# Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File: BG051336.D

Acq On : 3 Dec 2021 19:26

Operator : CG/JU Sample : M4833-10MSD

Misc

ALS Vial : 41 Sample Multiplier: 1

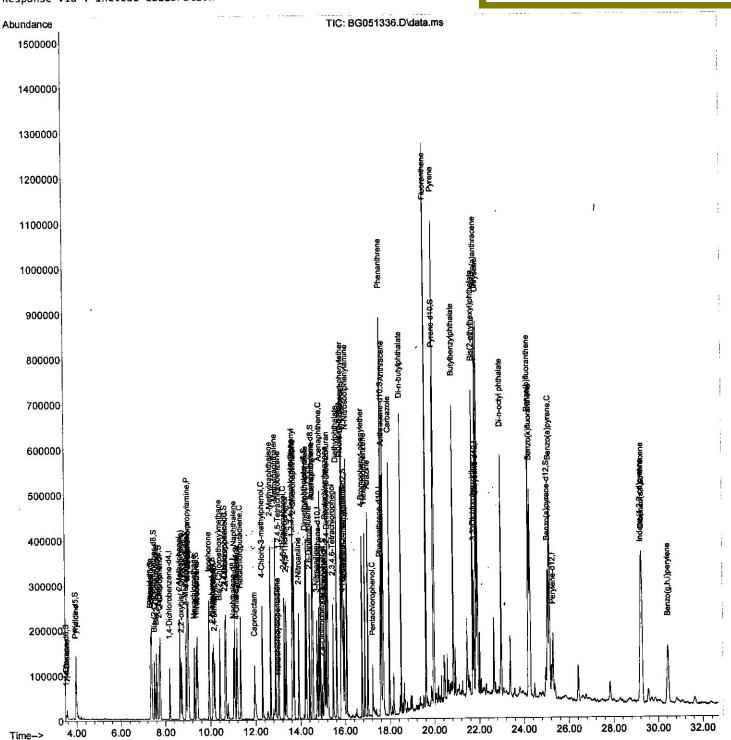
Quant Time: Dec 03 23:37:40 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument :
BNA\_G
ClientSampleId :
ESOM8MSD

### **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By :mohammad ahmed 12/07/2021



SFAM-EPA-BG112321.M Sat Dec 04 00:13:32 2021

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Misc

ALS Vial : 41 Sample Multiplier: 1

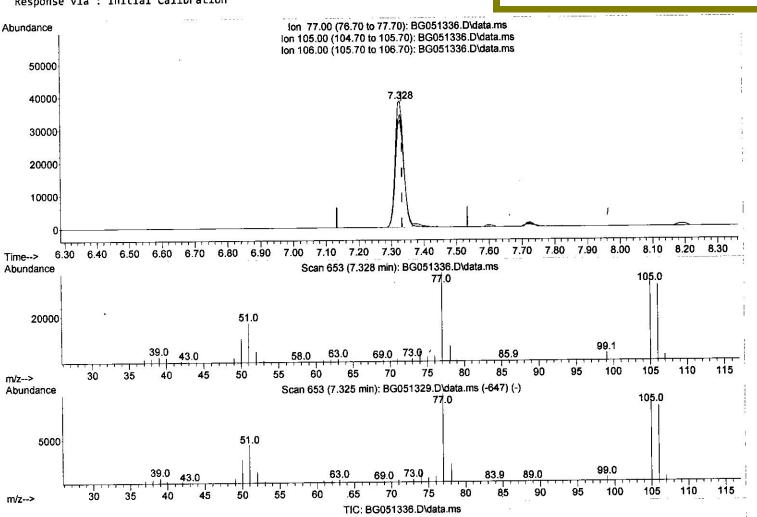
Quant Time: Dec 03 23:37:40 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration Instrument : BNA\_G ClientSampleId : ESQM8MSD

### **Manual IntegrationsAPPROVED**

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#### (6) Benzaldehyde

7.328min (-0.006) 36.85 ng/ul

response	69932	
Ion	Exp%	Act*
77.00	100.00	100.00
105.00	88.00	89.57
106.00	76.50	85.03
0.00	0.00	0.00

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Misc

Sample Multiplier: 1 : 41 ALS Vial

Quant Time: Dec 03 23:37:40 2021

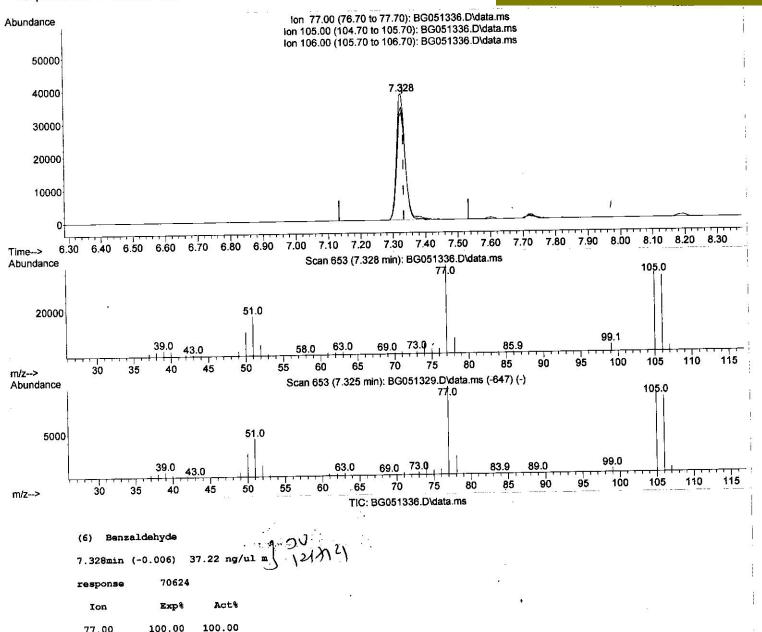
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100.00

88.00

76.50

0.00

89.57

85.03

0.00

77.00

105.00 106.00

0.00

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3 Dec 2021 19:26 Acq On

Operator : CG/JU : M4833-10MSD Sample

Misc

Sample Multiplier: 1 ALS Vial : 41

Quant Time: Dec 03 23:37:40 2021

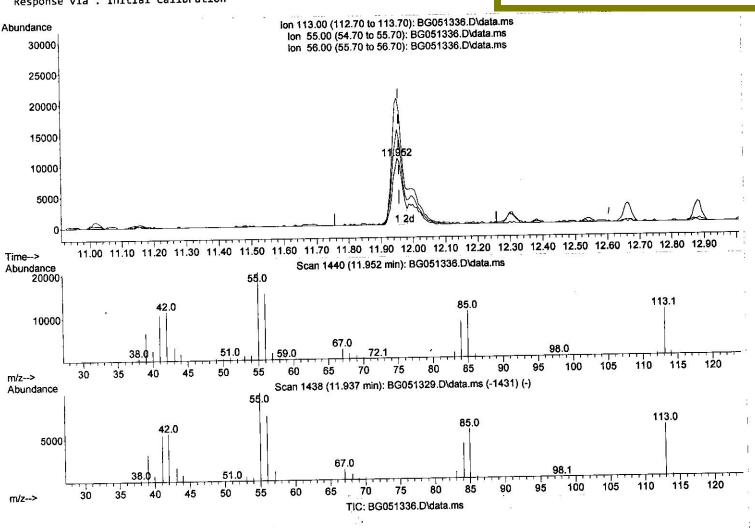
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Instrument: BNA\_G ClientSampleId: ESQM8MSD

### Manual IntegrationsAPPROVED

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#### (34) Caprolactam

28.40 ng/ul 11.952min (-0.006)

response	23410	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	183.80	188.93
56.00	136.50	142.36
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

: BG051336.D Data File

3 Dec 2021 19:26 Acq On

Operator : CG/JU : M4833-10MSD Sample

Misc ALS Vial

Sample Multiplier: 1 : 41

Quant Time: Dec 03 23:37:40 2021

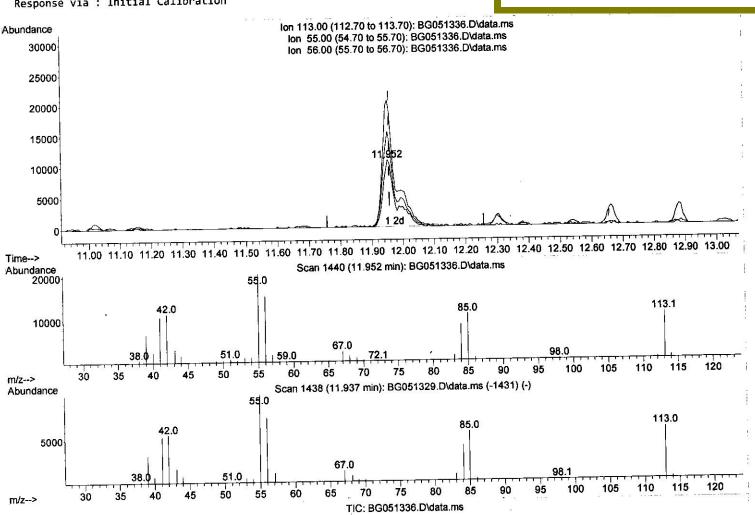
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Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via: Initial Calibration

Instrument: BNA\_G ClientSampleId: ESQM8MSD

### Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021 Supervised By :mohammad ahmed 12/07/2021



(34) Caprolactam 38.25 ng/ul 11.952min (-0.006)

response	31534		
Ion	Exp%	Act*	
113.00	100.00	100.00	
55.00	183.80	188.93	
56.00	136.50	142.36	
0.00	0.00	0.00	

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Misc

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Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Compound	R.T. Q	Ion	Response	Conc Units Dev(	Min)
Internal Standards				_	
1) 1,4-Dichlorobenzene-d4	8.191	152	30149	20.000 ng/ul	-0.01
20) Naphthalene-d8	11.018	136	131855	20.000 ng/ul	-0.01
38) Acenaphthene-d10	14.825	164	83549	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.575	188	178936	20.000 ng/ul	0.00
79) Chrysene-d12	21.881	240	149765	20.000 ng/ul	0.00
88) Perylene-d12	25.277	264	152645	20.000 ng/ul	0.00
System Monitoring Compounds		0 0		E 240 ==/ul	-0.01
3) 1,4-Dioxane-d8	3.532	96	4527	5.218 ng/uL	-0.02
4) Pyridine-d5	3.961	84	59329	23.305 ng/ul 33.155 ng/ul	0.00
7) Phenol-d5	7.351	99	98794	34.707 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth	7.510	67	64951	34.352 ng/ul	-0.01
11) 2-Chlorophenol-d4	7.721	132	73710 73980	30.767 ng/ul	0.00
15) 4-Methylphenol-d8	8.908	113 128	39798	35.756 ng/ul	0.00
21) Nitrobenzene-d5	9.372 10.095	143	46522	37.053 ng/ul	0.00
24) 2-Nitrophenol-d4	10.647	165	75877	35.618 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	11.159	131	70555	22.635 ng/ul	0.00
31) 4-Chloroaniline-d4	14.220	166	237721	36.979 ng/ul	0.00
46) Dimethylphthalate-d6	14.525	160	296016	36.516 ng/ul	0.00
49) Acenaphthylene-d8	15.048	143	35928	34.527 ng/ul	0.00
54) 4-Nitrophenol-d4 60) Fluorene-d10	15.818	176	212060	36.632 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.953	200	36038	32.639 ng/ul	0.00
73) Anthracene-d10	17.674	188	320898	37.497 ng/ul	0.00
81) Pyrene-d10	19.954	212	371490	40.995 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.048	264	320565	39.322 ng/ul	0.00
Target Compounds					alue
2) 1,4-Dioxane	3.573	88		11.558 ng/uL	94
5) Pyridine	3.985	79	72104		93
6) Benzaldehyde	7.328	7 <b>7</b>			
8) Phenol	7.381	94		36.539 ng/ul	98
10) Bis(2-Chloroethyl)ether	7.598	93		35.855 ng/ul	99
12) 2-Chlorophenol	7.757	128		35.765 ng/ul	100 99
13) 2-Methylphenol	8.638	108			98
14) 2,2'-oxybis(1-Chloropr	8.708			36.832 ng/ul	96
16) Acetophenone	9.020			36.137 ng/ul	99
17) N-Nitroso-di-n-propyla	8.996	70		35.706 ng/ul 84.557 ng/ul	95
<pre>18) 4-Methylphenol</pre>	8.973			85.037 ng/ul	92.
19) Hexachloroethane	9.278			39.010 ng/ul	94
22) Nitrobenzene	9.414			36.660 ng/ul	99
23) Isophorone	9.931	430		37.915 ng/ul	99
25) 2-Nitrophenol	10.124			18.611 ng/ul	97
26) 2,4-Dimethylphenol	10.177	1200		37.068 ng/ul	99
27) Bis(2-Chloroethoxy)met	. 10.406 10.671			36.390 ng/ul	98
29) 2,4-Dichlorophenol	11.070			37.584 ng/ul	98
30) Naphthalene	11.182			30.644 ng/ul	98
32) 4-Chloroaniline	11.335			35.477 ng/ul	96
33) Hexachlorobutadiene	11.952			The state of the s	
34) Caprolactam	12.304				94
35) 4-Chloro-3-methylphenol	12.304			=	

Instrument: BNA\_G ClientSampleId: ESQM8MSD

### **Manual IntegrationsAPPROVED**

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ALS Vial : 41 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION QLast Update : Fri Dec 03 15:23:09 2021 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units Dev(M	in)
36) 2-Methylnaphthalene	12.663	142	179428	36.768 ng/ul	98
37) 1-Methylnaphthalene	12.880	142	179941	35.840 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.027	216	101246	38.600 ng/ul	96
40) Hexachlorocyclopentadiene	12.992	237	8006	7.552 ng/ul	93
41) 2,4,6-Trichlorophenol	13.268	196	66223	40.233 ng/ul	96
42) 2,4,5-Trichlorophenol	13.356	196	71693	41.593 ng/ul	99
43) 1,1'-Biphenyl	13.656	154	236769	37.942 ng/ul	96
44) 2-Chloronaphthalene	13.709	162	189548	38.185 ng/ul	99
45) 2-Nitroaniline	13.920	65	72173	42.010 ng/ul	88
47) Dimethylphthalate	14.267	163	247858	38.091 ng/ul	100
48) 2,6-Dinitrotoluene	14.408	165	55594	40.673 ng/ul	94
50) Acenaphthylene	14.555	152	302737	37.800 ng/ul	98
51) 3-Nitroaniline	14.743	138	55070	40.760 ng/ul	97
52) Acenaphthene	14.889	153	208231	39.424 ng/ul	95
53) 2,4-Dinitrophenol	14.966	184	19902	26.342 ng/ul#	85
55) 4-Nitrophenol	15.066	109	36361	40.281 ng/ul	93
56) Dibenzofuran	15.224	168	290682	38.155 ng/ul	99
57) 2,4-Dinitrotoluene	15.201	165	81283	41.636 ng/ul	97
58) 2,3,4,6-Tetrachlorophenol	15.454	232	54666	40.387 ng/ul	99
59) Diethylphthalate	15.618	149	270984	39.674 ng/ul	99
61) Fluorene	15.871	166	243796	39.950 ng/ul	99
62) 4-Chlorophenyl-phenyle	15.853	204	123279	37.486 ng/ul	98
63) 4-Nitroaniline	15.906	138	58068	44.166 ng/ul	94
66) 4,6-Dinitro-2-methylph	15.965	198	40136	37.691 ng/ul	97
67) N-Nitrosodiphenylamine	16.070	169	214322	41.839 ng/ul	98
68) 4-Bromophenyl-phenylether	16.752	248	78522	40.944 ng/ul	93
69) Hexachlorobenzene	16.875	284	83456	42.678 ng/ul	99
70) Atrazine	17.016	200	85383	39.660 ng/ul	98
71) Pentachlorophenol	17.234	266	22056	25.454 ng/ul	94
72) Phenanthrene	17.622	178	571521	57.847 ng/ul	100
74) Anthracene	17.710	178	424867	43.300 ng/ul	98
75) 1,2,3,4-Tetrachloroben	13.632	216	103742	39.748 ng/uL	99
76) Pentachlorobenzene	15.142	250	96254	39.580 ng/uL	100
77) Carbazole	17.980	167	374088	<del></del>	99
78) Di-n-butylphthalate	18.509	149	459854	41.409 ng/ul	99
80) Fluoranthene	19.625	202	764945	68.728 ng/ul	97
82) Pyrene	19.989		698589	64.165 ng/ul	96
83) Butylbenzylphthalate	20.841			42.487 ng/ul	95
84) 3,3'-Dichlorobenzidine	21.764	252	100068	28.698 ng/ul	96
85) Benzo(a)anthracene	21.858	228	570443	56.158 ng/ul	98
<pre>86) Bis(2-ethylhexyl)phtha</pre>	21.717	149	278740	42.796 ng/ul	99
87) Chrysene	21.928	228	529992	The state of the s	98
89) Di-n-octyl phthalate	22.980	149	467923	42.313 ng/ul	100
90) Benzo(b)fluoranthene	24.196	252	561276	54.485 ng/ul	98
91) Benzo(k)fluoranthene	24.261	252	477002	49.343 ng/ul	98
93) Benzo(a)pyrene	25.125	252	518797	52.789 ng/ul	98
94) Indeno(1,2,3-cd)pyrene	29.214		543436	49.414 ng/ul	99
95) Dibenzo(a,h)anthracene	29.255		415710	44.556 ng/ul	97
96) Benzo(g,h,i)perylene	30.430		265786	28.725 ng/ul	95

### **Manual IntegrationsAPPROVED**

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Instrument: BNA\_G ClientSampleId: ESQM8MSD

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