Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc :

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

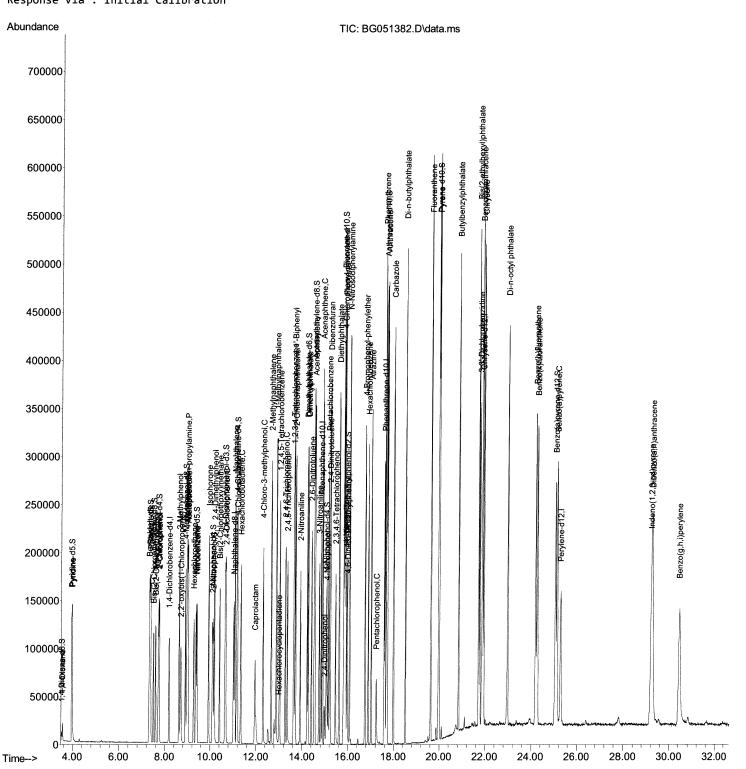
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleld : SLCS154

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 : BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc :

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

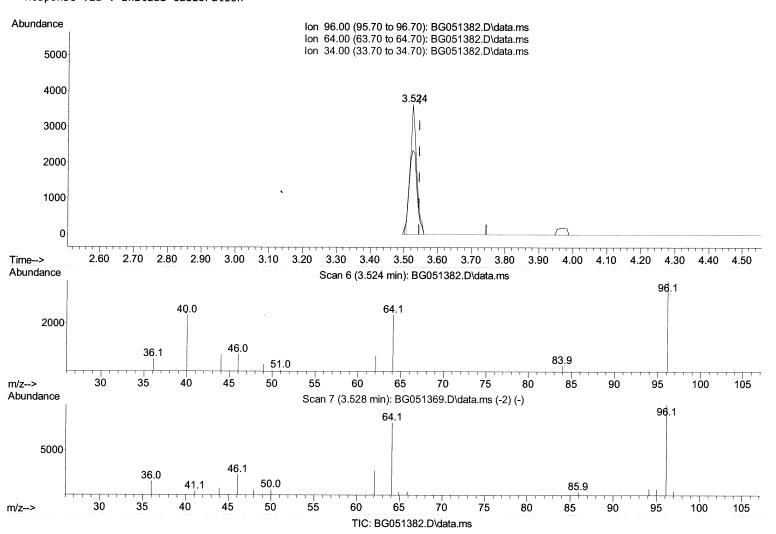
 $\label{lem:quant_method} {\tt Quant_Methods\SFAM-EPA-BG112321.M}$

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



(3) 1,4-Dioxane-d8 (S)

3.524min (-0.020) 5.62 ng/uL

response	4748	
Ion	Ехр%	Act%
96.00	100.00	100.00
64.00	77.60	64.41
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc :

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

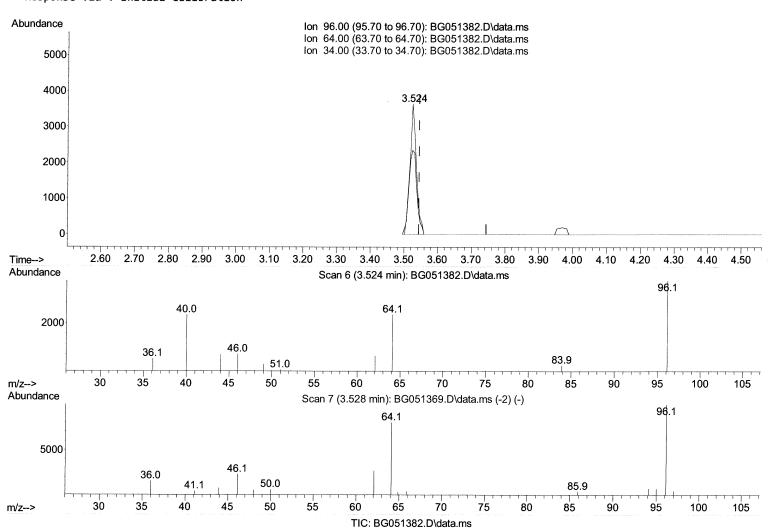
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument: BNA_G ClientSampleId: SLCS154

Manual IntegrationsAPPROVED

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



(3) 1,4-Dioxane-d8 (S)

3.524min (-0.020) 5.80 ng/uL m (2/10)JU

response	4901		
Ion	Ежр%	Act%	
96.00	100.00	100.00	
64.00	77.60	64.41	
34.00	0.00	0.00	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File: BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

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

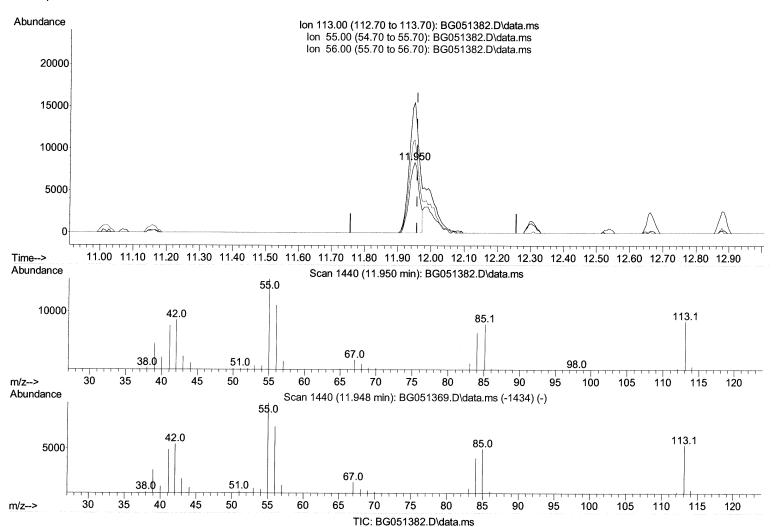
Quant Title : SVOA CALIBRATION

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

11.950min (-0.008) 22.01 ng/ul

response	17459				
Ion	Ехр%	Act%			
113.00	100.00	100.00			
55.00	183.80	185.40			
56.00	136.50	132.40			
0.00	0.00	0.00			

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

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

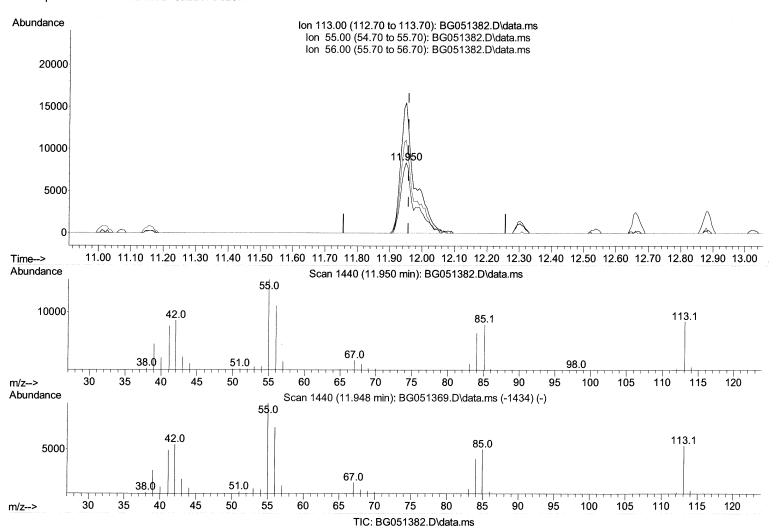
Quant Title : SVOA CALIBRATION

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

11.950min (-0.008) 31.44 ng/ul m \\(\frac{1}{2}\lambda \lambda \lambda

response	24935	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	183.80	185.40
56.00	136.50	132.40
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File: BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

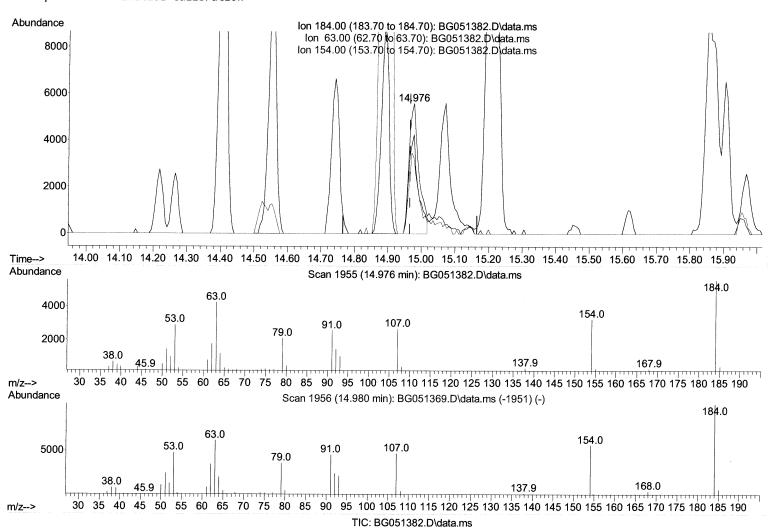
Quant Title : SVOA CALIBRATION

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



(53) 2,4-Dinitrophenol

14.976min (+ 0.010) 15.37 ng/ul

response	11517	
Ion	Ехр%	Act%
184.00	100.00	100.00
63.00	82.70	76.25
154.00	67.00	58.15
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

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

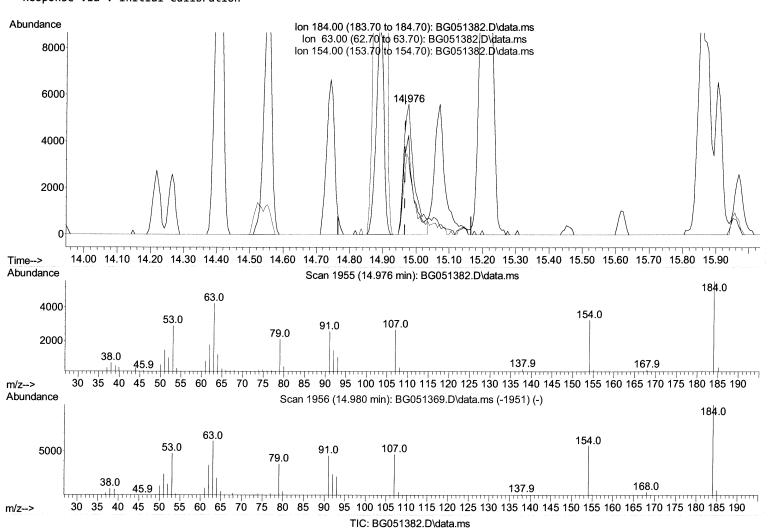
Quant Title : SVOA CALIBRATION

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



(53) 2,4-Dinitrophenol

14.976min (+ 0.010) 16.43 ng/ul m | 2/1(|2| JU

response	12311	
Ion	Exp%	Act%
184.00	100.00	100.00
63.00	82.70	76.25
154.00	67.00	58.15
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

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

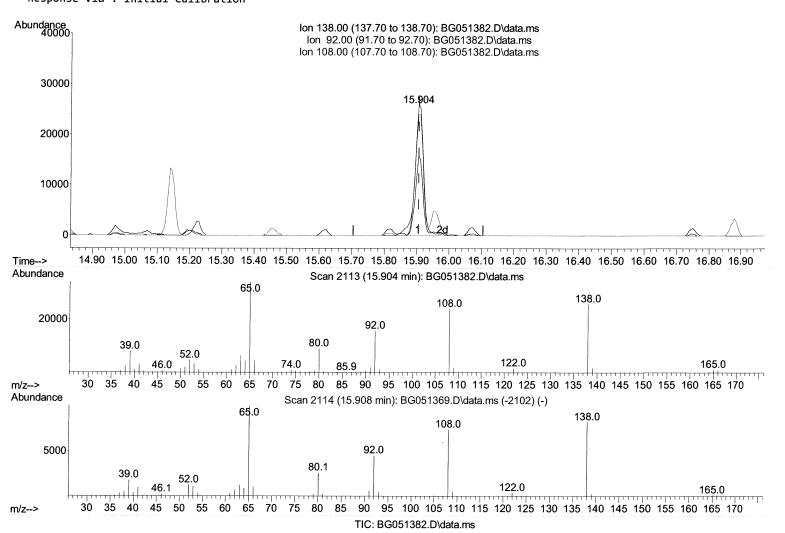
Quant Title : SVOA CALIBRATION

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



(63) 4-Nitroaniline

15.904min (-0.002) 36.85 ng/ul

response	48050	
Ion	Ехр%	Act%
138.00	100.00	100.00
92.00	61.60	60.80
108.00	90.70	92.72
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 2021

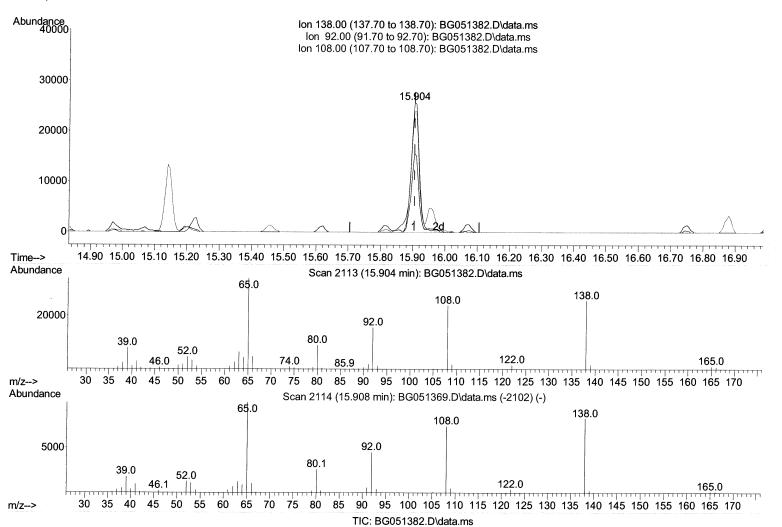
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument: BNA_G ClientSampleld: SLCS154

Manual IntegrationsAPPROVED

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



(63) 4-Nitroaniline

15.904min (-0.002) 36.93 ng/ul m (3/16/21JU

response	48160	
Ion	Ежр%	Act%
138.00	100.00	100.00
92.00	61.60	60.80
108.00	90.70	92.72
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051382.D

Acq On : 7 Dec 2021 13:51 Operator : CG/JU Sample : PB141154BS

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 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 : SLCS154

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)
Inte	rnal Standards						
1)	1,4-Dichlorobenzene-d4	8.189	152	29363	20.000	ng/ul	-0.01
20)	Naphthalene-d8	11.021	136	126856	20.000	-	0.00
38)	Acenaphthene-d10	14.823	164	82868	20.000		0.00
	Phenanthrene-d10	17.578	188	180047	20.000	ng/ul	0.00
79)	Chrysene-d12	21.879	240	151449	20.000	•	0.00
	Perylene-d12	25.281	264	150462	20.000	-	0.00
Syste	em Monitoring Compounds						116/21-
	1,4-Dioxane-d8	3.524	96	4901m ~	> 5.800	ng/uL>	-0.02/2/16/2/JC
	Pyridine-d5	3.953	84	70018	28.239	ng/ul	-0.03
7)	Phenol-d5	7.349	99	93200	32.115		0.00
9)	Bis-(2-Chloroethyl)eth	7.502	67	59014	32.379		-0.01
	2-Chlorophenol-d4	7.719	132	67726	32.409	_	-0.01
	4-Methylphenol-d8	8.906	113	73371	31.330		0.00
	Nitrobenzene-d5	9.370	128	34797	32.495	-	0.00
-	2-Nitrophenol-d4	10.093	143	39783	32.934	-	0.00
	2,4-Dichlorophenol-d3	10.645	165	67666	33.015		0.00
	4-Chloroaniline-d4	11.156	131	113011	37.685	•	0.00
	Dimethylphthalate-d6	14.218	166	205716	32.263	_	0.00
	Acenaphthylene-d8	14.523	160	265792	33.057		0.00
	4-Nitrophenol-d4	15.052	143	30446	29.499	-	0.00
	Fluorene-d10	15.816	176	188539	32.836		0.00
	4,6-Dinitro-2-methylph	15.957	200	31030	27.930	-	0.00
	Anthracene-d10	17.678	188	283114	32.878	•	0.00
•	Pyrene-d10	19.958	212	321333	35.065	_	0.00
	Benzo(a)pyrene-d12	25.052	264	270277	33.634	-	0.00
Targe	et Compounds					0va	lue
_	1,4-Dioxane	3.560	88	10206	10.710	-	94
	Pyridine	3.977	79	71492	27.710	_	95
	Benzaldehyde	7.320	77	66070	35.749		95
	Phenol	7.379	94	91340	30.382		99
	Bis(2-Chloroethyl)ether	7.596	93	68798	30.248	-	99
	2-Chlorophenol	7.755	128	64871	30.462	_	98
•	2-Methylphenol	8.642	108	67069	29.950	_	96
	2,2'-oxybis(1-Chloropr	8.712	45	101775	31.009		96
	Acetophenone	9.024	105	106359	29.362		97
	N-Nitroso-di-n-propyla	8.994	70	61346	29.471	-	97
	4-Methylphenol	8.971	108	71604	29.903	-	97
	Hexachloroethane	9.270	117	27377	30.436		96
	Nitrobenzene	9.411	77	90645	32.282		97
	Isophorone	9.929	82	167395	30.685		100
,	2-Nitrophenol	10.128	139	39007	31.176		96
	2,4-Dimethylphenol	10.128	107	76869	30.049	-	99
	Bis(2-Chloroethoxy)met	10.181	93	92508	30.717		
	2,4-Dichlorophenol				30.717	-	99
		10.675	162	62092		٠.	96
	Naphthalene	11.068	128	211740	30.676	_	98
	4-Chloroaniline	11.180	127	87304	28.999	•	99
	Hexachlorobutadiene	11.333	225	42048	30.216		99
	Caprolactam	11.950	113	24935m >	31.438		ionicus 27
35)	4-Chloro-3-methylphenol	12.302	107	77016	31.778	ng/ul	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\

Data File : BG051382.D

Acq On : 7 Dec 2021 13:51

Operator : CG/JU Sample : PB141154BS

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Dec 08 02:25:07 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: SLCS154

Manual IntegrationsAPPROVED

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

Compound	R.T.	QIon	Response	Conc Units De	v(Min)
36) 2-Methylnaphthalene	12.661	142	140389	29.902 ng/ul	100
37) 1-Methylnaphthalene	12.878	142	146083	30.243 ng/ul	95
39) 1,2,4,5-Tetrachloroben	13.025	216	82849	31.846 ng/ul	96
40) Hexachlorocyclopentadiene	12.990	237	5266	5.008 ng/ul	96
41) 2,4,6-Trichlorophenol	13.272	196	51729	31.686 ng/ul	96
42) 2,4,5-Trichlorophenol	13.354	196	55387	32.397 ng/ul	98
43) 1,1'-Biphenyl	13.659	154	191871	31.000 ng/ul	96
44) 2-Chloronaphthalene	13.712	162	153152	31.106 ng/ul	100
45) 2-Nitroaniline	13.918	65	57587	33.796 ng/ul	94
47) Dimethylphthalate	14.265	163	196996	30.523 ng/ul	99
48) 2,6-Dinitrotoluene	14.406	165	43752	32.273 ng/ul	96
50) Acenaphthylene	14.553	152	244493	30.778 ng/ul	99
51) 3-Nitroaniline	14.741	138	45647	34.064 ng/ul	93
52) Acenaphthene	14.887	153	162764	31.069 ng/ul	97
53) 2,4-Dinitrophenol	14.976	184	12311m>	16.429 ng/ul:	• •
55) 4-Nitrophenol	15.070	109	25377	28.344 ng/ul	96
56) Dibenzofuran	15.222	168	229719	30.400 ng/ul	100
57) 2,4-Dinitrotoluene	15.199	165	61650	31.839 ng/ul	99
58) 2,3,4,6-Tetrachlorophenol	15.457	232	39530	29.445 ng/ul	98
59) Diethylphthalate	15.616	149	211693	31.248 ng/ul	99
61) Fluorene	15.874	166	185616	30.666 ng/ul	97
62) 4-Chlorophenyl-phenyle	15.851	204	98915	30.324 ng/ul	> 13/16/13/27
63) 4-Nitroaniline	15.904	138	48160m>	36.931 ng/ul	> (3/16/2)
66) 4,6-Dinitro-2-methylph	15.974	198	27929	26.066 ng/ul#	
67) N-Nitrosodiphenylamine68) 4-Bromophenyl-phenylether	16.074	169	167295	32.457 ng/ul	98
69) Hexachlorobenzene	16.750 16.879	248 284	61792 62640	32.022 ng/ul	95 00
70) Atrazine	17.014	200	65197	31.835 ng/ul	99
71) Pentachlorophenol	17.238	266	15793	30.097 ng/ul 18.114 ng/ul	99 9 1
72) Phenanthrene	17.620	178	319100	32.099 ng/ul	100
74) Anthracene	17.714	178	310587	31.458 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.630	216	85986	32.742 ng/uL	98
76) Pentachlorobenzene	15.146	250	75694	30.934 ng/uL	97
77) Carbazole	17.984	167	285557	32.951 ng/ul	98
78) Di-n-butylphthalate	18.507	149	369330	33.052 ng/ul	99
80) Fluoranthene	19.623	202	373432	33.178 ng/ul	98
82) Pyrene	19.987	202	368347	33.456 ng/ul	96
83) Butylbenzylphthalate	20.845	149	154489	33.752 ng/ul	94
84) 3,3'-Dichlorobenzidine	21.762	252	116617	33.072 ng/ul	99
85) Benzo(a)anthracene	21.856	228	330551	32.179 ng/ul	100
86) Bis(2-ethylhexyl)phtha	21.715	149	217789	33.066 ng/ul	100
87) Chrysene	21.926	228	314082	31.828 ng/ul	99
89) Di-n-octyl phthalate	22.978	149	364843	33.470 ng/ul	100
<pre>90) Benzo(b)fluoranthene</pre>	24.194	252	322573	31.768 ng/ul	99
<pre>91) Benzo(k)fluoranthene</pre>	24.265	252	302983	31.797 ng/ul	99
93) Benzo(a)pyrene	25.128	252	311225	32.127 ng/ul	98
94) Indeno(1,2,3-cd)pyrene	29.218	276	344268	31.758 ng/ul	98
95) Dibenzo(a,h)anthracene	29.265	278	294646	32.039 ng/ul	98
96) Benzo(g,h,i)perylene	30.452	276	273602	29.998 ng/ul	94

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