#### Quantitation Report (QT Reviewed)

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

Data File : BG051242.D

Acq On : 25 Nov 2021 13:13

Operator : CG/JU Sample : P8140983BS

Misc

ALS Vial : 66 Sample Multiplier: 1

Quant Time: Nov 26 02:28:57 2021

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

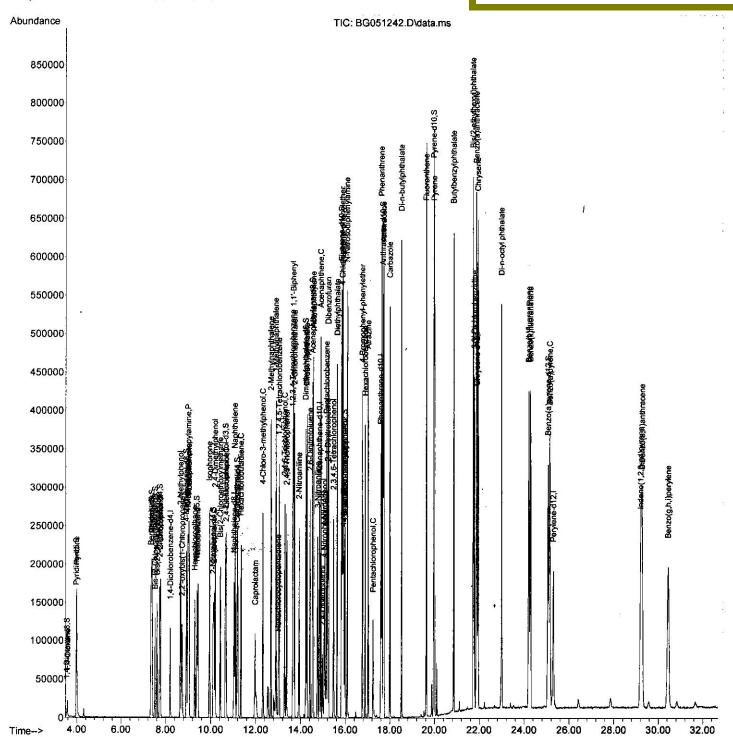
Quant Title : SVOA CALIBRATION

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



#### **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021



SFAM-EPA-BG112321.M Fri Nov 26 02:32:27 2021

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

Data File : BG051242.D

Acq On 25 Nov 2021 13:13

Operator : CG/JU Sample : PB140983BS

Misc

ALS Vial : 66 Sample Multiplier: 1

Quant Time: Nov 26 02:28:57 2021

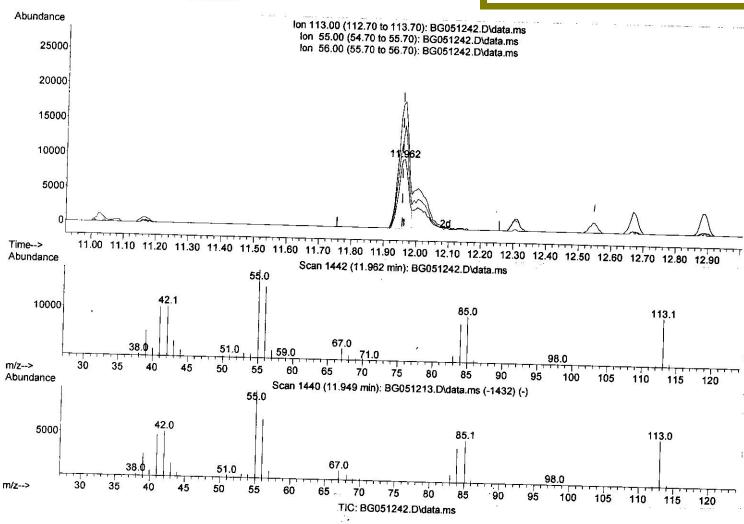
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

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 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021



### (34) Caprolactam

11.962min (+ 0.005) 22.34 ng/ul

response	21030	21030	
Ion	Exp&	Act%	
113.00	100.00	100.00	
55.00	183.80	182.44	
56.00	136.50	148.44	
0.00	0.00	0.00	

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

Data File : BG051242.D

Acq On : 25 Nov 2021 13:13

Operator : CG/JU Sample PB140983BS

Misc

ALS Vial : 66 Sample Multiplier: 1

Quant Time: Nov 26 02:28:57 2021

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

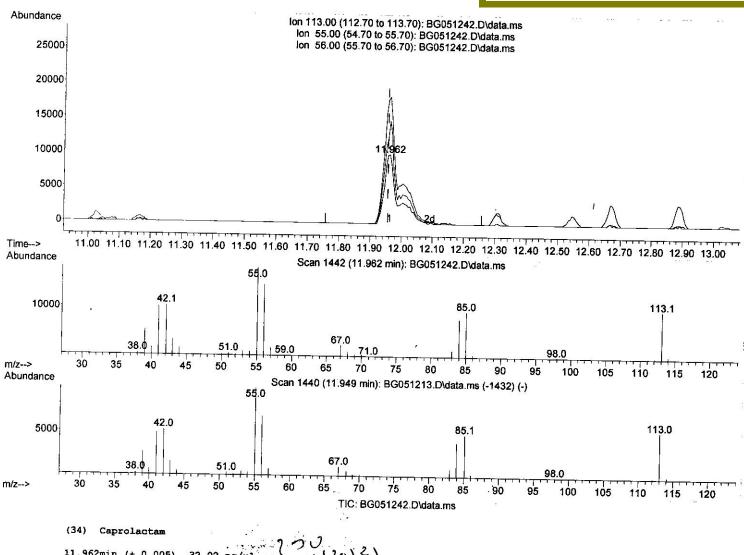
Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

Instrument: BNA\_G

ClientSampleId:

#### **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021



11/30121 11.962min (+ 0.005) 32.02 ng/ul

response	30138	
Ion	Expt	Act%
113.00	100.00	100.00
55.00	183.80	182.44
56.00	136.50	148.44
0.00	0.00	0.00

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

Data File : BG051242.D

Acq On : 25 Nov 2021 13:13

Operator : CG/JU Sample : PB140983BS

Misc

ALS Vial : 66 Sample Multiplier: 1

Quant Time: Nov 26 02:28:57 2021

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

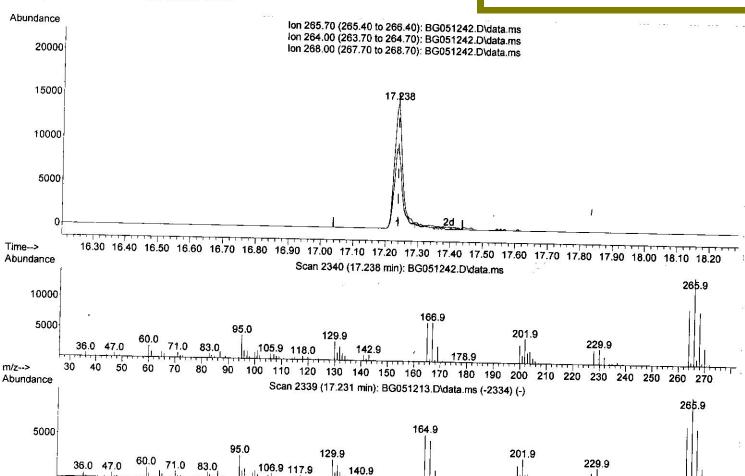
Quant Title : SVOA CALIBRATION

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

Instrument: BNA\_G ClientSampleId:

#### Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021



TIC: BG051242.D\data.ms

100 110 120 130 140 150 160 170 180 190 200 210 220 230 240 250 260 270

#### (71) Pentachlorophenol (C)

m/z-->

17.238min (-0.001) 23.90 ng/ul

80

90

response	25112	
Ion	Exp&	Act*
265.70	100.00	100.00
264.00	67.90	63.84
268,00	63.80	61.34
0.00	0.00	0.00

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

Data File : BG051242.D

Acq On : 25 Nov 2021 13:13

Operator : CG/JU Sample : P8140983BS

Misc

ALS Vial : 66 Sample Multiplier: 1

Quant Time: Nov 26 02:28:57 2021

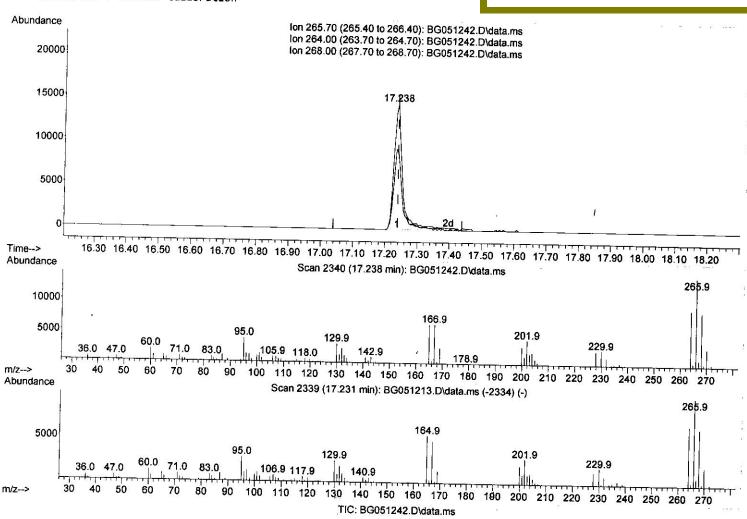
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

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 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021



### (71) Pentachlorophenol (C)

17.238min (-0.001) 24.30 ng/ul m 3 11 30 12

response	25535		
Ion	Expt	Act%	
265.70	100.00	100.00	
264.00	67.90	63.84	
268.00	63.80	61.34	
0.00	0.00	0.00	

#### Quantitation Report (QT Reviewed)

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

Data File : BG051242.D

Acq On : 25 Nov 2021 13:13

Operator : CG/JU Sample : PB140983BS

Misc

ALS Vial : 66 Sample Multiplier: 1

Quant Time: Nov 26 02:28:57 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:** SLCS983

#### **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021

1

Compound	R.T	. QIon	Response	Conc Units De	v(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.196	5 152	72164	22 222	
20) Naphthalene-d8	11.028			20.000 ng/ul	0.00
38) Acenaphthene-d10	14.836		150547	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.579		102530		0.00
79) Chrysene-d12	21.886		216967	20.000 ng/ul	0.00
88) Perylene-d12	25.288		185936	20.000 ng/ul	0.00
, , , , , , , , , , , , , , , , , , , ,	23.200	264	187707	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.537	96	5098	5.508 ng/uL	0.00
<ol><li>4) Pyridine-d5</li></ol>	3.966		75478	27.791 ng/ul	0.00
7) Phenol-d5	7.362		106692	33.563 ng/ul	-0.01
<ol><li>Bis-(2-Chloroethyl)eth</li></ol>	7.515		64790	32.452 ng/ul	0.00
11) 2-Chlorophenol-d4	7.732		76640	33.480 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.913		85166	33.200 ng/ul	0.00
21) Nitrobenzene-d5	9.377		41280	32.483 ng/ul	0.00
24) 2-Nitrophenol-d4	10.100		47651	33.240 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.652		81875	33.662 ng/ul	0.00
31) 4-Chloroaniline-d4	11.163	100000000000000000000000000000000000000	103560	29.099 ng/ul	0.00
46) Dimethylphthalate-d6	14,224		246637	31.263 ng/ul	0.00
49) Acenaphthylene-d8	14.530		319241	32.091 ng/ul	0.00
54) 4-Nitrophenol-d4	15.047		36926	28.917 ng/ul	0.00
60) Fluorene-d10	15.822	176	224311	31.575 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.952	200	40318	30.114 ng/ul	0.00
73) Anthracene-d10	17.685		330226	31.824 ng/ul	0.00
81) Pyrene-d10	19.959	212	374872		0.00
92) Benzo(a)pyrene-d12	25.053		329504	33.320 ng/ul 32.869 ng/ul	0.00 0.01
Target Compounds				2001	
2) 1,4-Dioxane	3.578	88	11777		lue
5) Pyridine	3.989	79	11773	11.278 ng/uL	94
6) Benzaldehyde	7.332	7 <b>7</b>	83662	29.603 ng/ul	97
8) Phenol	7.385		70358	34.754 ng/ul	97
10) Bis(2-Chloroethyl)ether	7.609	94 93	112681	34.217 ng/ul	98
12) 2-Chlorophenol	7.761		83455	33.497 ng/ul	96
13) 2-Methylphenol	8.649	128 108	79342	34.013 ng/ul	99
14) 2,2'-oxybis(1-Chloropr	8.725		84069	34.273 ng/ul	98
16) Acetophenone	9.030		122055 132367	33.950 ng/ul	98
17) N-Nitroso-di-n-propyla	9.007	705			97
18) 4-Methylphenol	8.978	108	75822	33.253 ng/ul	96
19) Hexachloroethane	9.283		90924	34.665 ng/ul	98
22) Nitrobenzene	9.418	117	32349	32.832 ng/ul	96
23) Isophorone	9.941	77 92	110634	33.201 ng/ul	99
25) 2-Nitrophenol	10.135	82	211169	32.618 ng/ul	99
26) 2,4-Dimethylphenol	10.133	139	49763	33.514 ng/ul	98
27) Bis(2-Chloroethoxy)met	10.417	107	99219	32.682 ng/ul	99
29) 2,4-Dichlorophenol	10.417	93	118213	33.076 ng/ul	98
30) Naphthalene		162	79581	33.238 ng/ul	96
32) 4-Chloroaniline	11.075 11.193	128	267193	32.618 ng/ul	98
33) Hexachlorobutadiene		127	106758	29.880 ng/ul	99 7 11 30 12 1
34) Caprolactam	11.339	225	51238	31.026 ng/ul	97 111 30 1
35) 4-Chloro-3-methylphenol	11.962 12.309	113	30138mS	32.019 ng/ul	
	±4.30 <del>3</del>	107	99793	34.697 ng/ul	97

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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 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021

Compound  R.T. QIon Response Conc Units Dev(Min)  36) 2-Methylnaphthalene 12.673 142 182182 32.697 ng/ul 99 37) 1-Methylnaphthalene 12.891 142 184600 32.203 ng/ul 98 39) 1,2,4,5-Tetrachloroben 13.032 216 106484 33.082 ng/ul 98 40) Hexachlorocyclopentadiene 12.996 237 19398 14.910 ng/ul 98 41) 2,4,6-Trichlorophenol 13.278 196 69355 34.335 ng/ul 99 42) 2,4,5-Trichlorophenol 13.361 196 72675 34.357 ng/ul 98
37) 1-Methylnaphthalene 12.891 142 184600 32.203 ng/ul 98 39) 1,2,4,5-Tetrachloroben 13.032 216 106484 33.082 ng/ul 98 40) Hexachlorocyclopentadiene 12.996 237 19398 14.910 ng/ul 98 41) 2,4,6-Trichlorophenol 13.278 196 69355 34.335 ng/ul 99 42) 2,4,5-Trichlorophenol 13.361 196 72675 34.357 ng/ul 98
37) 1-Methylnaphthalene 12.891 142 184600 32.203 ng/ul 98 39) 1,2,4,5-Tetrachloroben 13.032 216 106484 33.082 ng/ul 98 40) Hexachlorocyclopentadiene 12.996 237 19398 14.910 ng/ul 98 41) 2,4,6-Trichlorophenol 13.278 196 69355 34.335 ng/ul 99 42) 2,4,5-Trichlorophenol 13.361 196 72675 34.357 ng/ul 98
39) 1,2,4,5-Tetrachloroben 13.032 216 106484 33.082 ng/ul 98 40) Hexachlorocyclopentadiene 12.996 237 19398 14.910 ng/ul 98 41) 2,4,6-Trichlorophenol 13.278 196 69355 34.335 ng/ul 99 42) 2,4,5-Trichlorophenol 13.361 196 72675 34.357 ng/ul 98
40) Hexachlorocyclopentadiene 12.996 237 19398 14.910 ng/ul 98 41) 2,4,6-Trichlorophenol 13.278 196 69355 34.335 ng/ul 99 42) 2,4,5-Trichlorophenol 13.361 196 72675 34.357 ng/ul 98
41) 2,4,6-Trichlorophenol 13.278 196 69355 34.335 ng/ul 99 42) 2,4,5-Trichlorophenol 13.361 196 72675 34.357 ng/ul 98
42) 2,4,5-Trichlorophenol 13.361 196 72675 34.357 ng/ul 98
42\ 4 4  probable 7
43) 1,1'-Biphenyl 13.666 154 250317 32.687 ng/ul 97 44) 2-Chloronaphthalene 13.719 162 198704 32.619 ng/ul 100
4E\ 2 N\$+====214=
47\ Dimethilate 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1
40\ 0 C Di-11-17
EQ) Acompatibility
E1\ 2 Nd+ma==21
51) Assemble 192 .
E3\ 3 4 Di-11 32:000 lig/ul 33
EE) A Mitmode and 1
EC) Dibana C 91
56) Dibenzofuran 15.229 168 303075 32.417 ng/ul 98
57) 2,4-Dinitrotoluene 15.206 165 78678 32,841 ng/ul 94
58) 2,3,4,6-Tetrachlorophenol 15.458 232 53802 32.390 ng/ul 98
59) Diethylphthalate 15.629 149 266630 31.810 ng/ul 98
61) Fluorene 15.881 166 239591 31.993 ng/ul 100
62) 4-Chlorophenyl-phenyle 15.858 204 128221 31.771 ng/ul 99
63) 4-Nitroaniline 15.911 138 55636 34.482 ng/ul 96
66) 4,6-Dinitro-2-methylph 15.969 198 39990 30.971 ng/ul# 99
67) N-Nitrosodiphenylamine 16.075 169 214111 34.471 ng/ul 96
68) 4-Bromophenyl-phenylether 16.757 248 79846 34.337 ng/ul 94
69) Hexachlorobenzene 16.886 284 79682 33.605 ng/ul 98
70) Atrazine 17.021 200 83345 7 31.928 ng/ul 99
70) Acrazine 17.021 200 83345 1 31.928 ng/ul 99 71) Pentachlorophenol 17.238 266 25535m 24.304 ng/ul 72) Phenanthrene 17.636 170 201033
74) Anthonomic 17.020 1/8 394923 32.966 ng/ul 99
74) Anthracene 17.720 178 388490 32.653 ng/ul 99
75) 1,2,3,4-Tetrachloroben 13.637 216 110144 34.804 ng/uL 98
76) Pentachlorobenzene 15.153 250 97886 33.196 ng/uL 100
77) Carbazole 17.991 167 347002 33.227 ng/ul 98
78) Di-n-butylphthalate 18.513 149 450746 33.474 ng/ul 99 80) Fluoranthene 19.630 202 469103 33.948 ng/ul 97
03\ Dunana   105
93) Butulbassalatat 3
83) Butylbenzylphthalate 20.852 149 198488 35.322 ng/ul 93
84) 3,3'-Dichlorobenzidine 21.768 252 148462 34.294 ng/ul 99
85) Benzo(a)anthracene 21.862 228 429659 34.070 ng/ul 99
86) Bis(2-ethylhexyl)phtha 21.721 149 280887 34.736 ng/ul 100
87) Chrysene 21.933 228 408857 33.747 ng/ul 99
89) Di-n-octyl phthalate 22.985 149 477461 35.111 ng/ul 100
90) Benzo(b)fluoranthene 24.195 252 426962 33.705 ng/ul 97
91) Benzo(k)fluoranthene 24.271 252 402980 33.900 ng/ul 98
93) Benzo(a)pyrene 25.129 252 408892 33.834 ng/ul 99
94) Indeno(1,2,3-cd)pyrene 29.207 276 462937 34.231 ng/ul 98
95) Dibenzo(a,h)anthracene 29.266 278 388525 33.864 ng/ul 98
96) Benzo(g,h,i)perylene 30.441 276 386659 33.982 ng/ul 97

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