560	90.415 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.060	216	221214	93.369 ng/ul	96
40) Hexachlorocyclopentadiene	13.037	237	121452	106.749 ng/ul	99
41) 2,4,6-Trichlorophenol	13.301	196	156462	100.923 ng/ul	99
42) 2,4,5-Trichlorophenol	13.378	196	162623	97.707 ng/ul	99
43) 1,1' -Biphenyl	13.695	154	537298	90.403 ng/ul	100
44) 2-Chloronaphthalene	13.748	162	425703	91.402 ng/ul	98
45) 2-Nitroaniline	13.948	65	179140	96.810 ng/ul	96
47) Dimethylphthalate	14.306	163	562803	90.480 ng/ul	99
48) 2,6-Dinitrotoluene	14.435	165	127315	97.793 ng/ul	97
50) Acenaphthylene	14.588	152	699789	90.089 ng/ul	98
51) 3-Nitroaniline	14.764	138	128042	95.094 ng/ul	89
52) Acenaphthene	14.923	153	470131	92.044 ng/ul	96
53) 2,4-Dinitrophenol	14.976	184	76043	105.881 ng/ul	96
55) 4-Nitrophenol	15.070	109	100802	97.444 ng/ul	98
56) Dibenzofuran	15.258	168	652678	89.273 ng/ul	97
57) 2,4-Dinitrotoluene	15.223	165	178339	95.986 ng/ul	97
58) 2,3,4,6-Tetrachlorophenol	15.481	232	134210	102.606 ng/ul	99
59) Diethylphthalate	15.657	149	606802	91.137 ng/ul	98
61) Fluorene	15.904	166	512182	88.511 ng/ul	100
62) 4-Chlorophenyl-phenyle	15.887	204	275733	91.507 ng/ul	99
63) 4-Nitroaniline	15.934	138	127917	95.762 ng/ul	95
66) 4,6-Dinitro-2-methylph	15.986	198	108154	98.177 ng/ul#	99
67) N-Nitrosodiphenylamine	16.104	169	470945	90.424 ng/ul	99
68) 4-Bromophenyl-phenylether	16.785	248	175162	94.509 ng/ul	95
69) Hexachlorobenzene	16.909	284	176400	92.580 ng/ul	98
70) Atrazine	17.044	200	201949	91.438 ng/ul	99
71) Pentachlorophenol	17.250	266	97395	111.337 ng/ul	95
72) Phenanthrene	17.649	178	879097	88.388 ng/ul	97
74) Anthracene	17.743	178	861883	86.368 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.666	216	236347	93.161 ng/uL	97
76) Pentachlorobenzene	15.176	250	220732	93.901 ng/uL	99
77) Carbazole	18.008	167	812259	90.812 ng/ul	97
78) Di-n-butylphthalate	18.542	149	1023947	87.122 ng/ul	99
80) Fluoranthene	19.647	202	1047953	90.408 ng/ul	99
82) Pyrene	20.011	202	998813	88.188 ng/ul	98
83) Butylbenzylphthalate	20.875	149	453646	93. 1 44 ng/ul	99
84) 3,3'-Dichlorobenzidine	21.791	252	343392	94.292 ng/ul	100
85) Benzo(a)anthracene	21.885	228	942053	90.998 ng/ul	98
86) Bis(2-ethylhexyl)phtha	21.750	149	645624	92.350 ng/ul	100
87) Chrysene	21.956	228	897381	90.741 ng/ul	98
89) Di-n-octyl phthalate	23.025	149	1098238	90.522 ng/ul	100
90) Benzo(b)fluoranthene	24.224	252	953056	89.773 ng/ul	98
91) Benzo(k)fluoranthene	24.300	252	893529	89.693 ng/ul	98
93) Benzo(a)pyrene	25.164	252	914312	90.424 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	29.247	276	1044221	92.685 ng/ul	98
95) Dibenzo(a,h)anthracene	29.306	278	872339	91.541 ng/ul	100
96) Benzo(g,h,i)perylene	30.481	276	872500	92.710 ng/ul	98

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