Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111121\

Data File : BG051004.D

Acq On : 12 Nov 2021 18:41

Operator : CG/JU Sample : PB140711BS

Misc

ALS Vial : 43 Sample Multiplier: 1

Quant Time: Nov 15 00:37:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110321.M

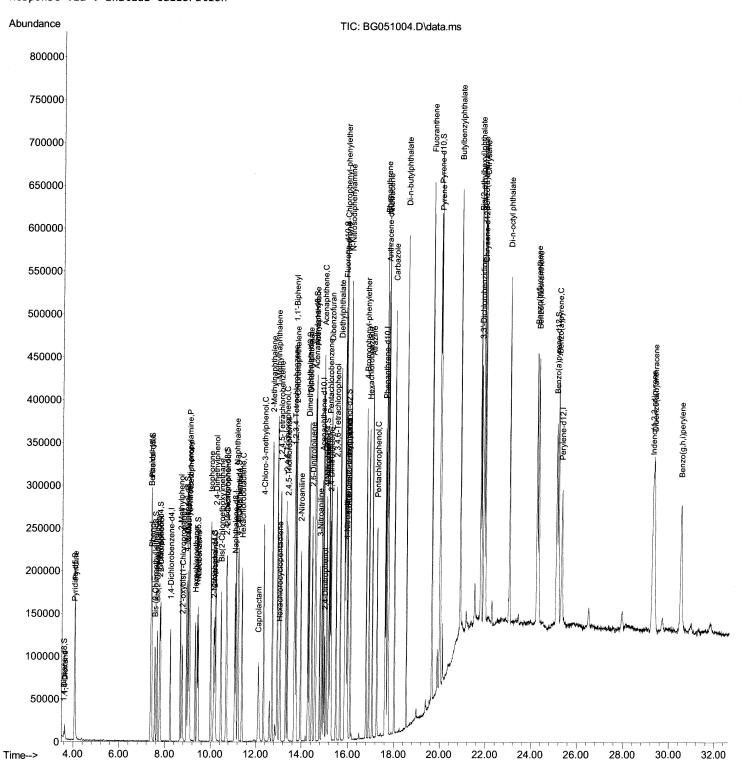
Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 15 00:27:19 2021 Response via : Initial Calibration



# Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/15/2021 Supervised By :mohammad ahmed 11/17/2021



### Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111121\

Data File: BG051004.D

Acq On : 12 Nov 2021 18:41

Operator : CG/JU Sample : PB140711BS

Misc

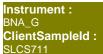
ALS Vial : 43 Sample Multiplier: 1

Quant Time: Nov 15 00:37:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110321.M

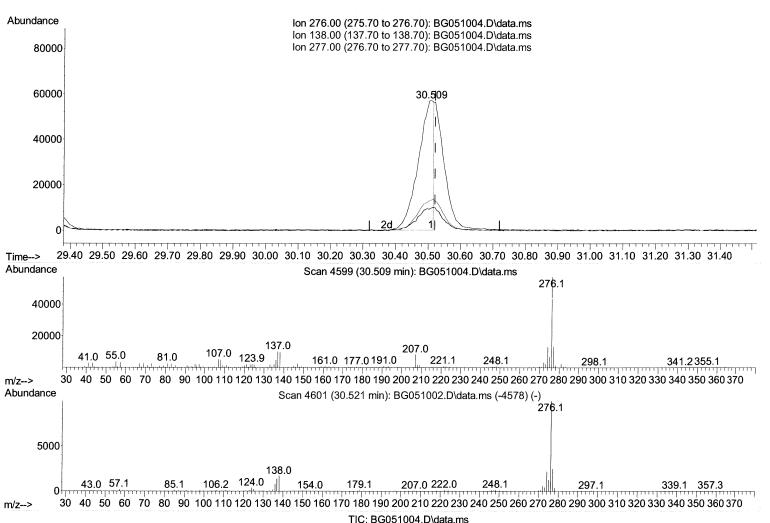
Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 15 00:27:19 2021 Response via : Initial Calibration



## **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 11/15/2021 Supervised By :mohammad ahmed 11/17/2021



### (96) Benzo(g,h,i)perylene

30.509min (-0.012) 17.67 ng/ul

response	186759	
Ion	Ехр%	Act%
276.00	100.00	100.00
138.00	20.70	17.27
277.00	22.00	23.57
0.00	0.00	0.00

#### Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111121\

Data File : BG051004.D

Acq On : 12 Nov 2021 18:41

Operator : CG/JU : PB140711BS Sample

Misc

ALS Vial : 43 Sample Multiplier: 1

Quant Time: Nov 15 00:37:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA G\Methods\SFAM-EPA-BG110321.M

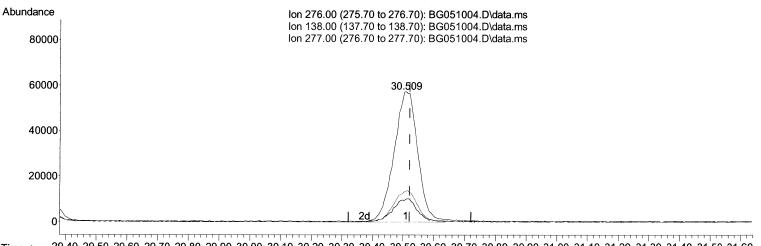
Quant Title : SVOA CALIBRATION

QLast Update: Mon Nov 15 00:27:19 2021 Response via : Initial Calibration

Instrument : BNA\_G ClientSampleId: SLCS711

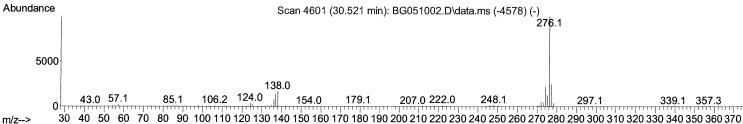
# **Manual Integrations APPROVED**

Reviewed By :Jagrut Upadhyay 11/15/2021 Supervised By:mohammad ahmed 11/17/2021



29.40 29.50 29.60 29.70 29.80 29.90 30.00 30.10 30.20 30.30 30.40 30.50 30.60 30.70 30.80 30.90 31.00 31.10 31.20 31.30 31.40 31.50 31.60 Time--> Abundance Scan 4599 (30.509 min): BG051004.D\data.ms 276.1 40000 20000 137.0 107.0 123.9 41.0 55.0 161.0 177.0 191.0 248.1

298.1 30 40 50 60 70 80 90 100 110 120 130 140 150 160 170 180 190 200 210 220 230 240 250 260 270 280 290 300 310 320 330 340 350 360 370 m/z--> Abundance Scan 4601 (30.521 min): BG051002.D\data.ms (-4578) (-)



TIC: BG051004.D\data.ms

## (96) Benzo(g,h,i)perylene

30.509min (-0.012) 29.29 ng/ul m \\\\7/2\J\

response	309633			
Ion	Ехр%	Act%		
276.00	100.00	100.00		
138.00	20.70	17.27		
277.00	22.00	23.57		
0.00	0.00	0.00		

339.1 357.3

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111121\

Data File : BG051004.D

Acq On : 12 Nov 2021 18:41

Operator : CG/JU Sample : PB140711BS

Misc

ALS Vial : 43 Sample Multiplier: 1

Quant Time: Nov 15 00:37:30 2021

 $\label{thm:lem1_BNA_GMethods\SFAM-EPA-BG110321.M} Quant \ \mbox{Methods\SFAM-EPA-BG110321.M}$ 

Quant Title : SVOA CALIBRATION QLast Update : Mon Nov 15 00:27:19 2021 Response via : Initial Calibration Instrument:
BNA\_G
ClientSampleId:
SLCS711

# **Manual IntegrationsAPPROVED**

Reviewed By: Jagrut Upadhyay 11/15/2021 Supervised By: mohammad ahmed 11/17/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.241	152	35090	20.000 ng/ul	0.00
20) Naphthalene-d8	11.067	136	159822	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.869	164	105013	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.612	188	223446	20.000 ng/ul	0.00
79) Chrysene-d12	21.913		168025	20.000 ng/ul	# 0.00
88) Perylene-d12	25.333	264	167202	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.599	96	4213	3.875 ng/uL	0.00
<ol><li>4) Pyridine-d5</li></ol>	4.052	84	76651	23.567 ng/ul	-0.01
7) Phenol-d5	7.389	99	93941	25.095 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.554	67	57095	23.611 ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.771	132	68005	26.213 ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.946	113	73803	25.043 ng/ul	0.00
21) Nitrobenzene-d5	9.416	128	35426	26.083 ng/ul	0.00
24) 2-Nitrophenol-d4	10.145	143	42211	27.951 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.685	165	71305	28.030 ng/ul	0.00
31) 4-Chloroaniline-d4	11.202	131	98947	25.684 ng/ul	0.00
46) Dimethylphthalate-d6	14.263	166	233674	29.085 ng/ul	0.00
49) Acenaphthylene-d8	14.563	160	292291	29.201 ng/ul	0.00
54) 4-Nitrophenol-d4	15.062	143	43221	29.670 ng/ul	0.00
60) Fluorene-d10	15.856	176	204621	28.750 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.973	200	39820	29.394 ng/ul	0.00
73) Anthracene-d10	17.712	188	305929	28.959 ng/ul	0.00
<ul><li>81) Pyrene-d10</li><li>92) Benzo(a)pyrene-d12</li></ul>	19.986 25.098	212 264	314195 257578	28.951 ng/ul 27.867 ng/ul	0.00 0.00
Target Compounds				Ov	alue
2) 1,4-Dioxane	3.635	88	9858	8.255 ng/uL	96
5) Pyridine	4.075	79	79302	23.554 ng/ul	94
6) Benzaldehyde	7.377	77	63057	26.706 ng/ul	99
8) Phenol	7.413	94	100860	26.046 ng/ul	99
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.648	93	71230	24.574 ng/ul	98
12) 2-Chlorophenol	7.800	128	69494	26.382 ng/ul	98
13) 2-Methylphenol	8.676	108	73584	25.708 ng/ul	98
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.764	45	108926	23.858 ng/ul	99
16) Acetophenone	9.075	105	120948	26.418 ng/ul	99
17) N-Nitroso-di-n-propyla	9.058	70	70961	25.690 ng/ul	99
<pre>18) 4-Methylphenol</pre>	9.011	108	79590	26.115 ng/ul	96
19) Hexachloroethane	9.328	117	27877	25.312 ng/ul	97
22) Nitrobenzene	9.463	77	100340	26.489 ng/ul	99
23) Isophorone	10.004	82	199092	27.082 ng/ul	97
25) 2-Nitrophenol	10.174	139	43119	28.460 ng/ul	98
26) 2,4-Dimethylphenol	10.221	107	90022	26.998 ng/ul	99
27) Bis(2-Chloroethoxy)met	10.456	93	103717	26.184 ng/ul	97
29) 2,4-Dichlorophenol	10.709	162	71325	28.744 ng/ul	97
30) Naphthalene	11.120	128	240638	27.535 ng/ul	97
32) 4-Chloroaniline	11.226	127	96603	25.254 ng/ul	99
33) Hexachlorobutadiene	11.384	225	46025	28.260 ng/ul	97
34) Caprolactam	12.101	113	28756	27.297 ng/ul	96 100
35) 4-Chloro-3-methylphenol	12.330	107	91508	28.868 ng/ul	100

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111121\

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Misc :

ALS Vial : 43 Sample Multiplier: 1

Quant Time: Nov 15 00:37:30 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 15 00:27:19 2021 Response via : Initial Calibration

Instrument:
BNA\_G
ClientSampleId:
SLCS711

## Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/15/2021 Supervised By: mohammad ahmed 11/17/2021

Compound	R.T.	QIon	Response	Conc Un:	its Dev	(Min)
36) 2-Methylnaphthalene	12.706	142	164796	27.679	ng/u1	99
37) 1-Methylnaphthalene	12.924	142	167586	27.779	-	97
39) 1,2,4,5-Tetrachloroben	13.071		92308	30.168		95
40) Hexachlorocyclopentadiene	13.035	237	28279	19.246	-	98
41) 2,4,6-Trichlorophenol	13.306		66223	33.075	-	99
42) 2,4,5-Trichlorophenol	13.382		71200	33.123		100
43) 1,1'-Biphenyl	13.699	154	227240	29.605	_	98
44) 2-Chloronaphthalene	13.752	162	178527	29.680	-	98
45) 2-Nitroaniline	13.952	65	68828	28.801	ng/ul	94
47) Dimethylphthalate	14.310	163	237161	29.522	ng/ul	99
48) 2,6-Dinitrotoluene	14.440	165	49900	29.678		96
50) Acenaphthylene	14.592	152	296541	29.560	ng/ul	99
51) 3-Nitroaniline	14.775	138	49747	28.607	ng/ul	97
52) Acenaphthene	14.933	153	191784	29.074	ng/ul	95
53) 2,4-Dinitrophenol	14.986	184	27090	29.206		90
55) 4-Nitrophenol	15.074	109	39645	29.675	_	97
56) Dibenzofuran	15.262	168	274127	29.033		99
57) 2,4-Dinitrotoluene	15.227	165	70261	29.281	•	97
58) 2,3,4,6-Tetrachlorophenol	15.486	232	58145	34.420	-	98
59) Diethylphthalate	15.662	149	248190	28.863		99
61) Fluorene	15.914	166	218446	29.230	-	99
62) 4-Chlorophenyl-phenyle	15.897	204	117198	30.116	•	93
63) 4-Nitroaniline	15.938	138	50133	29.060	_	95
66) 4,6-Dinitro-2-methylph	15.991	198	39762	30.097		98
67) N-Nitrosodiphenylamine	16.108	169	195684	31.330		98
68) 4-Bromophenyl-phenylether	16.790	248	72719	32.717	_	93
69) Hexachlorobenzene	16.913	284	73997	32.384	_	97
70) Atrazine	17.054	200	74595	28.164	-	98
<ul><li>71) Pentachlorophenol</li><li>72) Phenanthrene</li></ul>	17.260	266 170	44126	42.062 30.212		97 99
74) Anthracene	17.654 17.748	178 178	360349 351641	29.383		99
75) 1,2,3,4-Tetrachloroben	13.670	216	97117	31.921	_	98
76) Pentachlorobenzene	15.180	250	89268	31.666	_	99
77) Carbazole	18.018	167	321449	29.968	_	99
78) Di-n-butylphthalate	18.547	149	401107	28.458	-	99
80) Fluoranthene	19.657	202	392916	30.168		96
82) Pyrene	20.015	202	379256	29.802	•	97
83) Butylbenzylphthalate	20.879	149	152777	27.918	_	97
84) 3,3'-Dichlorobenzidine	21.802	252	115104	28.129	_	99
85) Benzo(a)anthracene	21.896	228	340508	29.273		99
86) Bis(2-ethylhexyl)phtha	21.760	149	216044	27.503	_	100
87) Chrysene	21.966	228	329084	29.615	_	99
89) Di-n-octyl phthalate	23.030	149	367981	27.033	-	100
90) Benzo(b)fluoranthene	24.240	252	340292	28.569		99
91) Benzo(k)fluoranthene	24.310	252	317785	28.431	_	99
93) Benzo(a)pyrene	25.180	252	325489	28.691	_	98
94) Indeno(1,2,3-cd)pyrene	29.275	276	369433	29.253	ng/ul	96
95) Dibenzo(a,h)anthracene	29.340	278	322232	30.156		97
96) Benzo(g,h,i)perylene	30.509	276	309633m >	29.290		11117/21 Ju

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