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

Data File : BG051005.D

Acq On : 12 Nov 2021 19:22

Operator : CG/JU Sample : PB140665BS

Misc :

ALS Vial : 44 Sample Multiplier: 1

Quant Time: Nov 15 00:42:15 2021

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

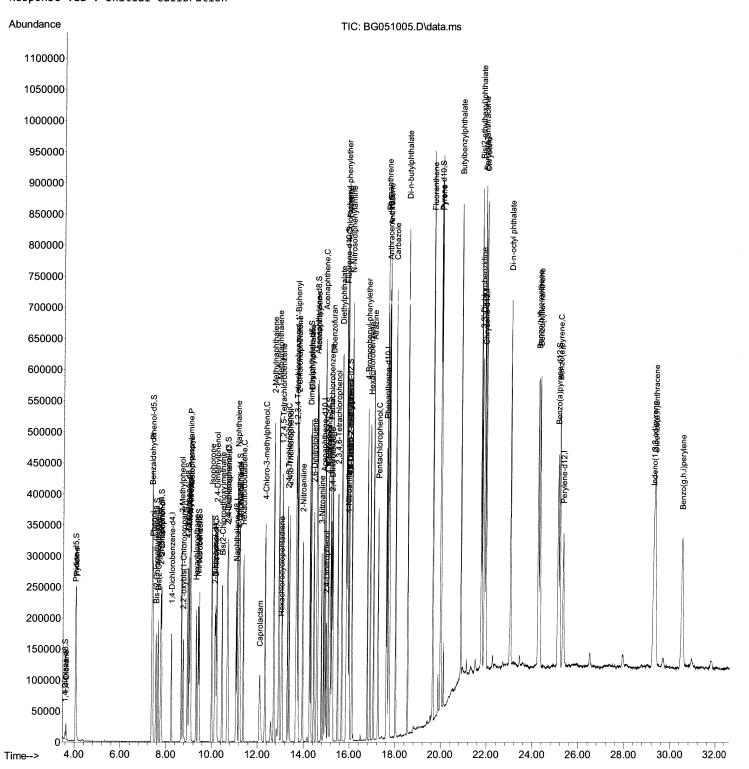
Quant Title : SVOA CALIBRATION

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: BG051005.D

Acq On : 12 Nov 2021 19:22

Operator : CG/JU Sample : PB140665BS

Misc :

ALS Vial : 44 Sample Multiplier: 1

Quant Time: Nov 15 00:42:15 2021

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

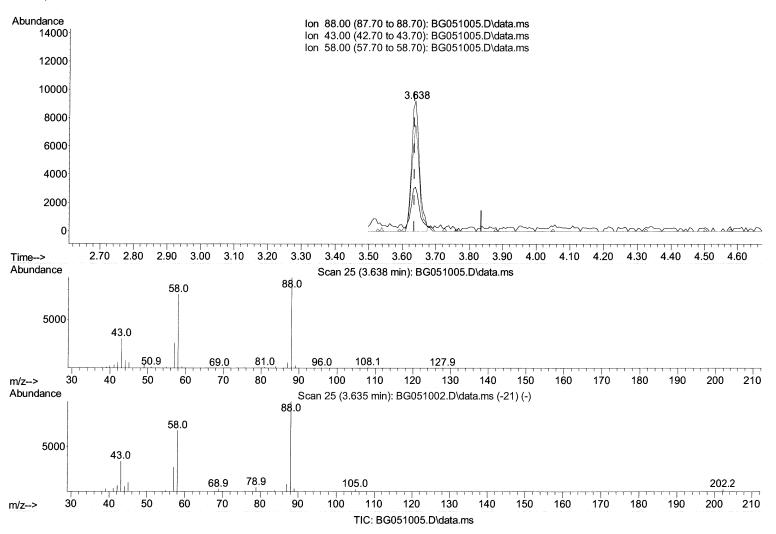
Quant Title : SVOA CALIBRATION

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



# (2) 1,4-Dioxane

3.638min (+ 0.003) 9.43 ng/uL

response	15028	
Ion	Ехр%	Act%
88.00	100.00	100.00
43.00	28.70	33.85
58.00	78.00	81.88
0.00	0.00	0.00

#### Quantitation Report (Qedit)

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

Data File : BG051005.D

Acq On : 12 Nov 2021 19:22

Operator : CG/JU Sample : PB140665BS

Misc :

ALS Vial : 44 Sample Multiplier: 1

Quant Time: Nov 15 00:42:15 2021

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

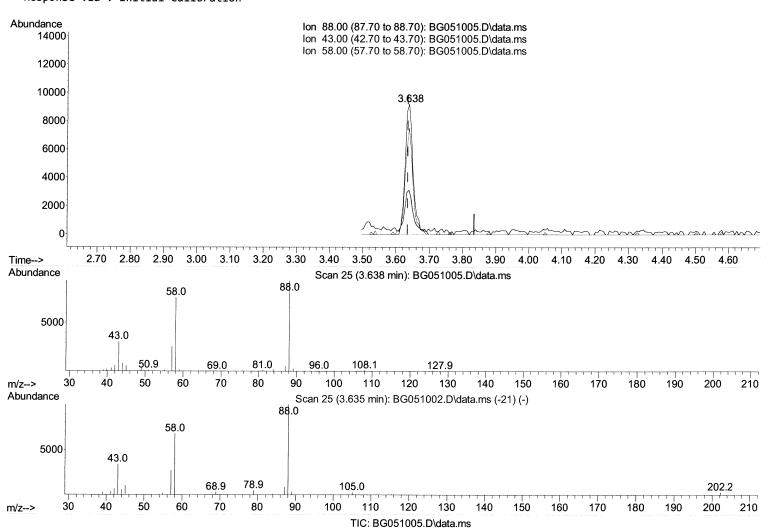
Quant Title : SVOA CALIBRATION

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



## (2) 1,4-Dioxane

3.638min (+ 0.003) 9.61 ng/uL m \\\\7/d\J\

response	15309	
Ion	Ехр%	Act%
88.00	100.00	100.00
43.00	28.70	33.85
58.00	78.00	81.88
0.00	0.00	0.00

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

Data File : BG051005.D

Acq On : 12 Nov 2021 19:22

Operator : CG/JU Sample : PB140665BS

Misc

ALS Vial : 44 Sample Multiplier: 1

Quant Time: Nov 15 00:42:15 2021

 $\label{thm:lem1_BNA_GMethods} \mbox{Quant Method}: Z:\svoasrv\end{\mbox{HPCHEM1\_BNA\_G\end{\mbox{Methods}}} \mbox{EPA-BG110321.M}$ 

Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 15 00:27:19 2021 Response via : Initial Calibration Instrument : BNA\_G ClientSampleId : SLCS665

# **Manual IntegrationsAPPROVED**

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

Compound	R.T.	QIon	Response	Conc Ur	its Dev(	Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.239	152	46830	20.000	ng/ul	0.00
20) Naphthalene-d8	11.070		212238		ng/ul	0.00
38) Acenaphthene-d10	14.866		137320		ng/ul	0.00
64) Phenanthrene-d10	17.610		294822		ng/ul	0.00
79) Chrysene-d12	21.917		230093		ng/ul	0.00
88) Perylene-d12	25.330		225524		ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.597	96	7000	4.824	ng/uL	0.00
<ol><li>4) Pyridine-d5</li></ol>	4.049	84	114155	26.299	ng/ul	-0.01
7) Phenol-d5	7.387	99	145848	29.193	ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.557	67	87810	27.209	ng/ul	0.00
<pre>11) 2-Chlorophenol-d4</pre>	7.768	132	104874	30.290	ng/ul	0.00
<pre>15) 4-Methylphenol-d8</pre>	8.944	113	113081	28.751	ng/ul	0.00
21) Nitrobenzene-d5	9.414	128	55094	30.546	_	0.00
24) 2-Nitrophenol-d4	10.142	143	63952	31.889	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.683	165	107041	31.686	_	0.00
31) 4-Chloroaniline-d4	11.200	131	145911	28.521		0.00
46) Dimethylphthalate-d6	14.267	166	328519	31.270		0.00
49) Acenaphthylene-d8	14.566	160	413975	31.628		0.00
54) 4-Nitrophenol-d4	15.066	143	63791	33.488	•	0.00
60) Fluorene-d10	15.859	176	291124	31.281		0.00
65) 4,6-Dinitro-2-methylph	15.976	200	58351	32.645	_	0.00
73) Anthracene-d10	17.716	188	435372	31.234		0.00
81) Pyrene-d10	19.989	212	469147	31.568	•	0.00
92) Benzo(a)pyrene-d12	25.101	264	384372	30.830	-	0.00
Target Compounds					Qva:	lue
2) 1,4-Dioxane	3.638	88	15309m>	9.606	ng/uL >	1117/21 Ju
5) Pyridine	4.073	79	119764	26.655		96
<ol><li>Benzaldehyde</li></ol>	7.375	77	97260	30.866		95
8) Phenol	7.416	94	151434	29.302		98
<pre>10) Bis(2-Chloroethyl)ether</pre>	7.651	93	110590	28.588		94
12) 2-Chlorophenol	7.798	128	105893	30.123		96
13) 2-Methylphenol	8.679	108	113081	29.603	-	96
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.756	45	163179	26.781	ng/ul	98
16) Acetophenone	9.079	105	179077	29.309		99
17) N-Nitroso-di-n-propyla	9.055	70	102876	27.907		99
18) 4-Methylphenol	9.014	108	119515	29.385		94
19) Hexachloroethane	9.325	117	43088	29.316	-	92
22) Nitrobenzene	9.461	77	147832	29.389	-	98
23) Isophorone	10.001	82	286967	29.395	_	99
25) 2-Nitrophenol	10.172	139	64287	31.952	-	99
26) 2,4-Dimethylphenol	10.219	107	124378	28.089		100
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.459	93	152538	28.999		97
29) 2,4-Dichlorophenol	10.712	162	107009	32.474	•	97
30) Naphthalene	11.117	128	354752	30.567		98
32) 4-Chloroaniline	11.229	127	147406	29.018	-	100
33) Hexachlorobutadiene	11.382	225	68445	31.647	•	99
34) Caprolactam	12.105	113	42075	30.076		97
35) 4-Chloro-3-methylphenol	12.334	107	129462	30.755	_	96

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

Data File: BG051005.D

Aca On : 12 Nov 2021 19:22

Operator : CG/JU Sample : PB140665BS

Misc

ALS Vial : 44 Sample Multiplier: 1

Quant Time: Nov 15 00:42:15 2021

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

Quant Title : SVOA CALIBRATION

Compound

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

R.T. QIon Response Conc Units Dev(Min) 36) 2-Methylnaphthalene 12.710 142 30.133 ng/ul 238244 37) 1-Methylnaphthalene 12.927 142 240969 30.078 ng/ul 98 39) 1,2,4,5-Tetrachloroben... 13.068 216 135229 33.797 ng/ul 40) Hexachlorocyclopentadiene 13.033 237 44022 22.911 ng/ul 41) 2,4,6-Trichlorophenol 13.309 196 95214 36.367 ng/ul 99 42) 2,4,5-Trichlorophenol 13.380 196 100919 35.903 ng/ul 96 43) 1,1'-Biphenyl 13.703 154 317083 31.591 ng/ul 98 44) 2-Chloronaphthalene 13.750 162 254394 32.343 ng/ul 98 45) 2-Nitroaniline 13.955 65 97263 31.124 ng/ul 90 47) Dimethylphthalate 31.280 ng/ul 14.314 163 328582 99 48) 2,6-Dinitrotoluene 165 32.473 ng/ul 14.437 71395 97 14.590 50) Acenaphthylene 152 410351 31.281 ng/ul 99 51) 3-Nitroaniline 14.772 138 72705 31.973 ng/ul 94 52) Acenaphthene 14.931 153 273812 31.743 ng/ul 97 53) 2,4-Dinitrophenol 14.984 184 41694 34.376 ng/ul 92 55) 4-Nitrophenol 15.078 109 59120 33.841 ng/ul 93 56) Dibenzofuran 15.266 168 386875 31.334 ng/ul 99 57) 2,4-Dinitrotoluene 15.224 165 100434 32.008 ng/ul# 99 58) 2,3,4,6-Tetrachlorophenol 15.489 232 83713 37.897 ng/ul 98 59) Diethylphthalate 15.665 149 348630 31.005 ng/ul 99 61) Fluorene 15.912 302896 166 30.995 ng/ul 99 62) 4-Chlorophenyl-phenyle... 15.894 204 164246 32.276 ng/ul 96 33.337 ng/ul 63) 4-Nitroaniline 15.935 138 75204 97 66) 4,6-Dinitro-2-methylph... 15.988 198 59340 34.042 ng/ul# 98 67) N-Nitrosodiphenylamine 16.106 169 274312 33.286 ng/ul 97 68) 4-Bromophenyl-phenylether 16.787 248 101821 34.720 ng/ul 95 69) Hexachlorobenzene 16.911 284 105938 35.138 ng/ul 70) Atrazine 17.058 200 107226 30.683 ng/ul 71) Pentachlorophenol 17.257 266 65268 47.153 ng/ul 72) Phenanthrene 17.657 178 505073 32.094 ng/ul 100 74) Anthracene 17.745 178 493013 31.223 ng/ul 99 75) 1,2,3,4-Tetrachloroben... 13.673 216 139551 34.764 ng/uL 97 76) Pentachlorobenzene 15.177 250 125330 33.695 ng/uL 97 77) Carbazole 18.015 167 32.457 ng/ul 459360 99 78) Di-n-butylphthalate 18.544 149 581184 31.251 ng/ul 100 80) Fluoranthene 19.655 202 568648 31.883 ng/ul 98 82) Pyrene 20.019 202 552737 31.717 ng/ul 96 83) Butylbenzylphthalate 20.877 149 228097 30.438 ng/ul 98 84) 3,3'-Dichlorobenzidine 21.799 252 178558 31.865 ng/ul 97 85) Benzo(a)anthracene 21.893 228 497543 31.235 ng/ul 98 86) Bis(2-ethylhexyl)phtha... 21.758 149 320678 29.811 ng/ul 99 87) Chrysene 21.964 228 481165 31.621 ng/ul 99 89) Di-n-octyl phthalate 23.033 149 29.612 ng/ul 543690 100 90) Benzo(b)fluoranthene 24.243 252 491292 30.579 ng/ul 98 91) Benzo(k)fluoranthene 24.314 252 473020 31.375 ng/ul 100 93) Benzo(a)pyrene 25.177 252 476786 31.159 ng/ul 99 94) Indeno(1,2,3-cd)pyrene 29.279 276 545004 31.995 ng/ul 96 29.343 278 95) Dibenzo(a,h)anthracene 468116 32.479 ng/ul 96 96) Benzo(g,h,i)perylene 30.512 276 451391 31.658 ng/ul 96

Instrument: BNA\_G

ClientSampleId:

SLCS665

## **Manual Integrations APPROVED**

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

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