Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

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

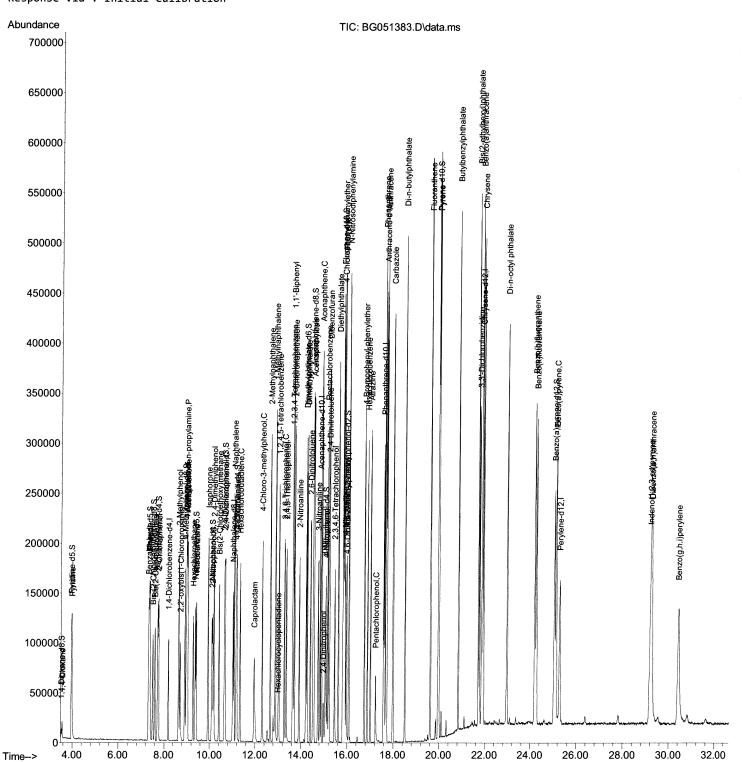
Quant Title : SVOA CALIBRATION

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



# Manual IntegrationsAPPROVED

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

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

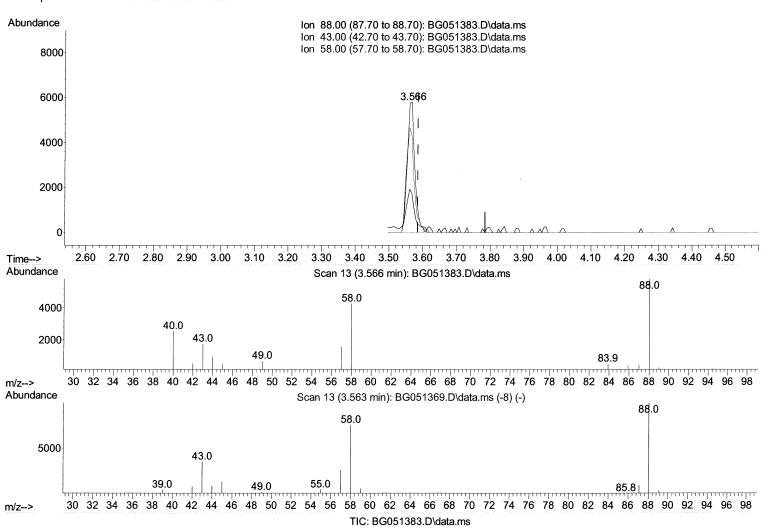
Quant Title : SVOA CALIBRATION

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



### (2) 1,4-Dioxane

3.566min (-0.019) 10.19 ng/uL

response	9219	
Ion	Ехр%	Act%
88.00	100.00	100.00
43.00	28.70	30.32
58.00	78.00	73.62
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

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

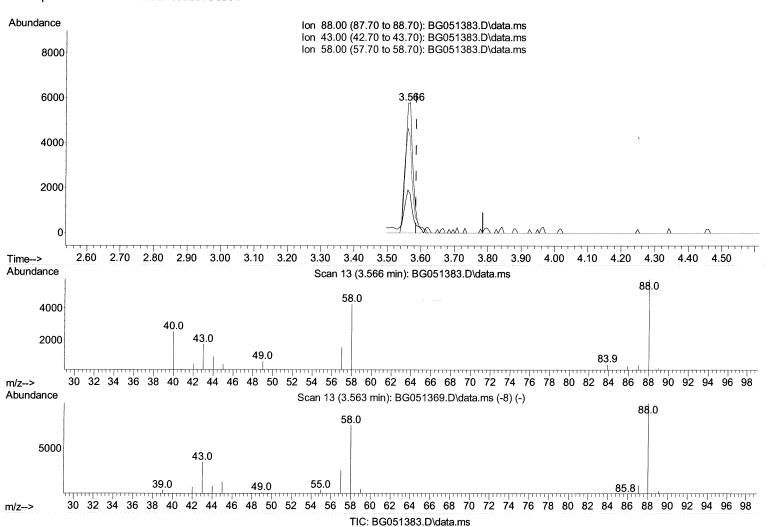
Quant Title : SVOA CALIBRATION

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

Instrument: BNA\_G ClientSampleld: SLCS156

# Manual IntegrationsAPPROVED

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



#### (2) 1,4-Dioxane

3.566min (-0.019) 10.27 ng/uL m \2////21 JU

response	9295		
Ion	Ежр%	Act%	
88.00	100.00	100.00	
43.00	28.70	30.32	
58.00	78.00	73.62	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

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

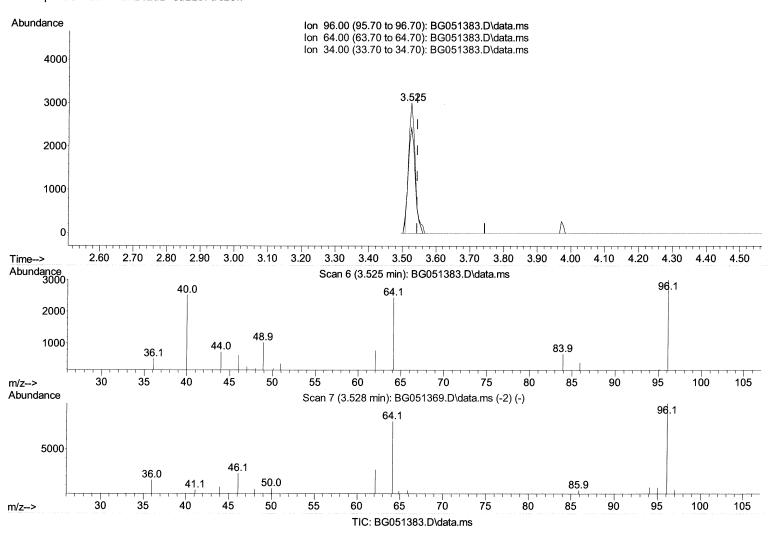
Quant Title : SVOA CALIBRATION

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)

3.525min (-0.019) 5.16 ng/uL

response	4141			
Ion	Ехр%	Act%		
96.00	100.00	100.00		
64.00	77.60	81.26		
34.00	0.00	0.00		
0.00	0.00	0.00		

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

 $\label{thm:constraint} Quant \ \mbox{Methods\SFAM-EPA-BG112321.M}$ 

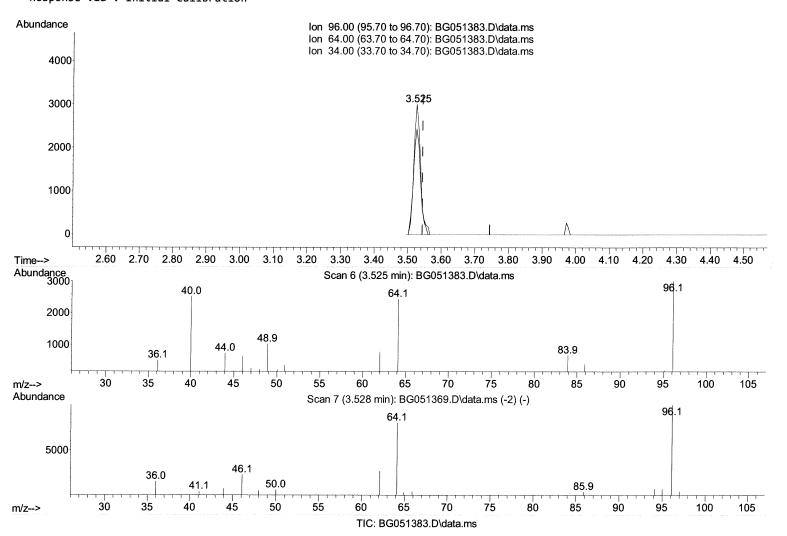
Quant Title : SVOA CALIBRATION

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)

3.525min (-0.019) 5.34 ng/uL m | 2//(/2\JU

response	4280	
Ion	Ехр%	Act%
96.00	100.00	100.00
64.00	77.60	81.26
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

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

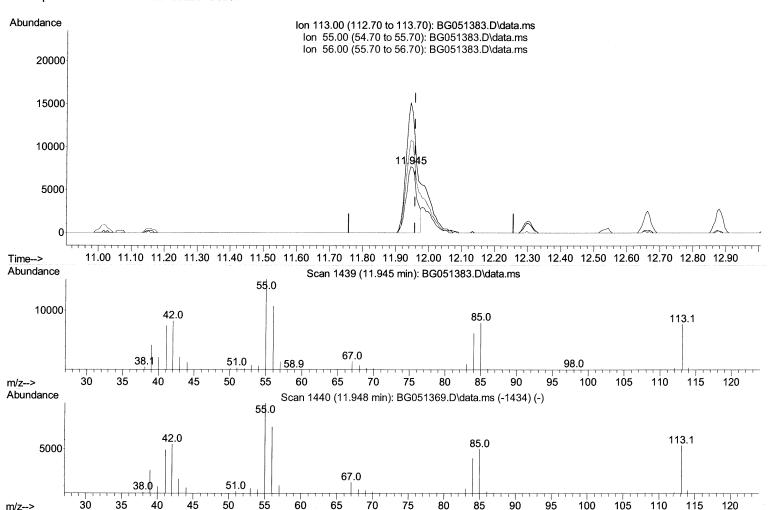
Quant Title : SVOA CALIBRATION

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: BG051383.D\data.ms

#### (34) Caprolactam

11.945min (-0.013) 21.80 ng/ul

response	17380		
Ion	Exp%	Act%	
113.00	100.00	100.00	
55.00	183.80	196.39	
56.00	136.50	139.32	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

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

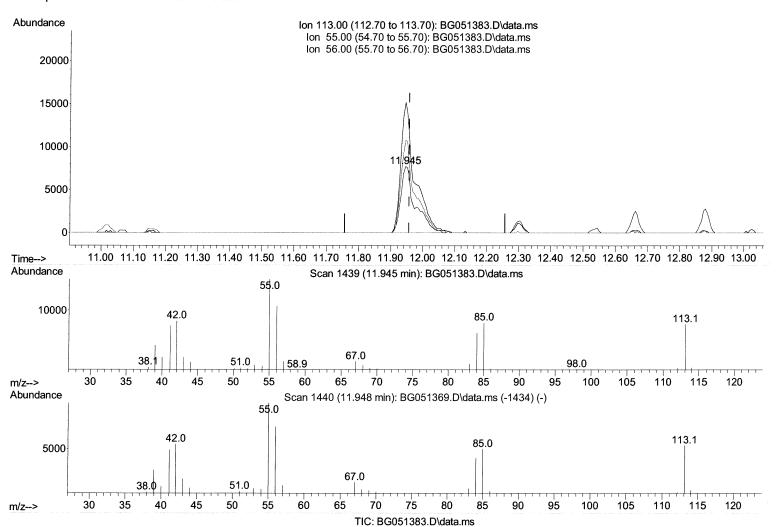
Quant Title : SVOA CALIBRATION

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

11.945min (-0.013) 30.28 ng/ul m 2///2/JU

response	24148	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	183.80	196.39
56.00	136.50	139.32
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

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

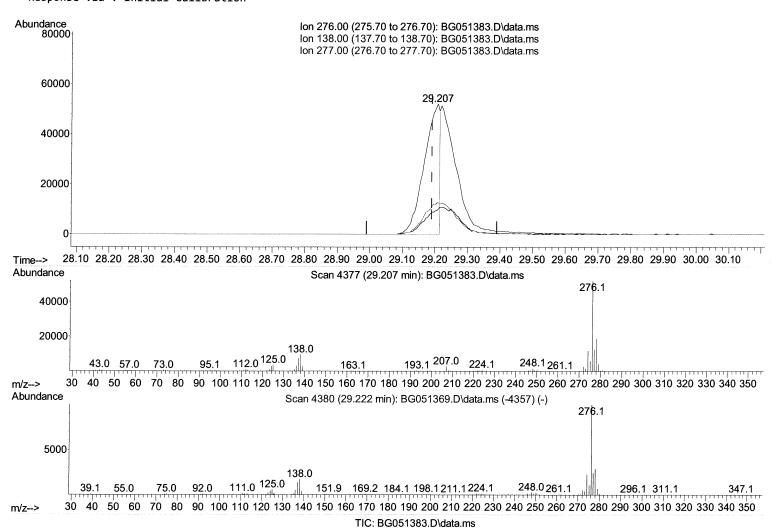
Quant Title : SVOA CALIBRATION

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



#### (94) Indeno (1,2,3-cd) pyrene

29.207min (+ 0.017) 15.85 ng/ul

response	173138	
Ion	Ежр%	Act%
276.00	100.00	100.00
138.00	19.40	18.47
277.00	25.60	24.42
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\

Data File: BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 2021

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

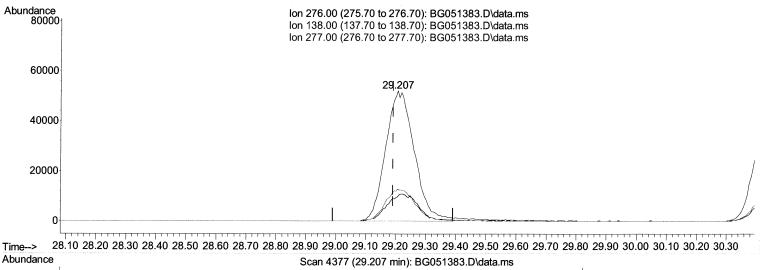
Quant Title : SVOA CALIBRATION

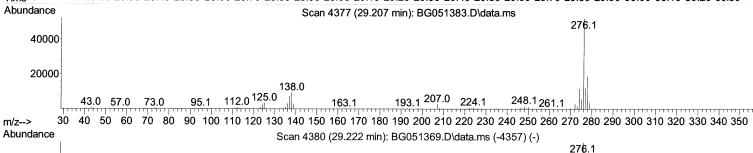
QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

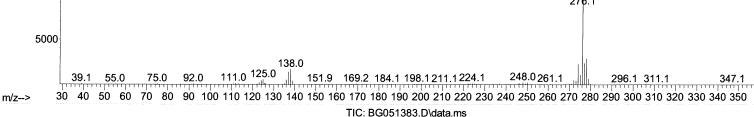


# Manual IntegrationsAPPROVED

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







### (94) Indeno (1,2,3-cd) pyrene

29.207min (+ 0.017) 31.34 ng/ul m \\ \( \) \( \) \( \)

response	342267	
Ion	Ехр%	Act%
276.00	100.00	100.00
138.00	19.40	18.47
277.00	25.60	24.42
0.00	0.00	0.00

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 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 ClientSampleId : SLCS156

# Manual IntegrationsAPPROVED

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

	Compound	R.T.	QIon	Response	Conc Un	its Dev	/(Min)
Interna	al Standards						
1) 1,	4-Dichlorobenzene-d4	8.190	152	27875	20.000	ng/ul	-0.01
20) Na	phthalene-d8	11.016	136	127538	20.000		-0.01
38) Ac	enaphthene-d10	14.824	164	84982	20.000	ng/ul	0.00
64) Ph	enanthrene-d10	17.573	188	185343	20.000	ng/ul	0.00
79) Ch	rysene-d12	21.874	240	155022	20.000	ng/ul	0.00
88) Pe	rylene-d12	25.276	264	151585	20.000	ng/ul	0.00
System	Monitoring Compounds						1116 t-
3) 1,	4-Dioxane-d8	3.525	96	4280m	> 5.336	ng/uL;	>-0.02 12/1(121J'
4) Py	ridine-d5	3.954	84	60289	25.614	ng/ul	-0.02
7) Ph	enol-d5	7.350	99	84655	30.728	ng/ul	0.00
9) Bi	s-(2-Chloroethyl)eth	7.503	67	53101	30.690	ng/ul	-0.01
11) 2-	Chlorophenol-d4	7.720	132	61475	30.988	ng/ul	-0.01
15) 4-	Methylphenol-d8	8.907	113	67907	30.545	ng/ul	0.00
21) Ni	trobenzene-d5	9.371	128	32660	30.336		0.00
24) 2-	Nitrophenol-d4	10.094	143	37464	30.848	ng/ul	0.00
28) 2,	4-Dichlorophenol-d3	10.646	165	64267	31.189	_	0.00
31) 4-	Chloroaniline-d4	11.157	131	79921	26.508	ng/ul	0.00
46) Di	methylphthalate-d6	14.218	166	199096	30.448	ng/ul	0.00
49) Ac	enaphthylene-d8	14.518	160	256959	31.164	ng/ul	-0.01
54) 4-	Nitrophenol-d4	15.053	143	28535	26.960	ng/ul	0.00
60) Fl	uorene-d10	15.817	176	185251	31.461	ng/ul	0.00
65) 4,	6-Dinitro-2-methylph	15.958	200	28723	25.114	ng/ul	0.00
73) An	thracene-d10	17.673	188	270900	30.561	ng/ul	0.00
81) Py	rene-d10	19.953	212	305544	32.574	ng/ul	0.00
92) Bei	nzo(a)pyrene-d12	25.041	264	255012	31.500	ng/ul	0.00
Target (	Compounds					Qv	alue
2) 1,4	4-Dioxane	3.566	88	9295m	> 10.275	-	. 11/10/3/1
5) Pyi	ridine	3.978	79	62195	25.393	_	98
6) Bei	nzaldehyde	7.321	77	54447	31.033	ng/ul	94
8) Phe		7.379	94	88117	30.875		96
10) Bis	s(2-Chloroethyl)ether	7.597	93	65013	30.110	-	97
	Chlorophenol	7.755	128	62334	30.834	_	97
13) 2-1	Methylphenol	8.637	108	63628	29.930	ng/ul	95
14) 2,2	2'-oxybis(1-Chloropr	8.707	45	97580	31.318	ng/ul	98
16) Ace	etophenone	9.019	105	102047	29.676		96
17) N-N	Nitroso-di-n-propyla	8.989	70	60539	30.636	ng/ul	97
18) 4-N	Methylphenol	8.972	108	69948	30.771		94
19) Hex	xachloroethane	9.271	117	26070	30.530	-	96
22) Nit	trobenzene	9.412	77	88497	31.349	ng/ul	99
23) Iso	ophorone	9.929	82	166510	30.360	ng/ul	98
	Nitrophenol	10.123	139	37581	29.875		97
26) 2,4	4-Dimethylphenol	10.176	107	74848	29.103	ng/ul	99
27) Bis	s(2-Chloroethoxy)met	10.405	93	90949	30.038		99
29) 2,4	4-Dichlorophenol	10.670	162	61738	30.438	ng/ul	94
	hthalene	11.069	128	206467	29.752	ng/ul	98
30) Nap	JIICHUICHC					-	
30) Nap	Chloroaniline	11.181	127	82576	27.281	ng/ul	98
30) Nap 32) 4-0		11.181 11.328	127 225	82576 40926	27.281 29.252		97
30) Nap 32) 4-0 33) Hex 34) Cap	Chloroaniline				29.252	ng/ul	97

Data File : BG051383.D

Acq On : 7 Dec 2021 14:32

Operator : CG/JU Sample : PB141156BS

Misc :

ALS Vial : 41 Sample Multiplier: 1

Quant Time: Dec 08 02:25:18 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 ClientSampleId : SLCS156

## 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)
36) 2-Methylnaphthalene	12.661	142	144036	30.515 ng/ul	99
37) 1-Methylnaphthalene	12.879	142	143187	29.485 ng/ul	98
39) 1,2,4,5-Tetrachloroben	13.026	216	82836	31.049 ng/ul	94
40) Hexachlorocyclopentadiene	12.990	237	4898	4.542 ng/ul	88
41) 2,4,6-Trichlorophenol	13.273	196	51927	31.016 ng/ul	99
42) 2,4,5-Trichlorophenol	13.355	196	52495	29.942 ng/ul	96
43) 1,1'-Biphenyl	13.654	154	194058	30.573 ng/ul	96
44) 2-Chloronaphthalene	13.707	162	153810	30.463 ng/ul	100
45) 2-Nitroaniline	13.919	65	59102	33.822 ng/ul	94
47) Dimethylphthalate	14.265	163	198392	29.975 ng/ul	100
48) 2,6-Dinitrotoluene	14.406	165	43297	31.143 ng/ul	95
50) Acenaphthylene	14.553	152	248176	30.465 ng/ul	99
51) 3-Nitroaniline	14.741	138	45611	33.190 ng/ul	93
52) Acenaphthene	14.888	153	165064	30.724 ng/ul	95
53) 2,4-Dinitrophenol	14.970	184	12347	16.067 ng/ul	90
55) 4-Nitrophenol	15.065	109	25597	27.878 ng/ul	96
56) Dibenzofuran	15.223	168	233946	30.190 ng/ul	97
57) 2,4-Dinitrotoluene	15.200	165	61733	31.089 ng/ul#	97
58 <del>) 2,</del> 3,4,6-Tetrachlorophenol	15.452	232	39233	28.497 ng/ul	98
59) Diethylphthalate	15.617	149	211451	30.436 ng/ul	99
61) Fluorene	15.869	166	190498	30.690 ng/ul	100
62) 4-Chlorophenyl-phenyle	15.852	204	100247	29.968 ng/ul	97
63) 4-Nitroaniline	15.905	138	46503	34.773 ng/ul	94
66) 4,6-Dinitro-2-methylph	15.969	198	27379	24.822 ng/ul#	97
67) N-Nitrosodiphenylamine	16.069	169	169671	31.977 ng/ul	98
68) 4-Bromophenyl-phenylether	16.751	248	62145	31.285 ng/ul	91
69) Hexachlorobenzene	16.880	284	63184	31.194 ng/ul	97
70) Atrazine	17.009	200	57264	25.679 ng/ul	99
71) Pentachlorophenol	17.238	266	15673	17.462 ng/ul	96
72) Phenanthrene	17.620	178	318211	31.095 ng/ul	100
74) Anthracene	17.708	178	311456	30.645 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.631	216	84832	31.379 ng/uL	94
76) Pentachlorobenzene	15.141	250	77807	30.889 ng/uL	100
77) Carbazole	17.985	167	281912	31.600 ng/ul	99
78) Di-n-butylphthalate	18.502	149	364363	31.676 ng/ul	99
80) Fluoranthene	19.624	202	376456	32.676 ng/ul	97
82) Pyrene	19.982	202	361345	32.064 ng/ul	97
83) Butylbenzylphthalate	20.840	149	151535	32.344 ng/ul	96
84) 3,3'-Dichlorobenzidine	21.763	252	113734	31.511 ng/ul	99
85) Benzo(a)anthracene	21.857	228	326212	31.025 ng/ul	98
86) Bis(2-ethylhexyl)phtha	21.710	149	213499	31.668 ng/ul	99
87) Chrysene	21.927	228	314870	31.172 ng/ul	100
89) Di-n-octyl phthalate	22.973	149	364109	33.156 ng/ul	100
90) Benzo(b)fluoranthene	24.189	252	322248	31.500 ng/ul	99
91) Benzo(k)fluoranthene 93) Benzo(a)pyrene	24.260	252 252	299550	31.204 ng/ul 31.734 ng/ul	98 97
94) Indeno(1,2,3-cd)pyrene	25.123 29.207	276	309708	•	1211/2174
95) Dibenzo(a,h)anthracene	29.267	278	288651	31.339 ng/ul > 31.154 ng/ul	97
96) Benzo(g,h,i)perylene	30.435	276	269185	29.295 ng/ul	97 97

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