Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File: BG050815.D

: 2 Nov 2021 12:52 Acq On

Operator : CG/JU Sample : SSTD16017

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 02 14:27:26 2021

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

Quant Title : SVOA CALIBRATION

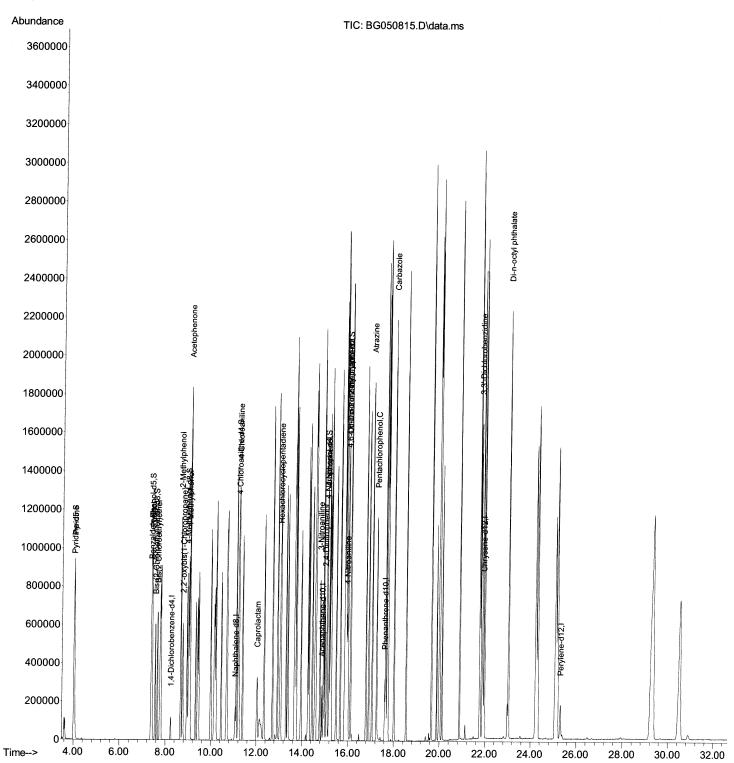
QLast Update : Tue Nov 02 14:25:39 2021

Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050815.D

Acq On : 2 Nov 2021 12:52

Operator : CG/JU Sample : SSTD16017

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 02 14:27:26 2021

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

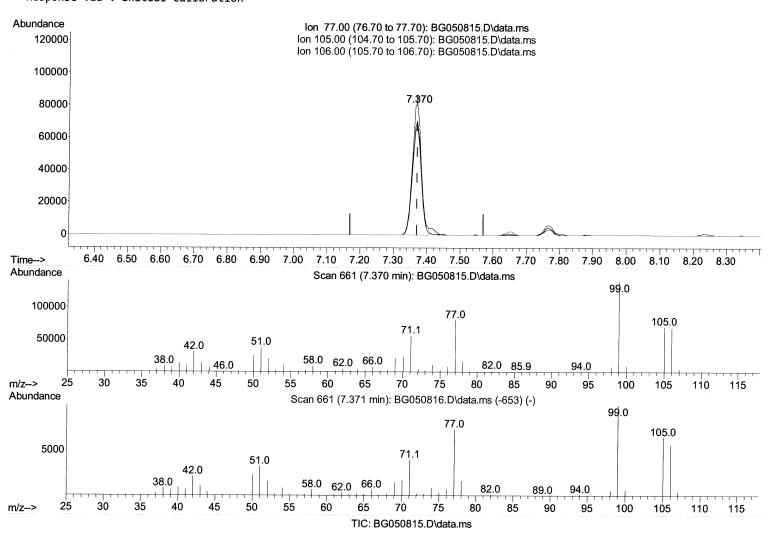
Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/02/2021 Supervised By: mohammad ahmed 11/08/2021



(6) Benzaldehyde

7.370min (-0.000) 71.39 ng/ul

response	151048	
Ion	Ехр%	Act%
77.00	100.00	100.00
105.00	88.00	87.11
106.00	76.50	84.51
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File: BG050815.D

Acq On : 2 Nov 2021 12:52

Operator : CG/JU Sample : SSTD16017

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 02 14:27:26 2021

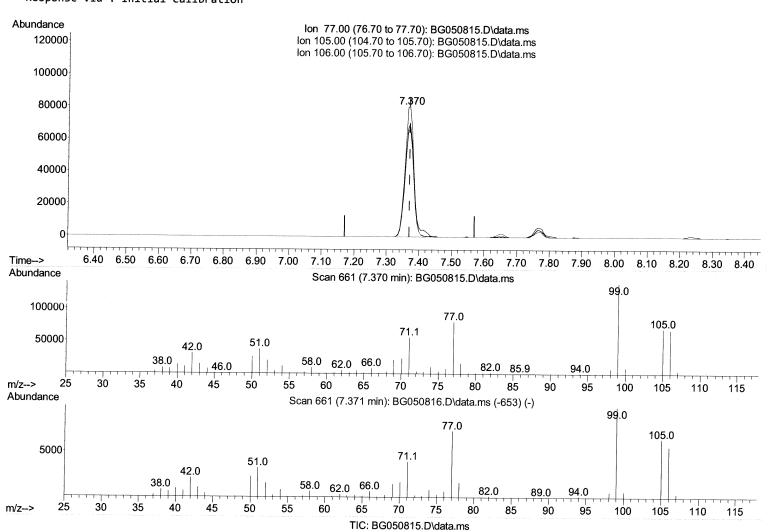
 $\label{lem:quant_method} \mbox{Quant Methods\SFAM-EPA-BG110321.M}$

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



(6) Benzaldehyde

7.370min (-0.000) 73.68 ng/ul m WOY/ALTA

response	155890	
Ion	Ехр%	Act%
77.00	100.00	100.00
105.00	88.00	87.11
106.00	76.50	84.51
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050815.D

Acq On : 2 Nov 2021 12:52

Operator : CG/JU Sample : SSTD16017

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 02 14:27:26 2021

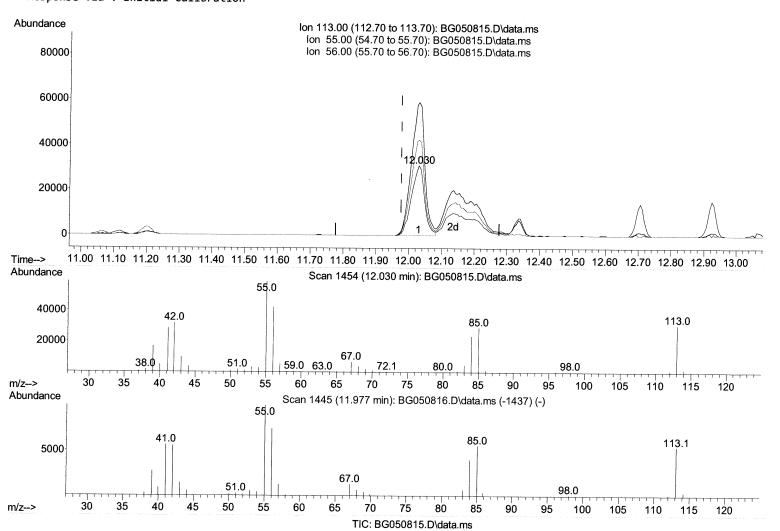
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



(34) Caprolactam

12.030min (+ 0.053) 86.21 ng/ul

response	84790		
Ion	Ехр%	Act%	
113.00	100.00	100.00	
55.00	183.80	190.67	
56.00	136.50	136.89	
0.00	0.00	0.00	

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File: BG050815.D

Acq On : 2 Nov 2021 12:52

Operator : CG/JU Sample : SSTD16017

Misc

ALS Vial : 7 Sample Multiplier: 1

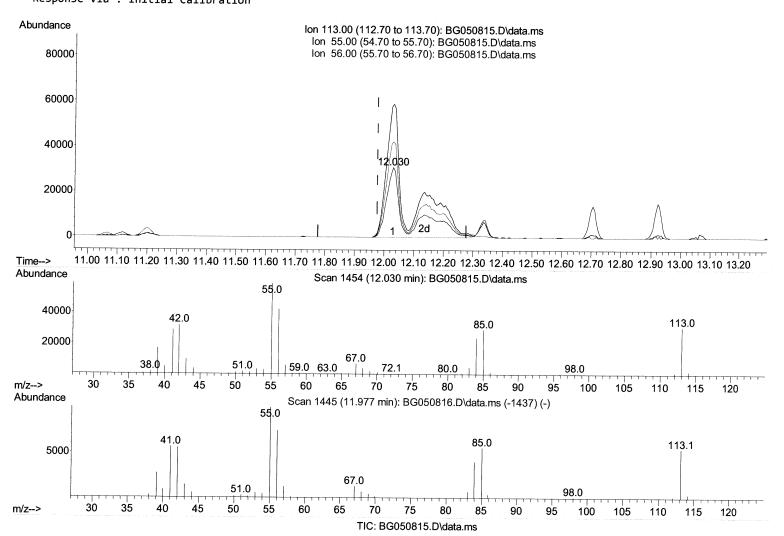
Quant Time: Nov 02 14:27:26 2021

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

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration Instrument :
BNA_G
ClientSampleId :
SSTD160417

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By :mohammad ahmed 11/08/2021



(34) Caprolactam

12.030min (+ 0.053) 153.62 ng/ul m Nloy/2134

response	151082			
Ion	Ехр%	Act%		
113.00	100.00	100.00		
55.00	183.80	190.67		
56.00	136.50	136.89		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050815.D

Acq On : 2 Nov 2021 12:52 Operator : CG/JU

Sample : SSTD16017

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 02 14:27:26 2021

 $\label{lem:quant_method} \mbox{Quant Methods} : \mbox{Z:\svoasrv\hPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M}$

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration

Instrument : BNA_G ClientSampleId : SSTD160417

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By:mohammad ahmed 11/08/2021

Internal Standards 1) 1,4-Dichlorobenzene-d4 2) Naphthalene-d8 3) 1,6-Dichlorobenzene-d8 3) 1,6-Dichlorobenzene-d9 4) Phenanthrene-d10 11,666 1366 149124 20,000 ng/ul 0,000 20,000 ng/ul 0,000 20,000 ng/ul 0,000 21,000 ng/ul 21,000 ng/ul 22,000 ng/ul 0,000 23,000 ng/ul 0,000 24,000 ng/ul 0,000 25,324 ng/ul 0,000 26,000 ng/ul 0,000 26,	Compound	R.T.		Response			(Min)
1) 1,4-Dichlorobenzene-d4 20) Naphthalene-d8 11.066 136 14.862 164 199218 20.000 ng/ul 0.00 040 14.862 164 17.611 188 126596 20.000 ng/ul 0.00 070 164) Phenanthrene-d10 17.611 188 126596 20.000 ng/ul 0.00 089) Perylene-d12 25.320 264 176739 20.000 ng/ul 0.00 089) Perylene-d12 25.320 264 176739 20.000 ng/ul 0.00 0.00 0.001 0.00 0.001 0.001 0.001 0.001 0.001 0.001 0.002 0.000 ng/ul 0.001 0.001 0.001 0.001 0.001 0.001 0.001 0.001 0.001 0.001 0.002 0.000 ng/ul 0.001 0	Internal Standards						
20) Naphthalene-d8 38) Acenaphthene-d10 14.862 164 99218 20.000 ng/ul 0.00 64) Phenanthrene-d10 17.611 188 216596 20.000 ng/ul 0.00 67) Chrysene-d12 21.912 240 176748 20.000 ng/ul 0.00 68) Perylene-d12 21.912 240 176748 20.000 ng/ul 0.00 69 60 60 60 60 60 60 60 60 60 60 60 60 60		8 234	152	31///3	20 000	na/ul	0 00
38) Acenaphthene-d10							
64) Phenanthrene-d10						-	
79 Chrysene-d12							
System Monitoring Compounds 3	· ·						
System Monitoring Compounds 3							
3) 1,4-Dioxane-d8 4) Pyridine-d5 4.015 84 450529 154.543 ng/ul 0.00 7) Phenol-d5 7) Bis-(2-Chloroethyl)eth 7.552 67 320003 147.681 ng/ul 0.00 11) 2-Chlorophenol-d4 0.000 120 0d 0.000 ng/ul 0.000 11) 2-Chlorophenol-d8 8.951 113 40546 153.534 ng/ul 0.02 21) Nitrobenzene-d5 0.000 128 0d 0.000 ng/ul 0.000 129 22-Nitrophenol-d4 0.000 128 0d 0.000 ng/ul 0.000 0g/ul 0.00						Q,	
A) Pyridine-dS	9 1						
7) Phenol-d5 9) Bis-(2-Chloroethyl)eth 7.552 67 320003 147.681 ng/ul 0.00 11) 2-Chlorophenol-d4 0.000 132 0d 0.000 ng/ul 15) 4-Methylphenol-d8 8.951 113 405446 153.534 ng/ul 0.02 21) Nitrobenzene-d5 0.000 128 0d 0.000 ng/ul 0.000 ng/ul 28 2,4-Dichlorophenol-d4 11.201 131 548711 152.651 ng/ul 0.01 28) 2,4-Dichlorophenol-d4 11.201 131 548711 152.651 ng/ul 0.01 31) 4-Chloroaniline-d4 11.201 131 548711 152.651 ng/ul 0.01 49) Acenaphthylene-d8 0.000 166 0d 0.000 ng/ul 0.01 49) Acenaphthylene-d8 0.000 160 0d 0.000 ng/ul 0.01 49) Acenaphthylene-d8 0.000 160 0d 0.000 ng/ul 0.03 65) 4,6-Dinitro-2-methylph 15.934 200 222291 169.278 ng/ul 0.03 66) Fluorene-d10 0.000 188 0d 0.000 ng/ul 0.02 73) Anthracene-d10 0.000 188 0d 0.000 ng/ul 0.02 73) Anthracene-d10 0.000 264 0d 0.000 ng/ul 0.02 74) Pyrene-d10 0.000 264 0d 0.000 ng/ul 0.		0.000	96		0.000	ng/uL	
9) Bis-(2-Chloroethyl)eth 7.552 67 320003 147.681 ng/ul 0.00 11) 2-Chlorophenol-d4 0.000 132 0d 0.000 ng/ul 15) 4-Methylphenol-d8 8.951 113 405446 153.534 ng/ul 0.02 21) Nitrobenzene-d5 0.000 128 0d 0.000 ng/ul 24) 2-Nitrophenol-d4 0.000 143 0d 0.000 ng/ul 28) 2,4-Dichlorophenol-d3 0.000 165 0d 0.000 ng/ul 31) 4-Chloroaniline-d4 11.201 131 548711 152.651 ng/ul 0.01 46) Dimethylphthalate-d6 0.000 166 0d 0.000 ng/ul 47) Acenaphthylpene-d8 0.000 160 0d 0.000 ng/ul 48) Acenaphthylpene-d8 0.000 160 0d 0.000 ng/ul 54) 4-Nitrophenol-d4 15.073 143 217725 158.192 ng/ul 0.03 60) Fluorene-d10 0.000 176 0d 0.000 ng/ul 61) 4-Nitrophenol-d4 15.073 143 217725 158.192 ng/ul 0.03 60) Fluorene-d10 0.000 176 0d 0.000 ng/ul 61) 4-Oinitro-2-methylph 15.984 200 222291 169.278 ng/ul 0.02 73) Anthracene-d10 0.000 212 0d 0.000 ng/ul 81) Pyrene-d10 0.000 264 0d 0.000 ng/ul 92) Benzo(a)pyrene-d12 0.000 264 0d 0.000 ng/ul 81) Pyrene-d10 0.000 264 0d 0.000 ng/ul 82) Phenol 7.417 94 529265 152.527 ng/ul 96 83 Benzaldehyde 7.370 77 155890m 73.682 ng/ul 98 84		4.015	84	450529	154.543	ng/ul	0.00
11) 2-Chlorophenol-d4		7.388	99	531401	158.419	ng/ul	0.01
15) 4-Methylphenol-d8		7.552	67	320003	147.681	ng/ul	0.00
21) Nitrobenzene-d5		0.000	132	0d	0.000	ng/ul	
24) 2-Nitrophenol-d4		8.951	113	405446	153.534	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	•	0.000	128	0d	0.000	ng/ul	
31) 4-Chloroaniline-d4 46) Dimethylphthalate-d6 47) Acenaphthylene-d8 48) 0.000 48) Acenaphthylene-d8 49) Acenaphthylene-d8 49) Acenaphthylene-d8 40) 0.000		0.000	143	0 d	0.000	ng/ul	
46) Dimethylphthalate-d6 49) Acenaphthylene-d8 49) Acenaphthylene-d8 40		0.000	165	0d	0.000	ng/ul	
49) Acenaphthylene-d8 40 000 160 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	· ·	11.201	131	548711	152 .651	ng/ul	0.01
15.073	- ·	0.000	166	0d	0.000	ng/ul	
60) Fluorene-d10		0.000	160	0d	0.000	ng/ul	
65) 4,6-Dinitro-2-methylph 15.984 200 222291 169.278 ng/ul 0.02 73) Anthracene-d10 0.000 188 0d 0.000 ng/ul 81) Pyrene-d10 0.000 212 0d 0.000 ng/ul 92) Benzo(a)pyrene-d12 0.000 264 0d 0.000 ng/ul 92) Benzo(a)pyrene-d12 0.000 264 0d 0.000 ng/ul 93) Pyridine 4.033 79 458599 151.462 ng/ul 96 94 6) Benzaldehyde 7.370 77 155890m 73.682 ng/ul 96 95 Pyridine 7.417 94 529265 152.527 ng/ul 96 10) Bis(2-Chloroethyl)ether 7.652 93 381142 146.745 ng/ul 98 13) 2-Methylphenol 8.675 108 400596 156.189 ng/ul 97 14) 2,2'-oxybis(1-Chloropr 8.763 45 597568 146.066 ng/ul 98 16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m 153.617 ng/ul 98 34) Caprolactam 12.030 113 151082m 153.617 ng/ul 98 35) 2,4-Dinitrophenol 14.783 138 206268 125.544 ng/ul 92 53) 2,4-Dinitrophenol 14.985 184 166308 189.774 ng/ul 93 55) 4-Nitroaniline 15.948 138 188568 115.690 ng/ul 93 63) 4-Nitrophenol 15.948 138 188568 115.690 ng/ul 95 66) 4,6-Dinitro-2-methylph 15.995 198 212426 165.879 ng/ul# 94 70) Atrazine 17.059 200 372072 144.920 ng/ul 98 71) Pentachlorophenol 17.253 266 204531 201.130 ng/ul 98 77) Carbazole 18.017 167 1432315 137.754 ng/ul 94 88) 3,3'-Dichlorobenzidine 21.795 252 544440 126.484 ng/ul 97 89) Di-n-octyl phthalate 23.028 149 1927616 133.968 ng/ul 100		15.073	143	217725	158.192	ng/ul	0.03
73) Anthracene-d10		0.000	176	0d	0.000	ng/ul	
81) Pyrene-d10	65) 4,6-Dinitro-2-methylph	15.984	200	222291	169.278	ng/ul	0.02
Target Compounds Qvalue O.000 264 Od O.000 ng/ul Ovalue	73) Anthracene-d10	0.000	188	0d	0.000	ng/ul	
Target Compounds 5) Pyridine 4.033 79 458599 151.462 ng/ul 96 6) Benzaldehyde 7.370 77 155890m > 73.682 ng/ul > \\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	81) Pyrene-d10	0.000	212	0d	0.000	ng/ul	
5) Pyridine 4.033 79 458599 151.462 ng/ul 96 6) Benzaldehyde 7.370 77 155890m > 73.682 ng/ul > NG4 3\J\ 8) Phenol 7.417 94 529265 152.527 ng/ul 96 10) Bis(2-Chloroethyl)ether 7.652 93 381142 146.745 ng/ul 98 13) 2-Methylphenol 8.675 108 400596 156.189 ng/ul 97 14) 2,2'-oxybis(1-Chloropr 8.763 45 597568 146.066 ng/ul 98 16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m >153.617 ng/ul > NO4/3\J\ 40) Hexachlorocyclopentadiene 13.040 237 261887 188.642 ng/ul 98 51) 3-Nitroaniline 14.773 138 206268 125.544 ng/ul 92 53) 2,4-Dinitrophenol 14.985 184 166308 189.774 ng/ul 93 55) 4-Nitrophenol 15.091 109 199297 157.889 ng/ul 93 63) 4-Nitroaniline 15.948 138 188568 115.690 ng/ul 95 66) 4,6-Dinitro-2-methylph 15.995 198 212426 165.879 ng/ul 98 71) Pentachlorophenol 17.253 266 204531 201.130 ng/ul 98 77) Carbazole 18.017 167 1432315 137.754 ng/ul 94 84) 3,3'-Dichlorobenzidine 21.795 252 544440 126.484 ng/ul 97 89) Di-n-octyl phthalate 23.028 149 1927616 133.968 ng/ul 100	92) Benzo(a)pyrene-d12	0.000	264	0d	0.000	ng/ul	
5) Pyridine 4.033 79 458599 151.462 ng/ul 96 6) Benzaldehyde 7.370 77 155890m > 73.682 ng/ul > NG4 3\J\ 8) Phenol 7.417 94 529265 152.527 ng/ul 96 10) Bis(2-Chloroethyl)ether 7.652 93 381142 146.745 ng/ul 98 13) 2-Methylphenol 8.675 108 400596 156.189 ng/ul 97 14) 2,2'-oxybis(1-Chloropr 8.763 45 597568 146.066 ng/ul 98 16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m >153.617 ng/ul > NO4/3\J\ 40) Hexachlorocyclopentadiene 13.040 237 261887 188.642 ng/ul 98 51) 3-Nitroaniline 14.773 138 206268 125.544 ng/ul 92 53) 2,4-Dinitrophenol 14.985 184 166308 189.774 ng/ul 93 55) 4-Nitrophenol 15.091 109 199297 157.889 ng/ul 93 63) 4-Nitroaniline 15.948 138 188568 115.690 ng/ul 95 66) 4,6-Dinitro-2-methylph 15.995 198 212426 165.879 ng/ul 98 71) Pentachlorophenol 17.253 266 204531 201.130 ng/ul 98 77) Carbazole 18.017 167 1432315 137.754 ng/ul 94 84) 3,3'-Dichlorobenzidine 21.795 252 544440 126.484 ng/ul 97 89) Di-n-octyl phthalate 23.028 149 1927616 133.968 ng/ul 100	Target Compounds					Ova	luo
6) Benzaldehyde 7.370 77 155890m > 73.682 ng/ul > 10 Phenol 7.417 94 529265 152.527 ng/ul 96 10 Bis(2-Chloroethyl)ether 7.652 93 381142 146.745 ng/ul 98 13) 2-Methylphenol 8.675 108 400596 156.189 ng/ul 97 14) 2,2'-oxybis(1-Chloropr 8.763 45 597568 146.066 ng/ul 98 16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m > 153.617 ng/ul > 10 U/A) JW 40 Hexachlorocyclopentadiene 13.040 237 261887 188.642 ng/ul 98 51) 3-Nitroaniline 14.773 138 206268 125.544 ng/ul 92 53) 2,4-Dinitrophenol 14.985 184 166308 189.774 ng/ul 93 55) 4-Nitrophenol 15.091 109 199297 157.889 ng/ul 93 63) 4-Nitroaniline 15.948 138 188568 115.690 ng/ul 95 66) 4,6-Dinitro-2-methylph 15.995 198 212426 165.879 ng/ul# 94 70) Atrazine 17.059 200 372072 144.920 ng/ul 98 71) Pentachlorophenol 17.253 266 204531 201.130 ng/ul 98 77) Carbazole 18.017 167 1432315 137.754 ng/ul 94 84) 3,3'-Dichlorobenzidine 21.795 252 544440 126.484 ng/ul 97 89) Di-n-octyl phthalate 23.028 149 1927616 133.968 ng/ul 100	-	4 033	79	158500	151 /62	-	
8) Phenol 7.417 94 529265 152.527 ng/ul 96 10) Bis(2-Chloroethyl)ether 7.652 93 381142 146.745 ng/ul 98 13) 2-Methylphenol 8.675 108 400596 156.189 ng/ul 97 14) 2,2'-oxybis(1-Chloropr 8.763 45 597568 146.066 ng/ul 98 16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m >153.617 ng/ul > 1104/31 JW 40) Hexachlorocyclopentadiene 13.040 237 261887 188.642 ng/ul 98 51) 3-Nitroaniline 14.773 138 206268 125.544 ng/ul 92 53) 2,4-Dinitrophenol 14.985 184 166308 189.774 ng/ul 93 55) 4-Nitrophenol 15.091 109 199297 157.889 ng/ul 93 63) 4-Nitroaniline 15.948 138 188568 115.690 ng/ul 95 66) 4,6-Dinitro-2-methylph 15.995 198 212426 165.879 ng/ul# 94 70) Atrazine 17.059 200 372072 144.920 ng/ul 98 71) Pentachlorophenol 17.253 266 204531 201.130 ng/ul 98 77) Carbazole 18.017 167 1432315 137.754 ng/ul 94 84) 3,3'-Dichlorobenzidine 21.795 252 544440 126.484 ng/ul 97 89) Di-n-octyl phthalate 23.028 149 1927616 133.968 ng/ul 100						-	
10) Bis(2-Chloroethyl)ether 7.652 93 381142 146.745 ng/ul 98 13) 2-Methylphenol 8.675 108 400596 156.189 ng/ul 97 14) 2,2'-oxybis(1-Chloropr 8.763 45 597568 146.066 ng/ul 98 16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m >153.617 ng/ul > 1104/31 JW 40) Hexachlorocyclopentadiene 13.040 237 261887 188.642 ng/ul 98 51) 3-Nitroaniline 14.773 138 206268 125.544 ng/ul 92 53) 2,4-Dinitrophenol 14.985 184 166308 189.774 ng/ul 93 55) 4-Nitrophenol 15.091 109 199297 157.889 ng/ul 93 63) 4-Nitroaniline 15.948 138 188568 115.690 ng/ul 95 66) 4,6-Dinitro-2-methylph 15.995 198 212426 165.879 ng/ul# 94 70) Atrazine 17.059 200 372072 144.920 ng/ul 98 71) Pentachlorophenol 17.253 266 204531 201.130 ng/ul 98 77) Carbazole 18.017 167 1432315 137.754 ng/ul 94 84) 3,3'-Dichlorobenzidine 21.795 252 544440 126.484 ng/ul 97 89) Di-n-octyl phthalate 23.028 149 1927616 133.968 ng/ul 100	•					_	
13) 2-Methylphenol 8.675 108 400596 156.189 ng/ul 97 14) 2,2'-oxybis(1-Chloropr 8.763 45 597568 146.066 ng/ul 98 16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m >153.617 ng/ul > \(\begin{align*} \begin{align*}	•						
14) 2,2'-oxybis(1-Chloropr 8.763 45 597568 146.066 ng/ul 98 16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m >153.617 ng/ul > WIOUIN JU 40) Hexachlorocyclopentadiene 13.040 237 261887 188.642 ng/ul 98 51) 3-Nitroaniline 14.773 138 206268 125.544 ng/ul 92 53) 2,4-Dinitrophenol 14.985 184 166308 189.774 ng/ul 93 55) 4-Nitrophenol 15.091 109 199297 157.889 ng/ul 93 63) 4-Nitroaniline 15.948 138 188568 115.690 ng/ul 95 66) 4,6-Dinitro-2-methylph 15.995 198 212426 165.879 ng/ul# 94 70) Atrazine 17.059 200 372072 144.920 ng/ul 98 71) Pentachlorophenol 17.253 266 204531 201.130 ng/ul 98 77) Carbazole 18.017 167 1432315 137.754 ng/ul 94 84) 3,3'-Dichlorobenzidine 21.795 252 544440 126.484 ng/ul 97 89) Di-n-octyl phthalate 23.028 149 1927616 133.968 ng/ul 100	the state of the s						
16) Acetophenone 9.074 105 603825 147.191 ng/ul 94 18) 4-Methylphenol 9.021 108 418688 153.316 ng/ul 96 32) 4-Chloroaniline 11.225 127 536348 150.269 ng/ul 98 34) Caprolactam 12.030 113 151082m >153.617 ng/ul > \\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\							
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89) Di-n-octyl phthalate 23.028 149 1927616 133.968 ng/ul 100	•					_	
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Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050815.D

Acq On : 2 Nov 2021 12:52 Operator : CG/JU Sample : SSTD16017

Misc

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 02 14:27:26 2021

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

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:25:39 2021 Response via : Initial Calibration

Compound R.T. QIon Response Conc Units Dev(Min) -----(#) = qualifier out of range (m) = manual integration (+) = signals summed

Instrument : BNA_G ClientSampleId :

SSTD160417

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/02/2021 Supervised By:mohammad ahmed 11/08/2021