Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 2021

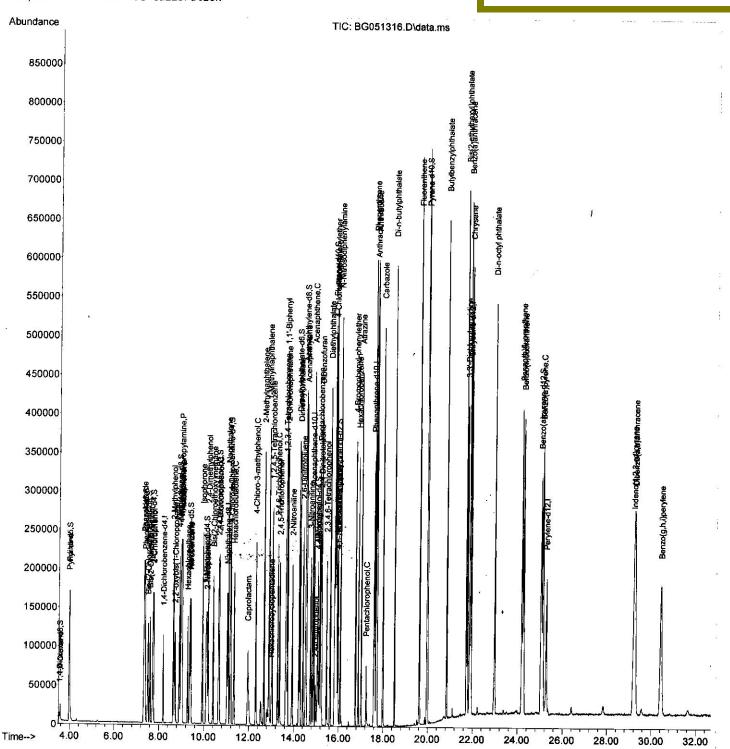
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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



Manual IntegrationsAPPROVED

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



SFAM-EPA-BG112321.M Fri Dec 03 00:59:47 2021

Page: 4

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File: BG051316.D

Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 2021

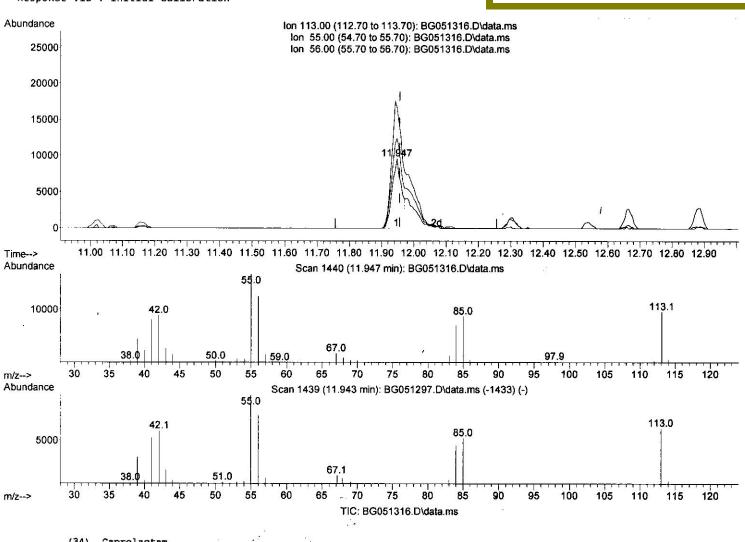
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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

Instrument :
BNA_G
ClientSampleId :

Manual IntegrationsAPPROVED

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(34) Caprolactam

11.947min (-0.010) 21.72 ng/ul

response	19329	9		
Ion	Ежр%	Act%		
113.00	100.00	100.00		
55.00	183.80	173.50		
56.00	136.50	129.78		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 2021

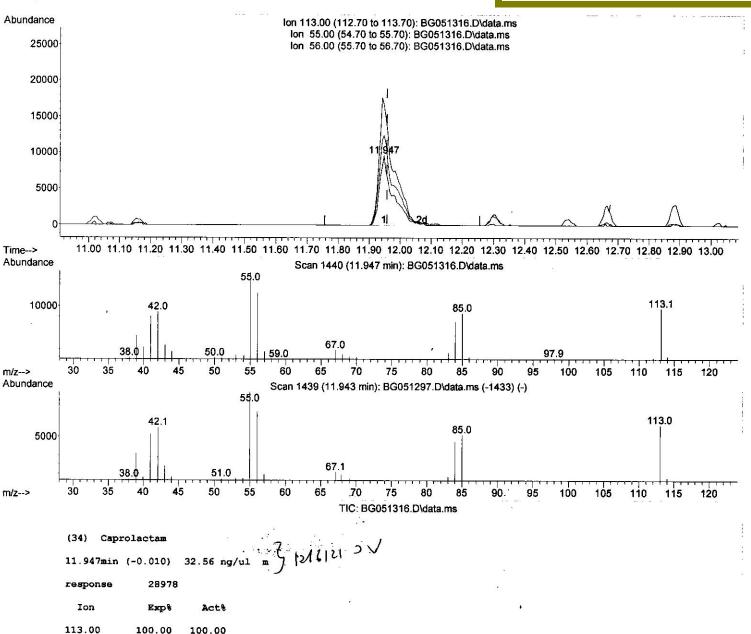
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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

Instrument :
BNA_G
ClientSampleId :

Manual IntegrationsAPPROVED

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183.80

136.50

0.00

173.50

129.78

0.00

55.00

56.00

0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 2021

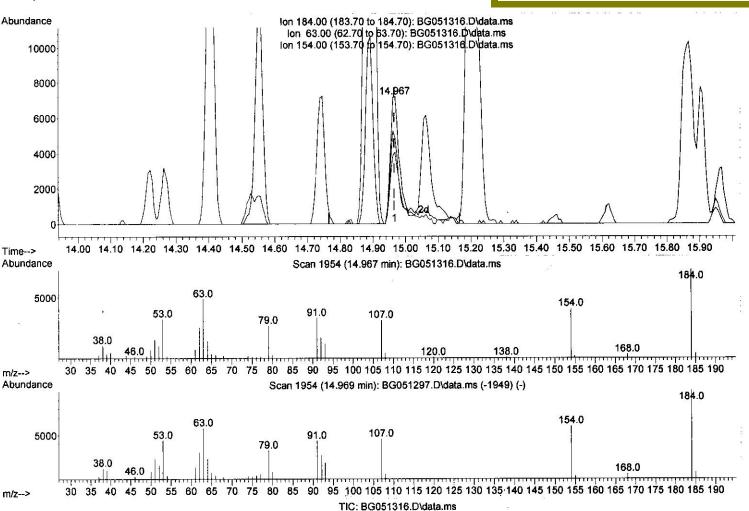
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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

Instrument :
BNA_G
ClientSampleId :

Manual IntegrationsAPPROVED

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(53) 2,4-Dinitrophenol

14.967min (+ 0.002) 16.61 ng/ul

response	14039			
Ion	Expt	Act*		
184.00	100.00	100.00		
63.00	82.70	67.65		
154.00	67.00	56.55		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 2021

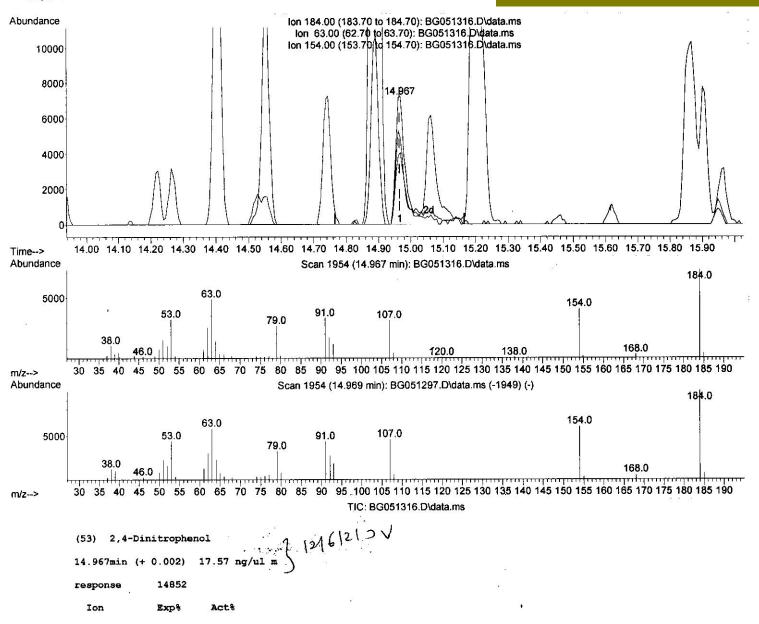
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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

82.70

67.00

0.00

100.00

67.65

56.55

0.00

184.00

63.00

154.00

0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 2021

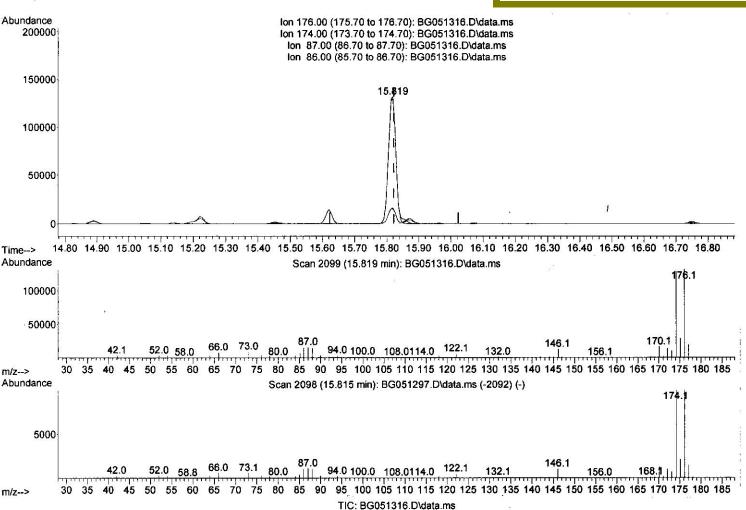
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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

Instrument :
BNA_G
ClientSampleId :

Manual IntegrationsAPPROVED

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(60) Fluorene-d10 (S)

15.819min (-0.004) 33.31 ng/ul

response	215790	
Ion	Exp%	Act%
176.00	100.00	100.00
174.00	97.50	98.11
87.00	10.60	12.11
86.00	10.90	11.36

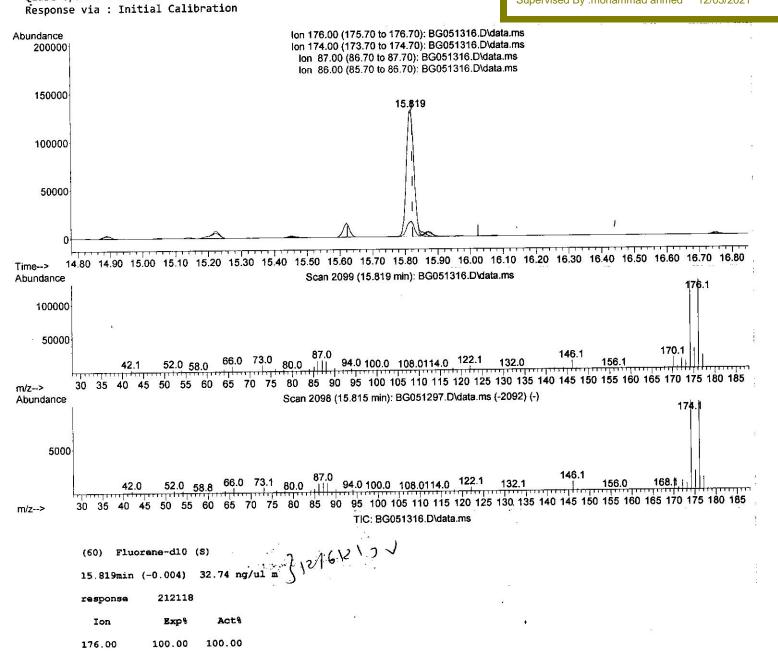
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\ Data File : BG051316.D : 2 Dec 2021 23:53 Acq On : CG/JU Operator : PB141136BS Sample Misc Sample Multiplier: 1 : 20 ALS Vial Quant Time: Dec 03 00:54:44 2021 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021

Instrument: BNA_G ClientSampleId:

Manual IntegrationsAPPROVED

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97.50

10.60

10.90

174.00 87.00

86.00

98.11

12.11 11.36

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

: 20 Sample Multiplier: 1 ALS Vial

Quant Time: Dec 03 00:54:44 2021

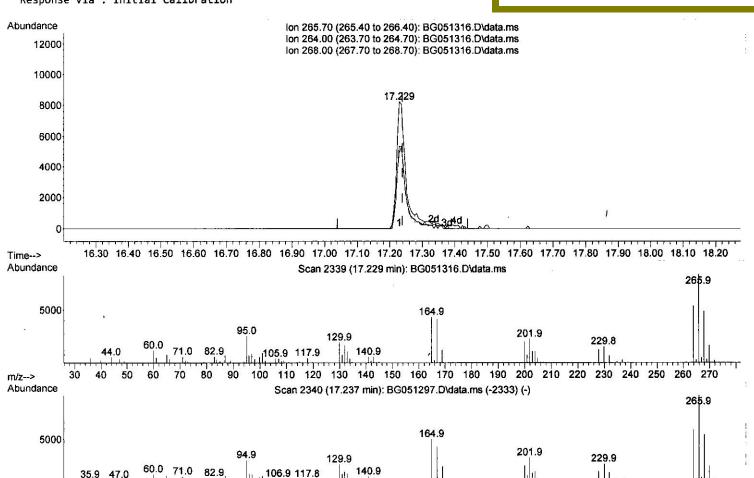
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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

Instrument: BNA_G ClientSampleId:

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TIC: BG051316.D\data.ms

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

106.9 117.8

(71) Pentachlorophenol (C)

35.9

m/z-->

47.0

17.229min (-0.010) 15.15 ng/ul

response	15088	
Ion	Ежр∜	Acts
265.70	100.00	100.00
264.00	67.90	65.07
268.00	63.80	59.46
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

: 2 Dec 2021 23:53 Acq On

: CG/JU Operator : PB141136BS Sample

Misc

Sample Multiplier: 1 ALS Vial : 20

Quant Time: Dec 03 00:54:44 2021

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

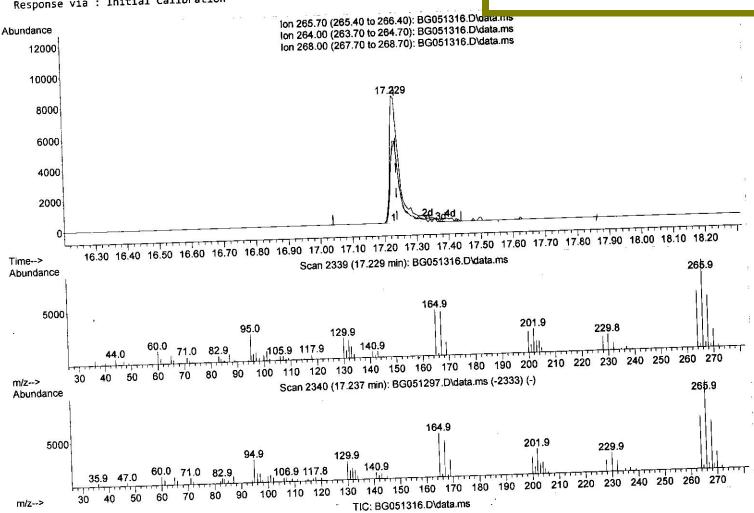
Quant Title : SVOA CALIBRATION

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

Instrument: BNA_G ClientSampleId:

Manual IntegrationsAPPROVED

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(71) Pentachlorophenol (C) 17.229min (-0.010) 16.38 ng/ul m

response	16320	
Ion	Exp8	Act*
265.70	100.00	100.00
264.00	67.90	65.07
268.00	63.80	59.46
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

: 2 Dec 2021 23:53 Acq On

: CG/JU Operator : PB141136BS Sample

Misc

Sample Multiplier: 1 ALS Vial : 20

Quant Time: Dec 03 00:54:44 2021

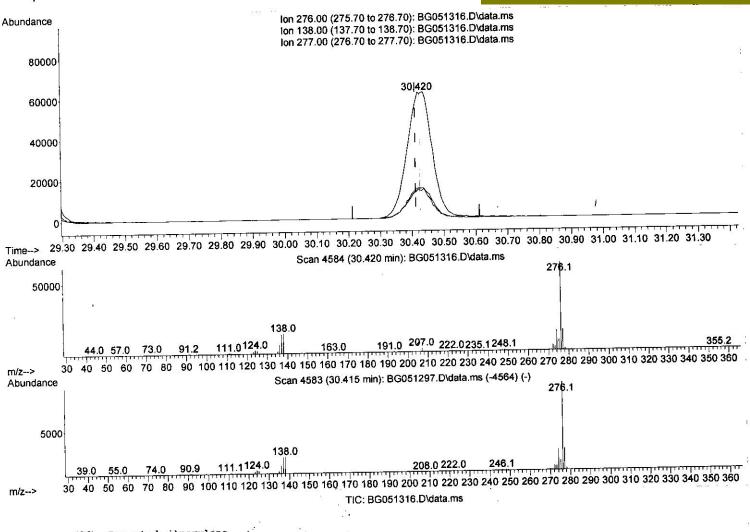
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

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



(96)	Benzo	(g,	h,ı)	pery.	Lene

30.420min (+ 0.008) 16.84 ng/ul .

response	179414	15
Ion	Ежр%	Act%
276.00	100.00	100.00
138.00	20.70	22.71
277.00	22.00	23.71
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 2021

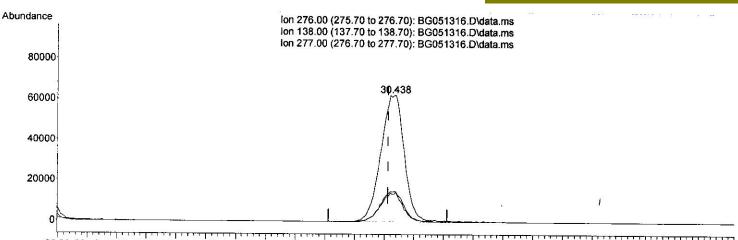
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

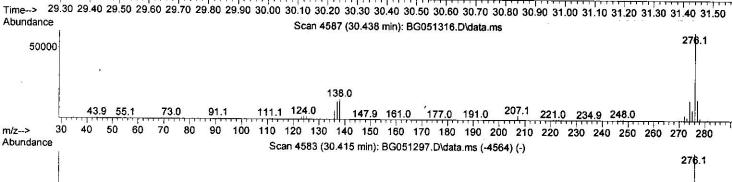
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
Response via : Initial Calibration

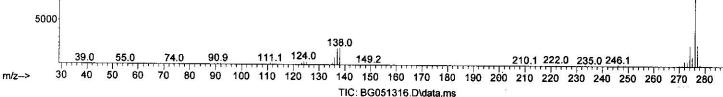
Instrument :
BNA_G
ClientSampleld :

Manual IntegrationsAPPROVED

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







(96) Benzo(g,h,i)perylene
30.438min (+ 0.025) 32.07 ng/ul m

response	341646	
Ion	Exp%	Act*
276.00	100.00	100.00
138.00	20.70	23.37
277.00	22.00	23,77
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 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: SLCS136

Manual IntegrationsAPPROVED

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

Compound	R.T.	QIon	Response	Conc Uni	its Dev	/(Min)		
Internal Standards								
1) 1,4-Dichlorobenzene-d4	8.193	152	31719	20.000	ng/ul	-0.01		
20) Naphthalene-d8	11.019	136	142348	20.000		0.00		
38) Acenaphthene-d10	14.826	164	93494	20.000		0.00		
64) Phenanthrene-d10	17.576	188	205688	20.000		0.00		
79) Chrysene-d12	21.877	240	174630	20.000		0.00		
88) Perylene-d12	25.279	264	175721	20.000		0.00		
System Monitoring Compounds								
3) 1,4-Dioxane-d8	3.534	96	5510	6.037	ng/uL	-0.01		
4) Pyridine-d5	3.963	84	80110	29.910	ng/ul	-0.02		
7) Phenol-d5	7.353	99	105771	33.740	ng/ul	0.00		
9) Bis-(2-Chloroethyl)eth	7.506	67	66643	33.848	ng/ul	-0.01	¥	1
11) 2-Chlorophenol-d4	7.723	132	75479	33.436	ng/ul	-0.01		
<pre>15) 4-Methylphenol-d8</pre>	8.910	113	84040	33.220	ng/ul	0.00		
21) Nitrobenzene-d5	9.368	128	39945	33.242	ng/ul	-0.01		
24) 2-Nitrophenol-d4	10.097	143	45692	33.709	ng/ul	0.00		
28) 2,4-Dichlorophenol-d3	10.643	165	76327	33.188	ng/ul	-0.01		
31) 4-Chloroaniline-d4	11.160	131	124946	37.130	ng/ul	0.00		
46) Dimethylphthalate-d6	14.221	166	240125	33.379	ng/ul	0.00		o ∨ -
49) Acenaphthylene-d8	14.527	160	305652	33.694	ng/ul	0.00	124121	
54) 4-Nitrophenol-d4	15.050	143	361037	31.005	ng/ul	0.00	10	
60) Fluorene-d10	15.819	176	212118m	32.744	ng/ul	0.00		
65) 4,6-Dinitro-2-methylph	15.949	200	37181	29.294		0.00		
73) Anthracene-d10	17.676	188	333515	33.903	ng/ul	0.00		
81) Pyrene-d10	19.956	212	379921	35.955	ng/ul	0.00		
92) Benzo(a)pyrene-d12	25.044	264	320381	34.139	ng/ul	0.00		
Target Compounds					Qv	/alue		
2) 1,4-Dioxane	3.569	88	11156	10.837	ng/uL	98		
5) Pyridine	3.980	79	83402	29.925	ng/ul	97		
Benzaldehyde	7.329	77	78807	39.474	ng/ul	94		
8) Phenol	7.382	94	105007	32.334	ng/ul	99		
10) Bis(2-Chloroethyl)ether	7.600	93	77535	31.558	ng/ul	100	e e	
<pre>12) 2-Chlorophenol</pre>	7.758	128	72750	31.625	ng/ul	98		
<pre>13) 2-Methylphenol</pre>	8.640	108	76416	31.590	ng/ul	98		
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.716	45	117217	33.061	ng/ul	99		
16) Acetophenone	9.021	105	122303	31.256	ng/ul	97		
17) N-Nitroso-di-n-propyla	8.998	70	72426	32.210	ng/ul	98		
<pre>18) 4-Methylphenol</pre>	8.974	108	82995	32.086	ng/ul	100		
19) Hexachloroethane	9.274	117	29817	30.687		98 -		
22) Nitrobenzene	9.415	77	102274.	32.460	ng/ul	97	72	
23) Isophorone	9.932	82	195635	31.959	ng/ul	99	<i>8</i> *	
<pre>25) 2-Nitrophenol</pre>	10.132	139	43957	31.308	ng/ul	98		
26) 2,4-Dimethylphenol	10.179	107	93336	32.515		100		
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.408	93	110999	32.846		100		2529
<pre>29) 2,4-Dichlorophenol</pre>	10.672	162	69993	30.917		96	12/6/21	7.1
30) Naphthalene	11.072	128	244124	31.518	ng/ul	98	,2 6 1°1	
32) 4-Chloroaniline	11.184	127	99692	29.510		98	(-)	
33) Hexachlorobutadiene	11.336	225	43497 7			96		
34) Caprolactam	11.947	113	28978m -					
35) 4-Chloro-3-methylphenol	12.300	107	88911	32.694	ng/ul	94		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051316.D

Acq On : 2 Dec 2021 23:53

Operator : CG/JU Sample : PB141136BS

Misc

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Dec 03 00:54:44 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:
SLCS136

Manual IntegrationsAPPROVED

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

Compound	R.T. QIon	Response Conc Units Dev(Min)
36) 2-Methylnaphthalene	12.664 142	164589 31.241 ng/ul 98
37) 1-Methylnaphthalene	12.882 142	169123 31.203 ng/ul 97
39) 1,2,4,5-Tetrachloroben	13.028 216	88656 30.205 ng/ul 98
40) Hexachlorocyclopentadiene	12.993 237	12364 10.422 ng/ul 98
41) 2,4,6-Trichlorophenol	13.269 196	57521 31.229 ng/ul 99
42) 2,4,5-Trichlorophenol	13.358 196	59595 30.897 ng/ul 96
43) 1,1'-Biphenyl	13.657 154	221245 31.683 ng/ul 99
44) 2-Chloronaphthalene	13.710 162	173381 31.213 ng/ul 98
45) 2-Nitroaniline	13.916 65	67649 35.188 ng/ul 94
47) Dimethylphthalate	14.268 163	228688 31.406 ng/ul 100
48) 2,6-Dinitrotoluene	14.403 165	49827 32.577 ng/ul 91 287570 32.087 ng/ul 98 55167 36.489 ng/ul 97
50) Acenaphthylene	14.556 152	287570 32.087 ng/ul 98 //) ン
51) 3-Nitroaniline	14.738 138	55167 36.489 ng/ul 97 1
52) Acenaphthene	14.891 153	186892 / 31.620 ng/ul 93
53) 2,4-Dinitrophenol	14.967 184	14852m \ 17.567 ng/ul
55) 4-Nitrophenol	15.061 109	28680 > 28.392 ng/ul 92
56) Dibenzofuran	15.226 168	264750 31.054 ng/ul 99
57) 2,4-Dinitrotoluene	15.197 165	71778 32.856 ng/ul 97
58) 2,3,4,6-Tetrachlorophenol	15.455 232	44316 29.258 ng/ul 97
59) Diethylphthalate	15.620 149	244486 31.987 ng/ul 99
61) Fluorene	15.872 166	213605 31.280 ng/ul 99
62) 4-Chlorophenyl-phenyle	15.855 204	108513 29.486 ng/ul 99
63) 4-Nitroaniline	15.907 138	58395 39.690 ng/ul 95
66) 4,6-Dinitro-2-methylph	15.966 198	33759 27.579 ng/ul# 97
67) N-Nitrosodiphenylamine	16.072 169	33759 27.579 ng/ul# 9/ 192369 32.669 ng/ul 99 67963 30.829 ng/ul 97 68988 30.691 ng/ul 97 78022 31.527 ng/ul 99
68) 4-Bromophenyl-phenylether	16.748 248	67963 30.829 ng/ul 97
69) Hexachlorobenzene	16.877 284	68988 30.691 ng/ul 97 12 1 78022 31.527 ng/ul 99
70) Atrazine	17.012 200	
71) Pentachlorophenol	17.229 266	
72) Phenanthrene	17.617 178	365870) 32.216 ng/ul 99 367342 32.568 ng/ul 98
74) Anthracene	17.711 178	
75) 1,2,3,4-Tetrachloroben	13.634 216	
76) Pentachlorobenzene	15.144 250	
77) Carbazole	17.987 167	
78) Di-n-butylphthalate	18.504 149	
80) Fluoranthene	19.621 202 19.985 202	435452 34.301 ng/ul 99
82) Pyrene	20.843 149	
83) Butylbenzylphthalate	21.759 252	
<pre>84) 3,3'-Dichlorobenzidine 85) Benzo(a)anthracene</pre>	21.859 228	
86) Bis(2-ethylhexyl)phtha	21.712 149	
87) Chrysene	21.930 228	
89) Di-n-octyl phthalate	22.976 149	
90) Benzo(b)fluoranthene	24.192 252	
91) Benzo(k)fluoranthene	24.262 252	358809 32.243 ng/ul 99 , 12\
93) Benzo(a)pyrene	25.120 252	371434 32.831 ng/ul 99
94) Indeno(1,2,3-cd)pyrene	29.204 276	411843 32.531 ng/ul 97 ,2
95) Dibenzo(a,h)anthracene	29.251 278	345192 32.139 ng/ul 98
96) Benzo(g,h,i)perylene	30.438 276	

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