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

Data File : BG051204.D

Acq On : 24 Nov 2021 9:22

Operator : CG/JU Sample : PB140881BS

Misc

ALS Vial : 25 Sample Multiplier: 1

Quant Time: Nov 24 12:45:43 2021

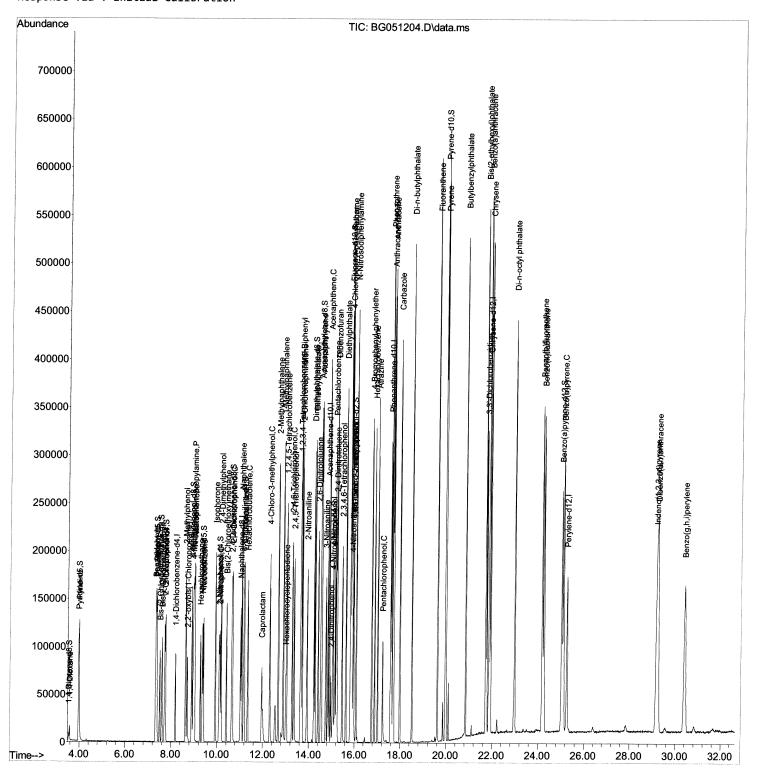
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument : BNA\_G ClientSampleld : SLCS881

## **Manual IntegrationsAPPROVED**

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



#### Quantitation Report (Qedit)

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

Data File: BG051204.D

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Operator : CG/JU Sample : PB140881BS

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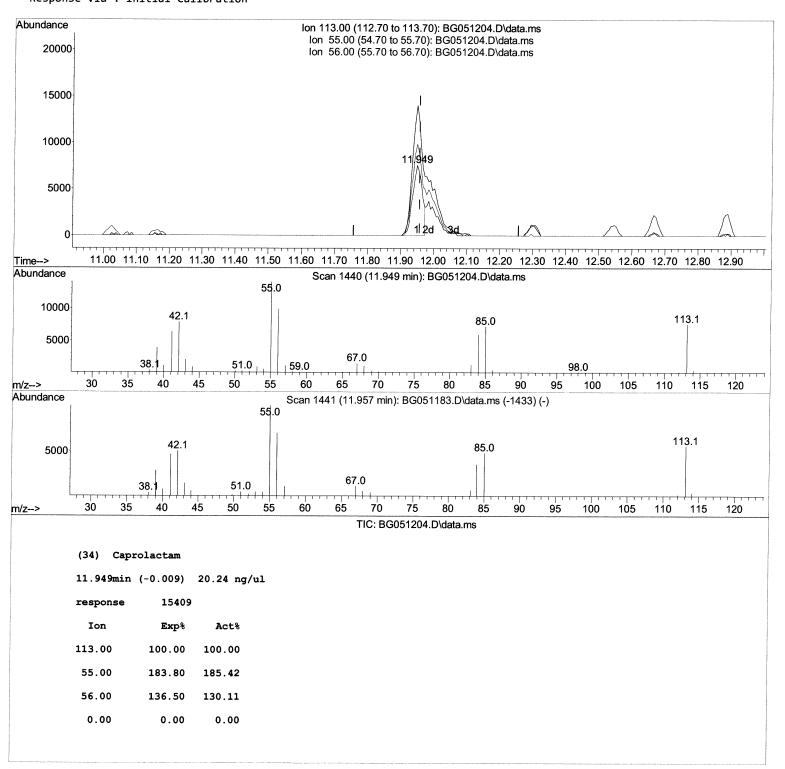
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
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Misc

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 $\label{thm:lem1_BNA_GMethods} \textbf{Quant Method}: Z:\\ \textbf{Z:}\\ \textbf{SFAM-EPA-BG112321.M}$ 

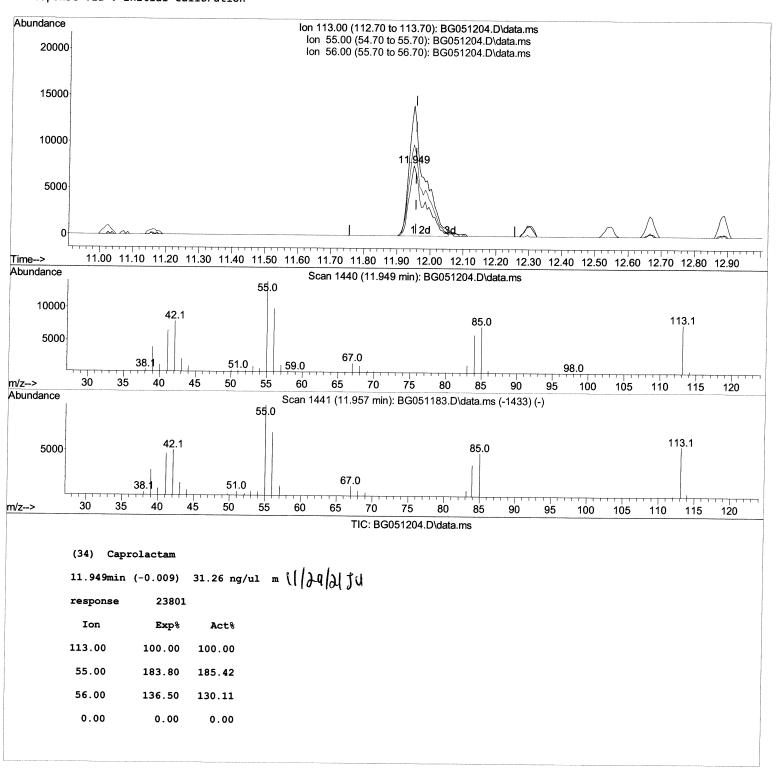
Quant Title : SVOA CALIBRATION

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



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# **Manual IntegrationsAPPROVED**

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

Compound				Conc Units Dev(	(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.200	152	26372	20.000 ng/ul	0.00
20) Naphthalene-d8	11.026	136	121770	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.828	164	86476	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.577		192395	20.000 ng/ul	0.00
79) Chrysene-d12	21.878	240	165452	20.000 ng/ul	0.00
88) Perylene-d12	25.274	264	167661	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.535	96	4183	5.512 ng/uL	0.00
4) Pyridine-d5	3.964	84	61026	27.404 ng/ul	-0.01
7) Phenol-d5	7.354	99	78516	30.124 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.513	67	49049	29.963 ng/ul	0.00
11) 2-Chlorophenol-d4	7.730	132	57689	30.737 ng/ul	0.00
15) 4-Methylphenol-d8	8.911	113	64548	30.689 ng/ul	0.00
21) Nitrobenzene-d5	9.375	128	30328	29.504 ng/ul	0.00
24) 2-Nitrophenol-d4	10.098	143	35571	30.677 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.644	165	62471	31.754 ng/ul	0.00
31) 4-Chloroaniline-d4	11.161	131	80348	27.912 ng/ul	0.00
46) Dimethylphthalate-d6	14.222	166	199263	29.947 ng/ul	0.00
49) Acenaphthylene-d8	14.528	160	257071	30.639 ng/ul	0.00
54) 4-Nitrophenol-d4	15.039	143	31971	29.684 ng/ul	0.00
60) Fluorene-d10	15.821	176	179619	29.977 ng/ul	0.00
65) 4,6-Dinitro-2-methylph		200	35434	29.847 ng/ul	0.00
73) Anthracene-d10	17.677	188	276078	30.003 ng/ul	0.00
81) Pyrene-d10	19.957	212	310233	30.989 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.039	264	276489	30.878 ng/ul	0.00
arget Compounds				Qva	lue
2) 1,4-Dioxane	3.576	88	8558	9.999 ng/uL	98
5) Pyridine	3.987	79	61532	26.554 ng/ul	99
6) Benzaldehyde	7.331	77	49915	30.071 ng/ul	93
8) Phenol	7.383	94	82462	30.540 ng/ul	99
10) Bis(2-Chloroethyl)ether	7.607	93	61805	30.256 ng/ul	96
12) 2-Chlorophenol	7.759	128	58567	30.621 ng/ul	98
13) 2-Methylphenol	8.641	108	61930	30.792 ng/ul	95
14) 2,2'-oxybis(1-Chloropr		45	89072	30.217 ng/ul	97
16) Acetophenone	9.029	105	97208	29.879 ng/ul	97
17) N-Nitroso-di-n-propyla	8.999	70	57809	30.922 ng/ul	97
18) 4-Methylphenol	8.970	108	66319	30.837 ng/ul	100
19) Hexachloroethane	9.281	117	23748	29.396 ng/ul	96
22) Nitrobenzene	9.416	77	82816	30.726 ng/ul	98
23) Isophorone	9.933	82	159620	30.482 ng/ul	99
25) 2-Nitrophenol	10.133	139	36617	30.488 ng/ul	96
26) 2,4-Dimethylphenol	10.180	107	73907	30.098 ng/ul	98
27) Bis(2-Chloroethoxy)met	10.409	93	88810	30.721 ng/ul	98
29) 2,4-Dichlorophenol	10.403	162	59615	30.721 ng/ul	96
30) Naphthalene	11.073	128	197890	29.867 ng/ul	98
32) 4-Chloroaniline		127	81212	28.102 ng/ul	100
22) 4-CHIOLOGHITTINE	11.185			_	
33) Heyachlorobutadione	11 229	フフに			
<ul><li>33) Hexachlorobutadiene</li><li>34) Caprolactam</li></ul>	11.338 11.949	225 113	37740 23801m 🦴	28.253 ng/ul 31.262 ng/ul >	100 (( 24 2)

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

Instrument : BNA\_G ClientSampleld : SLCS881

# 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)
36) 2-Methylnaphthalene	12.665	142	139316	30.913 ng/ul	99
37) 1-Methylnaphthalene	12.883	142	140841	30.376 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.030	216	82084	30.235 ng/ul	96
40) Hexachlorocyclopentadiene	13.000	237	18939	17.259 ng/ul	99
41) 2,4,6-Trichlorophenol	13.271	196	52568	30.856 ng/ul	97
42) 2,4,5-Trichlorophenol	13.353	196	56504	31.672 ng/ul	99
43) 1,1'-Biphenyl	13.664	154	192461	29.798 ng/ul	99
44) 2-Chloronaphthalene	13.711	162	152091	29.602 ng/ul	98
45) 2-Nitroaniline	13.923	65	57572	32.377 ng/ul	94
47) Dimethylphthalate	14.269	163	201827	29.967 ng/ul	99
48) 2,6-Dinitrotoluene	14.405	165	44802	31.668 ng/ul	97
50) Acenaphthylene	14.557	152	249890	30.145 ng/ul	97
51) 3-Nitroaniline	14.740	138	41851	29.928 ng/ul	90
52) Acenaphthene	14.892	153	167588	30.655 ng/ul	98
53) 2,4-Dinitrophenol	14.963	184	18687	23.897 ng/ul#	83
55) 4-Nitrophenol	15.057	109	28508	30.512 ng/ul	90
56) Dibenzofuran	15.227	168	235627	29.881 ng/ul	99
57) 2,4-Dinitrotoluene	15.198	165	63266	31.310 ng/ul	91
58) 2,3,4,6-Tetrachlorophenol	15.456	232	45636	32.575 ng/ul	97
59) Diethylphthalate	15.621	149	216495	30.624 ng/ul	99
61) Fluorene	15.874	166	194291	30.760 ng/ul	97
62) 4-Chlorophenyl-phenyle	15.856	204	101165	29.720 ng/ul	97
63) 4-Nitroaniline	15.903	138	43938	32.287 ng/ul	95
66) 4,6-Dinitro-2-methylph	15.962	198	33748	29.475 ng/ul#	96
67) N-Nitrosodiphenylamine	16.073	169	174437	31.670 ng/ul	97
68) 4-Bromophenyl-phenylether	16.755	248	65188	31.614 ng/ul	94
69) Hexachlorobenzene	16.878	284	65793	31.291 ng/ul	97
70) Atrazine	17.019	200	68432	29.563 ng/ul	99
71) Pentachlorophenol	17.231	266	22445	24.091 ng/ul	95
72) Phenanthrene	17.619	178	328818	30.954 ng/ul	99
74) Anthracene	17.713	178	324720	30.779 ng/ul	97
<ul><li>75) 1,2,3,4-Tetrachloroben</li><li>76) Pentachlorobenzene</li></ul>	13.635	216	84610	30.150 ng/uL	99
77) Carbazole	15.145	250	79192	30.286 ng/uL	97
78) Di-n-butylphthalate	17.983 18.512	167	288440	31.147 ng/ul	99
80) Fluoranthene	19.622	149 202	366613	30.703 ng/ul	99
82) Pyrene	19.986	202	388888 379050	31.627 ng/ul	98 97
83) Butylbenzylphthalate	20.844	149	157211	31.514 ng/ul 31.440 ng/ul	97 96
84) 3,3'-Dichlorobenzidine	21.767	252	113586	29.486 ng/ul	
85) Benzo(a)anthracene	21.861	228	350995	31.278 ng/ul	99 99
86) Bis(2-ethylhexyl)phtha	21.720	149	225796	31.380 ng/ul	100
87) Chrysene	21.931	228	331299	30.731 ng/ul	99
89) Di-n-octyl phthalate	22.983	149	389685	32.082 ng/ul	100
90) Benzo(b)fluoranthene	24.187	252	350588	30.985 ng/ul	99
91) Benzo(k)fluoranthene	24.258	252	326339	30.735 ng/ul	99
93) Benzo(a)pyrene	25.116	252	333169	30.864 ng/ul	99
94) Indeno(1,2,3-cd)pyrene	29.181	276	374953	31.040 ng/ul	99
95) Dibenzo(a,h)anthracene	29.240	278	314924	30.731 ng/ul	96
96) Benzo(g,h,i)perylene	30.415	276	315434	31.037 ng/ul	98

<sup>(#) =</sup> qualifier out of range (m) = manual integration (+) = signals summed