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

Data File : BG051192.D

Acq On : 23 Nov 2021 21:20

Operator : CG/JU Sample : PB140910BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 06:56:05 2021

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

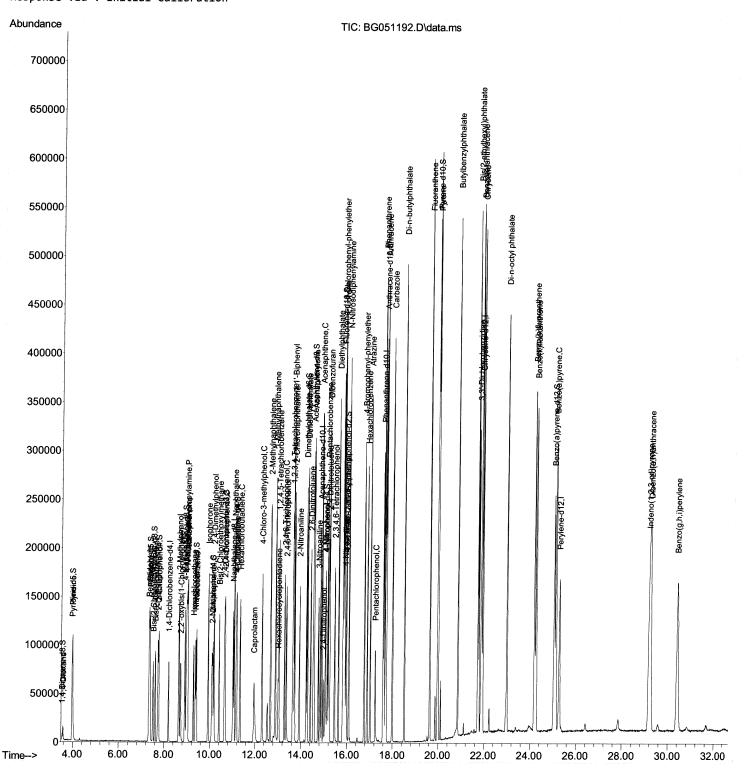
Quant Title : SVOA CALIBRATION

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



# **Manual IntegrationsAPPROVED**

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



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

Data File : BG051192.D

Acq On : 23 Nov 2021 21:20

Operator : CG/JU Sample : PB140910BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 06:56:05 2021

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

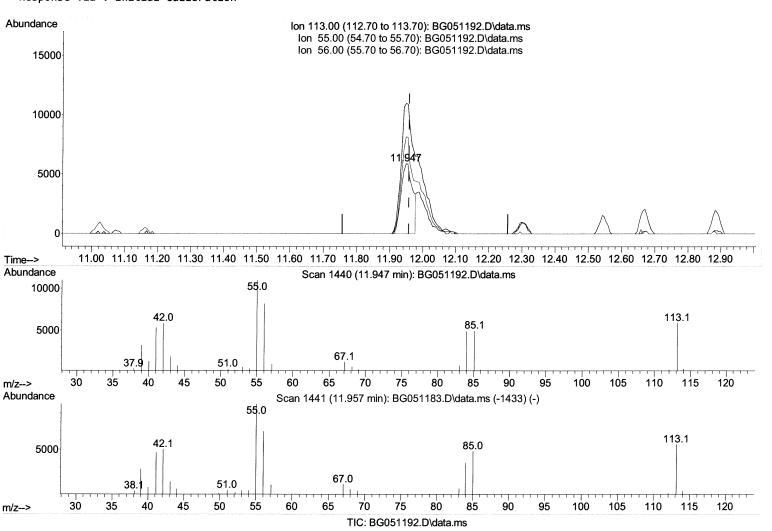
Quant Title : SVOA CALIBRATION

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



# **Manual IntegrationsAPPROVED**

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



### (34) Caprolactam

11.947min (-0.010) 20.65 ng/ul

response	14378				
Ion	Ехр%	Act%			
113.00	100.00	100.00			
55.00	183.80	186.15			
56.00	136.50	138.44			
0.00	0.00	0.00			

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

Data File : BG051192.D

Acq On : 23 Nov 2021 21:20

Operator : CG/JU Sample : PB140910BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 06:56:05 2021

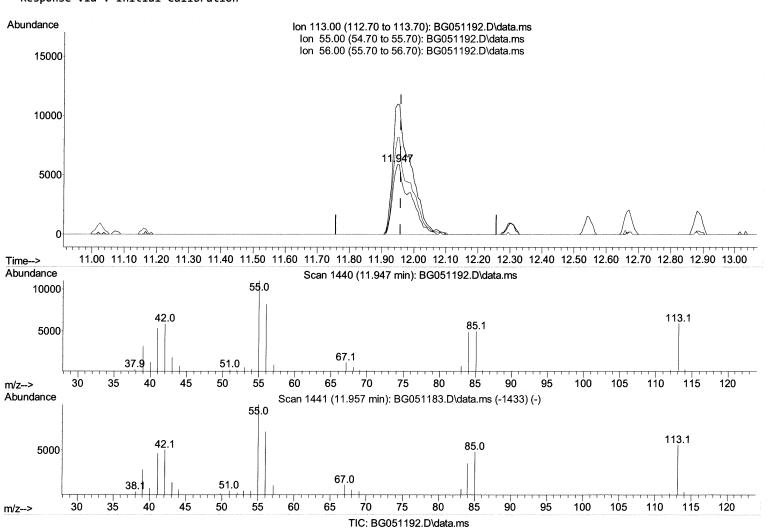
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument: BNA\_G ClientSampleld: SLCS910

## **Manual IntegrationsAPPROVED**

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



### (34) Caprolactam

11.947min (-0.010) 31.58 ng/ul m (1/24/2/14

response	21985	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	183.80	186.15
56.00	136.50	138.44
0.00	0.00	0.00

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

Data File : BG051192.D

Acq On : 23 Nov 2021 21:20

Operator : CG/JU Sample : PB140910BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 06:56:05 2021

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

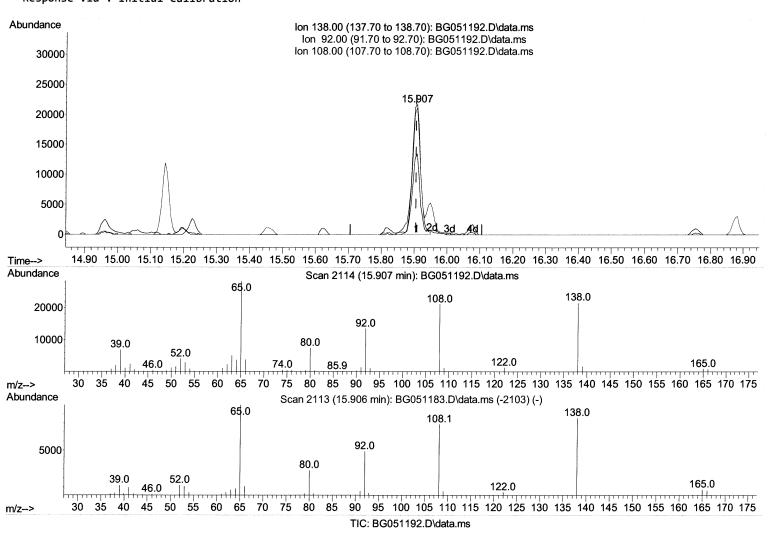
Quant Title : SVOA CALIBRATION

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



## **Manual IntegrationsAPPROVED**

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



### (63) 4-Nitroaniline

15.907min (+ 0.002) 31.91 ng/ul

response	40120				
Ion	Ехр%	Act%			
138.00	100.00	100.00			
92.00	61.60	62.19			
108.00	90.70	97.57			
0.00	0.00	0.00			

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

Data File: BG051192.D

Acg On : 23 Nov 2021 21:20

Operator : CG/JU Sample : PB140910BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 06:56:05 2021

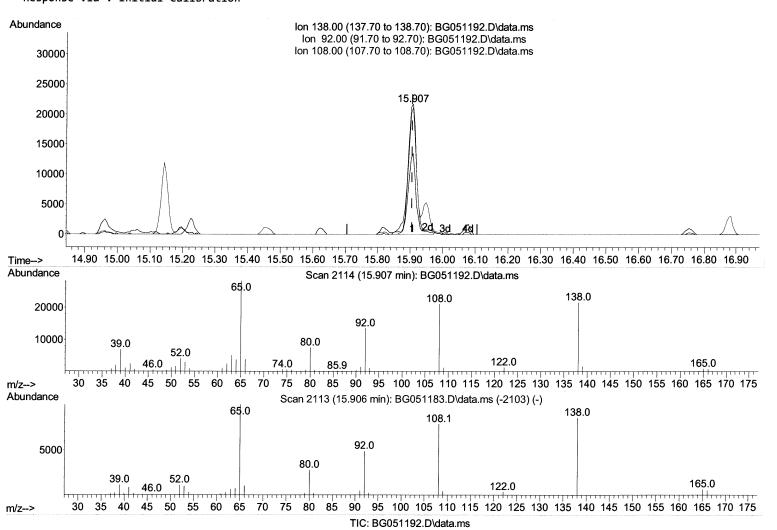
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument : BNA\_G ClientSampleId : SLCS910

## **Manual Integrations APPROVED**

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



## (63) 4-Nitroaniline

response	40572	
Ion	Ехр%	Act%
138.00	100.00	100.00
92.00	61.60	62.19
108.00	90.70	97.57
0.00	0.00	0.00

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

Data File : BG051192.D

Acq On : 23 Nov 2021 21:20

Operator : CG/JU Sample : PB140910BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 06:56:05 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 : SLCS910

# Manual IntegrationsAPPROVED

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

Compound	R.T.	QIon	Response	Conc Unit	ts Dev(	Min)
Internal Standards						
<ol> <li>1,4-Dichlorobenzene-d4</li> </ol>	8.199	152	24021	20.000 r	ng/ul	0.00
20) Naphthalene-d8	11.025	136	111356	20.000 r		0.00
38) Acenaphthene-d10	14.832	164	79891	20.000 r		0.00
64) Phenanthrene-d10	17.582	188	186837	20.000 r	ng/ul	0.00
79) Chrysene-d12	21.883	240	166683	20.000 r	ng/ul	0.00
88) Perylene-d12	25.279	264	172130	20.000 r	ng/ul	0.00
System Monitoring Compounds						
<ol><li>3) 1,4-Dioxane-d8</li></ol>	3.540	96	3790	5.483 r	ng/uL	0.00
4) Pyridine-d5	3.969	84	50990	25.139 r	ng/ul	-0.01
7) Phenol-d5	7.353	99	67397	28.389 n		0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.511	67	41400	27.766 n	-	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.729	132	48239	28.217 n	-	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.910	113	53071	27.702 n	g/ul	0.00
21) Nitrobenzene-d5	9.374	128	25723	27.365 n		0.00
24) 2-Nitrophenol-d4	10.102	143	29644	27.956 n	-	0.00
28) 2,4-Dichlorophenol-d3	10.649	165	50211	27.909 n	g/ul	0.00
31) 4-Chloroaniline-d4	11.166	131	65402	24.845 n	g/ul	0.00
46) Dimethylphthalate-d6	14.221	166	176110	28.649 n		0.00
49) Acenaphthylene-d8	14.527	160	212792	27.452 n	-	0.00
54) 4-Nitrophenol-d4	15.044	143	29682	29.831 n		0.00
60) Fluorene-d10	15.819	176	153756	27.776 n		0.00
65) 4,6-Dinitro-2-methylph	15.949	200	32532	28.217 n		0.00
73) Anthracene-d10	17.682	188	251175	28.109 n		0.00
81) Pyrene-d10	19.956	212	297422	29.490 n	_	0.00
92) Benzo(a)pyrene-d12	25.044	264	263236	28.635 n		0.00
arget Compounds					Qva:	lue
2) 1,4-Dioxane	3.575	88	7981	10.238 n	•	96
5) Pyridine	3.986	79	55679	26.380 n		98
6) Benzaldehyde	7.329	77	43149	28.539 n	-	94
8) Phenol	7.382	94	71579	29.104 n		98
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.605	93	53851	28.942 n	-	96
12) 2-Chlorophenol	7.758	128	51081	29.321 n		96
13) 2-Methylphenol	8.645	108	53735	29.332 n	_	98
14) 2,2'-oxybis(1-Chloropr	8.716	45	78826	29.358 n		98
16) Acetophenone	9.027	105	86469	29.180 n	g/ul	96
17) N-Nitroso-di-n-propyla	9.004	70	50713	29.781 n		96
18) 4-Methylphenol	8.974	108	58103	29.661 n	_	94
19) Hexachloroethane	9.286	117	20496	27.854 n		92
22) Nitrobenzene	9.415	77	72402	29.374 ng		95
23) Isophorone	9.932	82	141168	29.480 ng		99
25) 2-Nitrophenol	10.132	139	31675	28.840 ng		99
26) 2,4-Dimethylphenol	10.185	107	66726	29.715 ng		96
27) Bis(2-Chloroethoxy)met	10.414	93	77839	29.444 ng	-	99
29) 2,4-Dichlorophenol	10.672	162	51700	29.193 ng	•	95
30) Naphthalene	11.078	128	171658	28.331 ng		98
32) 4-Chloroaniline	11.189	127	71245	26.958 ng		95
33) Hexachlorobutadiene	11.342	225	32434	26.552 ng		95
34) Caprolactam	11.947	113	21985m >	-		iilaalal
35) 4-Chloro-3-methylphenol	12.300	107	65257	30.674 ng		97

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

Data File : BG051192.D

Acq On : 23 Nov 2021 21:20 Operator : CG/JU Sample : PB140910BS

Misc

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 24 06:56:05 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 : SLCS910

# **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	12.670	142	118694	28.800 ng/ul	98
37) 1-Methylnaphthalene	12.887	142	120882	28.509 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.028	216	69118	27.558 ng/ul	98
<ol><li>40) Hexachlorocyclopentadiene</li></ol>	12.999	237	17398	17.162 ng/ul#	98
<pre>41) 2,4,6-Trichlorophenol</pre>	13.275	196	45307	28.786 ng/ul	98
42) 2,4,5-Trichlorophenol	13.352	196	48432	29.385 ng/ul	96
43) 1,1'-Biphenyl	13.663	154	166729	27.942 ng/ul	97
44) 2-Chloronaphthalene	13.716	162	132585	27.933 ng/ul	99
45) 2-Nitroaniline	13.922	65	49474	30.116 ng/ul	91
47) Dimethylphthalate	14.268	163	184001	29.572 ng/ul	100
48) 2,6-Dinitrotoluene	14.409	165	39569	30.275 ng/ul	91
50) Acenaphthylene	14.556	152	218202	28.492 ng/ul	99
51) 3-Nitroaniline	14.744	138	37698	29.180 ng/ul#	98
52) Acenaphthene	14.891	153	143466	28.406 ng/ul	97
53) 2,4-Dinitrophenol	14.961	184	17457	24.164 ng/ul	87
55) 4-Nitrophenol	15.056	109	27474	31.829 ng/ul	94
56) Dibenzofuran	15.226	168	209151	28.710 ng/ul	99
57) 2,4-Dinitrotoluene	15.197	165	58608	31.396 ng/ul#	99
58) 2,3,4,6-Tetrachlorophenol	15.455	232	40893	31.595 ng/ul	97
59) Diethylphthalate	15.625	149	200388	30.682 ng/ul	100
61) Fluorene	15.878	166	169307	29.014 ng/ul	100
62) 4-Chlorophenyl-phenyle	15.860	204	90904	28.907 ng/ul	95 1001010
63) 4-Nitroaniline	15.907	138	40572m>	32.271 ng/ul	
66) 4,6-Dinitro-2-methylph	15.966	198	33068	29.741 ng/ul#	95
67) N-Nitrosodiphenylamine	16.078	169	155367	29.047 ng/ul	97
68) 4-Bromophenyl-phenylether	16.754	248	58184	29.056 ng/ul	95
69) Hexachlorobenzene 70) Atrazine	16.883	284	60024	29.397 ng/ul	95
71) Pentachlorophenol	17.018	200 266	66335	29.509 ng/ul	99
72) Phenanthrene	17.235	178	22328	24.678 ng/ul	97
74) Anthracene	17.623 17.717	178	305654 305661	29.629 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.634	216	71928	29.834 ng/ul	98
76) Pentachlorobenzene	15.144	250	67055	26.393 ng/uL 26.407 ng/uL	96 99
77) Carbazole	17.987	167	279422	31.071 ng/ul	
78) Di-n-butylphthalate	18.510	149	358597	30.925 ng/ul	100 99
80) Fluoranthene	19.627	202	382521	30.880 ng/ul	96
82) Pyrene	19.985	202	369858	30.523 ng/ul	99
83) Butylbenzylphthalate	20.849	149	156337	31.034 ng/ul	96
84) 3,3'-Dichlorobenzidine	21.765	252	112758	29.055 ng/ul	98
85) Benzo(a)anthracene	21.859	228	344622	30.483 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.724	149	224856	31.019 ng/ul	100
87) Chrysene	21.930	228	332454	30.611 ng/ul	99
89) Di-n-octyl phthalate	22.987	149	386305	30.978 ng/ul	100
90) Benzo(b)fluoranthene	24.192	252	352177	30.317 ng/ul	98
91) Benzo(k)fluoranthene	24.262	252	324752	29.791 ng/ul	99
93) Benzo(a)pyrene	25.120	252	332671	30.018 ng/ul	98
94) Indeno(1,2,3-cd)pyrene	29.186	276	374223	30.176 ng/ul	98
95) Dibenzo(a,h)anthracene	29.251	278	315791	30.015 ng/ul	98
96) Benzo(g,h,i)perylene	30.420	276	312635	29.963 ng/ul	96

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