Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051311.D

Acq On : 2 Dec 2021 19:45

Operator : CG/JU Sample : PB141107BS

Misc

ALS Vial : 16 Sample Multiplier: 1

Quant Time: Dec 03 00:02:12 2021

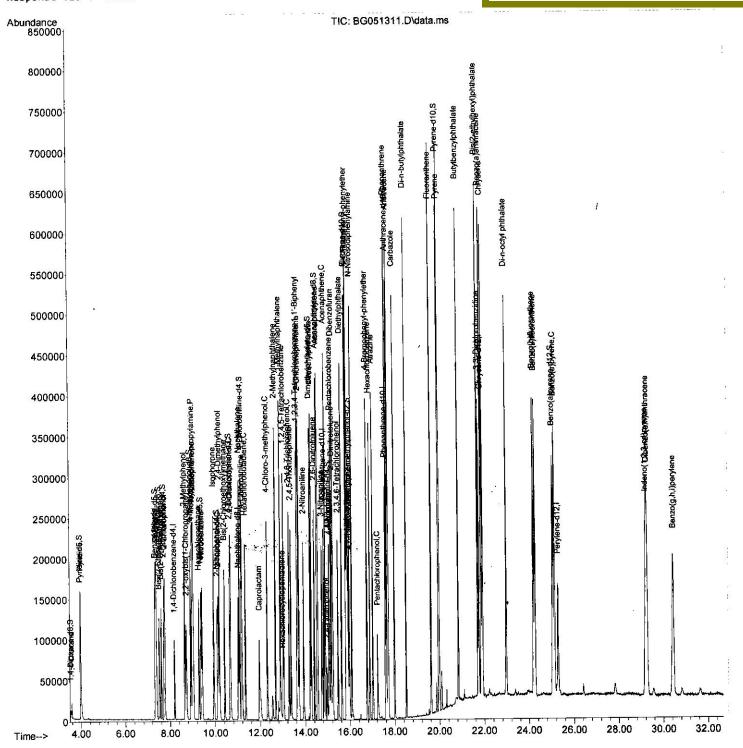
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
Response via : Initial Calibration

Instrument :
BNA\_G
ClientSampleId :

# **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



SFAM-EPA-BG112321.M Fri Dec 03 00:24:00 2021

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051311.D

Acq On : 2 Dec 2021 19:45

Operator : CG/JU Sample : PB141107BS

Misc

ALS Vial : 16 Sample Multiplier: 1

Quant Time: Dec 03 00:02:12 2021

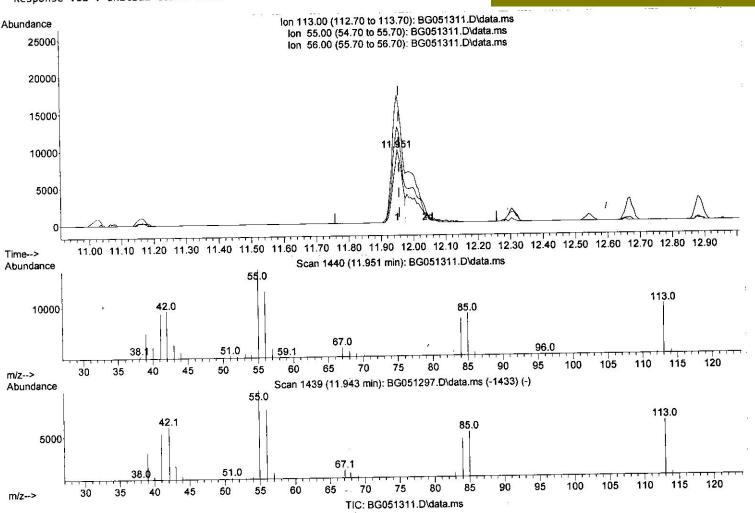
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
Response via : Initial Calibration

Instrument : BNA\_G ClientSampleId :

# Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



### (34) Caprolactam

11.951min (-0.006) 24.51 ng/ul

response	19062	
Ion	Exp&	Act%
113.00	100.00	100.00
55.00	183.80	174.44
56.00	136.50	131.77
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051311.D

Acq On : 2 Dec 2021 19:45

Operator : CG/JU Sample : PB141107BS

Misc

ALS Vial : 16 Sample Multiplier: 1

Quant Time: Dec 03 00:02:12 2021

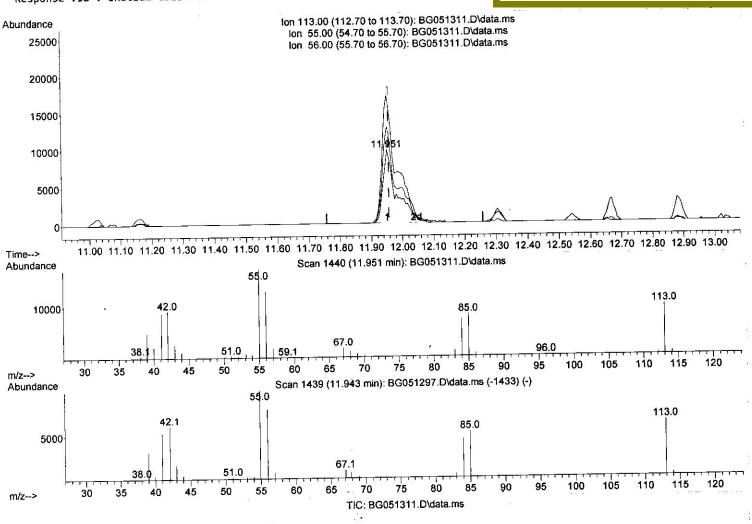
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
Response via : Initial Calibration

Instrument :
BNA\_G
ClientSampleId :

# Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



(34) Caprolactam

11.951min (-0.006) 37.91 ng/ul m \ \( \frac{1}{12} \)

response	29487	
Ion	Ежр%	Act%
113.00	100.00	100.00
55.00	183.80	174.44
56.00	136.50	131.77
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051311.D

Acq On : 2 Dec 2021 19:45

Operator : CG/JU Sample : PB141107BS

Misc

ALS Vial : 16 Sample Multiplier: 1

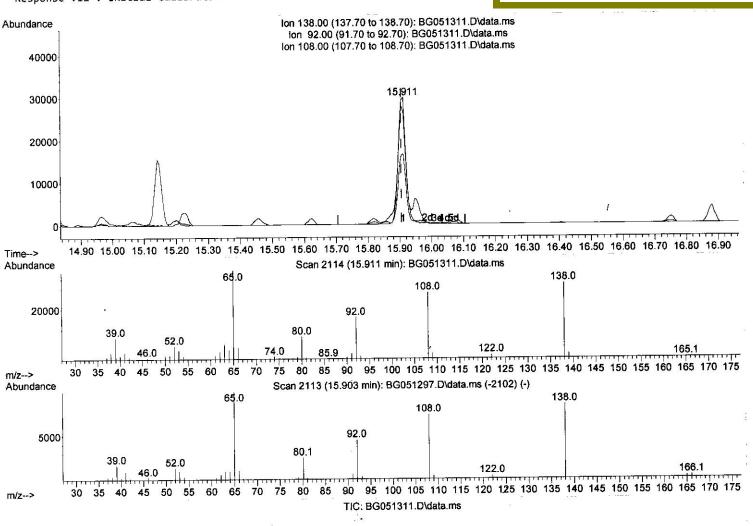
Quant Time: Dec 03 00:02:12 2021

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
ClientSampleId :

#### Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



#### (63) 4-Nitroaniline

15.911min (+ 0.006) 41.87 ng/ul

response	54396	
Ion	Exp%	Act%
138.00	100.00	100.00
92.00	61.60	55.60
108.00	90.70	88.10
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051311.D

Acq On : 2 Dec 2021 19:45

Operator : CG/JU Sample : PB141107BS

Misc

ALS Vial : 16 Sample Multiplier: 1

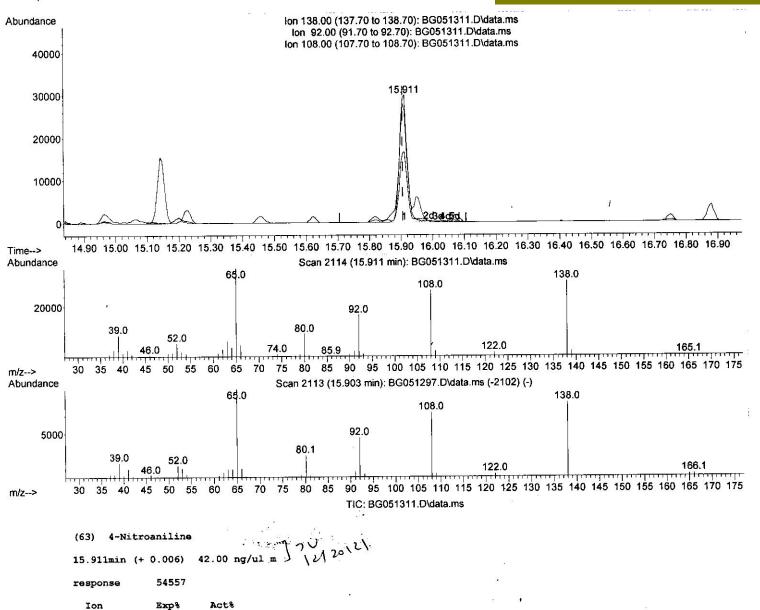
Quant Time: Dec 03 00:02:12 2021

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 ClientSampleId :

#### Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/03/2021 Supervised By: mohammad ahmed 12/05/2021



100.00

61.60

90.70

0.00

100.00

55.60

88.10

0.00

138.00

92.00

108.00

0.00

# Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051311.D

Acq On : 2 Dec 2021 19:45

Operator : CG/JU Sample : PB141107BS

Misc

ALS Vial : 16 Sample Multiplier: 1

Quant Time: Dec 03 00:02:12 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

nstrument :
NA_G
lientSampleId :
LCC107

# **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021

Compound			Response			
Internal Standards		<b>-</b>				
<ol> <li>1,4-Dichlorobenzene-d4</li> </ol>	8.191	152	27973	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.023	136	124405	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.830	164	82549	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.580	188	179425	20.000	ng/ul	0.00
79) Chrysene-d12	21.881	240	149239	20.000	ng/ul	0.00
88) Perylene-d12	25.283	264	148922	20.000	ng/ul	0.00
ystem Monitoring Compounds						
3) 1,4-Dioxane-d8	3.532	96	5323	6.613	ng/uL	-0.01
4) Pyridine-d5	3.961	84	75110	31.798	22 (32)	-0.02
7) Phenol-d5	7.357	99	102798	37.183	F1.200	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.509	67	62705	36.113		0.00
11) 2-Chlorophenol-d4	7.727		73848	37.094		0.00
15) 4-Methylphenol-d8	8.914	113	82111	36.805		0.00
21) Nitrobenzene-d5	9.372	128	38652	36.806		0.00
24) 2-Nitrophenol-d4	10.101	143	44303	37.398		0.00
28) 2,4-Dichlorophenol-d3	10.647		75490	37.559		0.00
31) 4-Chloroaniline-d4	11.164		143197	48.691		0.00
46) Dimethylphthalate-d6	14.219		234692	36.950		0.00
49) Acenaphthylene-d8	14.525		299240	37.361		0.00
54) 4-Nitrophenol-d4	15.054		35065	34.106	3.40	0.00
60) Fluorene-d10	15.817		207987	36.363		0.00
65) 4,6-Dinitro-2-methylph	15.953		35634	32.185		
73) Anthracene-d10	17,680		319597	37.244		
81) Pyrene-d10	19.960		358054	39.651	-	0.00
92) Benzo(a)pyrene-d12	25.048		297686	37.428		
Target Compounds					0	value
2) 1,4-Dioxane	3.573	88	10784	11.879		93
5) Pyridine	3.984		78602			96
6) Benzaldehyde	7.327		64749	36.776		
8) Phenol	7.386		107546	37.550		
10) Bis(2-Chloroethyl)ether	7.603		78349			
12) 2-Chlorophenol	7.756		75425			
13) 2-Methylphenol	8.643	108		37.575		
14) 2,2'-oxybis(1-Chloropr	8.720		117738	37.655		
16) Acetophenone				36.284		
17) N-Nitroso-di-n-propyla	9.002		74290	37.463		
18) 4-Methylphenol	8.972		85282	37.385		
19) Hexachloroethane	9.278		30475	35.564		
22) Nitrobenzene	9.419		103947	37.749		
23) Isophorone	9.936		201968	37.752		
25) 2-Nitrophenol	10.130		47163	38.437		
26) 2,4-Dimethylphenol	10.183		94742	37.766		
27) Bis(2-Chloroethoxy)met	10.412		112222	37.998		
29) 2,4-Dichlorophenol	10.412		75378	38.098		
30) Naphthalene	11.076		252519	37.305		
32) 4-Chloroaniline	11.188		100563	34.061		
33) Hexachlorobutadiene	11.334		48267			
34) Caprolactam	11.334		29487m			
35) 4-Chloro-3-methylphenol	12.304		92587	38.956		
33) 4-Chioro-3-methylphenol	12.304	10/	32301	30.220	⊞\ u1	70

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051311.D

Acq On : 2 Dec 2021 19:45

Operator : CG/JU : PB141107BS Sample

Misc

ALS Vial : 16 Sample Multiplier: 1

Quant Time: Dec 03 00:02:12 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

Compound	R.T.	QIon	Response	Conc Units Dev(	Min)
36) 2-Methylnaphthalene	12.668	142	173087	37.593 ng/ul	99
37) 1-Methylnaphthalene	12.886	142	172807	36.481 ng/ul	98
39) 1,2,4,5-Tetrachloroben	13.027	216	96879	37.383 ng/ul	96
40) Hexachlorocyclopentadiene	12.991	237	13727	13.105 ng/ul#	96
41) 2,4,6-Trichlorophenol	13.273	196	63131	38.819 ng/ul	100
42) 2,4,5-Trichlorophenol	13.356	196	64827	38.065 ng/ul	97
43) 1,1'-Biphenyl	13.661	154	232428	37.698 ng/ul	97
44) 2-Chloronaphthalene	13.714	162	182902	37.292 ng/ul	99
45) 2-Nitroaniline	13.920	65	68228	40.195 ng/ul	94
47) Dimethylphthalate	14.266	163	236372	36.766 ng/ul	100
48) 2,6-Dinitrotoluene	14.407	165	51739	38.312 ng/ul	93
50) Acenaphthylene	14.554	152	298709	37.749 ng/ul	98
51) 3-Nitroaniline	14.742	138	54011	40.461 ng/ul	97
52) Acenaphthene	14.889	153	199386	38.206 ng/ul	97
53) 2,4-Dinitrophenol	14.965	184	16540	22.158 ng/ul	86
55) 4-Nitrophenol	15.065	109	30682	34.401 ng/ul	94
56) Dibenzofuran	15.224	168	280264	37.233 ng/ul	100
57) 2,4-Dinitrotoluene	15.200	165	72779	37.732 ng/ul	94
58) 2,3,4,6-Tetrachlorophenol	15.459	232	50075	37.444 ng/ul	96
59) Diethylphthalate	15.624	149	248191	36.777 ng/ul	98
61) Fluorene	<b>1</b> 5.876	166	221393	36.719 ng/ul	100
62) 4-Chlorophenyl-phenyle	15.859	204	117070 1	36.029 ng/ul	96
63) 4-Nitroaniline	15.911	138	54557m	41.998 ng/ul	
66) 4,6-Dinitro-2-methylph	15.970		35455 ~	33,205 ng/ul#	93
67) N-Nitrosodiphenylamine	16.076		199955	38,927 ng/ul	98
68) 4-Bromophenyl-phenylether	16.752		73287	38,110 ng/ul	94
69) Hexachlorobenzene	16.881		75354	38.429 ng/ul	95
70) Atrazine	17.016		70915	32.850 ng/ul	99
71) Pentachlorophenol	17.233		21006	24.176 ng/ul	97
72) Phenanthrene	17.621	178	379066	38.263 ng/ul	100
74) Anthracene	17.715		370886	37.696 ng/ul	98
75) 1,2,3,4-Tetrachloroben	13.632	216	100448	38.381 ng/uL	99
76) Pentachlorobenzene	15.148	250	89992	36.904 ng/uL	98
77) Carbazole	17.985		334708	38.756 ng/ul	99
78) Di-n-butylphthalate	18.508		429157	38.539 ng/ul	99
80) Fluoranthene	19.625		441604	39.816 ng/ul	98
82) Pyrene	19.989		427788	39.430 ng/ul	97
83) Butylbenzylphthalate	20.847	• 0000000000000000000000000000000000000		39.413 ng/ul	95
<pre>84) 3,3'-Dichlorobenzidine</pre>	21.763	*	128391	36.951 ng/ul	99
<pre>85) Benzo(a)anthracene</pre>	21.857		384766	38.012 ng/ul	99
<pre>86) Bis(2-ethylhexyl)phtha</pre>	21.716		251155	38.696 ng/ul	99 .
87) Chrysene	21.928		370163 <sup>-</sup>	38.066 ng/ul	100
89) Di-n-octyl phthalate	22.980		428969	39.760 ng/ul	100
90) Benzo(b)fluoranthene	24.196		386480	38.455 ng/ul	100
91) Benzo(k)fluoranthene	24,266		352429	37.368 ng/ul	99
93) Benzo(a)pyrene	25.124		366148	38.188 ng/ul	98
94) Indeno(1,2,3-cd)pyrene	29.202		411978	38.397 ng/ul	97
95) Dibenzo(a,h)anthracene	29.255		349246	38.368 ng/ul	98
96) Benzo(g,h,i)perylene	30.441	. 276	348245	38.577 ng/ul	98

# **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By:mohammad ahmed 12/05/2021

72/20121

Instrument: BNA\_G ClientSampleId: SLCS107

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