







Data Path : Z:\svoasrv\HPCHEM1\ Data File : BG051414.D Acq On : 8 Dec 2021 19:45 Operator : CG/JU Sample : SSTD04031 Misc : ALS Vial : 6 Sample Multipli Quant Time: Dec 09 02:17:43 202 Quant Method : Z:\svoasrv\HPCHE Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 02:14	_ er: 1 !1 !M1\BNA_G	\Metho		A-BG120821.M	Instrument : BNA_G ClientSampleId : SSTD040431 Manual IntegrationsAPPROVED Reviewed By :Jagrut Upadhyay 12/09/2021 Supervised By :Yogesh Patel 12/16/2021	
Response via : Initial Calibrat						
Compound				Conc Units Dev		
Internal Standards						
 1,4-Dichlorobenzene-d4 	8.185	152	24564	20.00 ng/ul	0.00	
20) Naphthalene-d8	11.017	136	113431	20.00 ng/ul	0.00	
38) Acenaphthene-d10	14.824		75162	20.00 ng/ul	0.00	
64) Phenanthrene-d10	17.574		169655	20.00 ng/ul	0.00	
79) Chrysene-d12	21.875		141496	20.00 ng/ul	0.00	
88) Perylene-d12	25.277		141237	20.00 ng/ul	0.00	
	23.277	204	141207	20.00 hg/ui	0.00	
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.526	96	12017	17.00 ng/uL	0.00	
4) Pyridine-d5	3.961		87479	-		
7) Phenol-d5	7.357			42.17 ng/ul	0.00	
•			102004	42.02 ng/ul	0.00	
9) Bis-(2-Chloroethyl)eth			64827	42.52 ng/ul	0.00	
11) 2-Chlorophenol-d4	7.721		73769	42.20 ng/ul	0.00	
15) 4-Methylphenol-d8	8.914		79798	40.73 ng/ul	0.00	
21) Nitrobenzene-d5	9.366		39008	40.74 ng/ul	0.00	
24) 2-Nitrophenol-d4	10.095		45651	42.26 ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.647		75619	41.26 ng/ul	0.00	
31) 4-Chloroaniline-d4	11.158		106812	39.83 ng/ul	0.00	
46) Dimethylphthalate-d6	14.219		234346	40.52 ng/ul	0.00	
49) Acenaphthylene-d8	14.519		297079	40.74 ng/ul	0.00	
54) 4-Nitrophenol-d4	15.059		35508	37.93 ng/ul	0.00	
60) Fluorene-d10	15.811		208380	40.01 ng/ul	0.00	
65) 4,6-Dinitro-2-methylph			41961	40.08 ng/ul	0.00	
73) Anthracene-d10	17.674		320720	39.53 ng/ul	0.00	
81) Pyrene-d10	19.954		367150	42.88 ng/ul	0.00	
92) Benzo(a)pyrene-d12	25.048	264	310062	41.11 ng/ul	0.01	
Target Compounds				-	value	
2) 1,4-Dioxane	3.567	88	12891	16.17 ng/uL	95	
5) Pyridine	3.978	79	91105	42.21 ng/ul	95	
6) Benzaldehyde	7.327	77	70046	45.31 ng/ul	94	
8) Phenol	7.386	94	104393	41.51 ng/ul	99	
10) Bis(2-Chloroethyl)ether	7.597	93	79397	41.73 ng/ul	99	
<pre>12) 2-Chlorophenol</pre>	7.756	128	74793	41.98 ng/ul	98	
<pre>13) 2-Methylphenol</pre>	8.643	108	78203	41.75 ng/ul	97	
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.708	45	118891	43.30 ng/ul	96	
16) Acetophenone	9.019	105	126256	41.66 ng/ul	97	
17) N-Nitroso-di-n-propyla	8.996	70	74255	42.64 ng/ul	99	
18) 4-Methylphenol	8.972	108	83079	41.47 ng/ul	98	
19) Hexachloroethane	9.272	117	32034	42.57 ng/ul	97	
22) Nitrobenzene	9.413	77	105684	42.09 ng/ul	94	
23) Isophorone	9.930	82	206199	42.27 ng/ul	100	
25) 2-Nitrophenol	10.124	139	46041	41.15 ng/ul	99	
26) 2,4-Dimethylphenol	10.177	107	95535	41.77 ng/ul	99	
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.406	93	112047	41.61 ng/ul	98	
29) 2,4-Dichlorophenol	10.676	162	72806	40.36 ng/ul	98	
30) Naphthalene	11.070	128	247622	40.12 ng/ul	99	
32) 4-Chloroaniline	11.181	127	108014	40.12 ng/ul	98	
33) Hexachlorobutadiene	11.328	225	47530	38.20 ng/ul	98	
34) Caprolactam	11.957	113	29685m >		12/16/21 JU	
35) 4-Chloro-3-methylphenol	12.304	107	90191	41.62 ng/ul	99	

100

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Compound	R.T.	QIon	Response	Conc Units Dev	(Min)	
<pre>36) 2-Methylnaphthalene 37) 1-Methylnaphthalene 39) 1,2,4,5-Tetrachloroben 40) Hexachlorocyclopentadiene 41) 2,4,6-Trichlorophenol 42) 2,4,5-Trichlorophenol 43) 1,1'-Biphenyl 44) 2-Chloronaphthalene 45) 2-Nitroaniline 47) Dimethylphthalate 48) 2,6-Dinitrotoluene 50) Acenaphthylene 51) 3-Nitroaniline 52) Acenaphthene 53) 2,4-Dinitrophenol 55) 4-Nitrophenol 55) 4-Nitrophenol 56) Dibenzofuran 57) 2,4-Dinitrotoluene 58) 2,3,4,6-Tetrachlorophenol 59) Diethylphthalate 61) Fluorene 62) 4-Chlorophenyl-phenyle 63) 4-Nitroaniline 66) 4,6-Dinitro-2-methylph 67) N-Nitrosodiphenylamine 68) 4-Bromophenyl-phenylether 69) Hexachlorobenzene 70) Atrazine 71) Pentachlorophenol 72) Phenanthrene 74) Anthracene 75) 1,2,3,4-Tetrachloroben 76) Pentachlorobenzene 77) Carbazole</pre>	12.985 13.267 13.355 13.655 13.708 13.919 14.266 14.407 14.548 14.742 14.889 14.965 15.071 15.218 15.200 15.453 15.617 15.870 15.852 15.911 15.970 16.070 16.746 16.875 17.016 17.233 17.615 17.709 13.626 15.142 17.985	142 216 237 196 154 162 65 163 165 152 138 153 184 109 168 165 232 149 166 204 138 169 248 284 200 266 178 178 216 250 167	166685 172778 94083 39833 63171 64851 224693 177800 68811 234698 51838 288670 52301 191636 21338 31864 268935 72224 51778 250740 217178 114348 47910 40892 194567 71943 72510 83941 26557 374090 371073 99760 91448 335351	39.70 ng/ul 40.00 ng/ul 39.87 ng/ul 41.76 ng/ul 42.66 ng/ul 41.82 ng/ul 39.82 ng/ul 40.02 ng/ul 44.52 ng/ul 40.09 ng/ul 42.16 ng/ul 40.07 ng/ul 43.03 ng/ul 40.07 ng/ul 43.03 ng/ul 40.33 ng/ul 39.24 ng/ul 39.24 ng/ul 39.24 ng/ul 39.24 ng/ul 39.56 ng/ul 39.56 ng/ul 39.56 ng/ul 39.57 ng/ul	99 97 97 98 96 97 99 99 90 97 99 90 97 85 98 90 97 99 90 97 99 90 97 98 97 99 99 99 99 99 99 99 99 99 99 99 99	
 78) Di-n-butylphthalate 80) Fluoranthene 82) Pyrene 83) Butylbenzylphthalate 84) 3,3'-Dichlorobenzidine 85) Benzo(a)anthracene 86) Bis(2-ethylhexyl)phtha 87) Chrysene 	21.757 21.857 21.710 21.928	149 202 202 149 252 228 149 228	439930 454255 437555 187022 135889 397288 265648 377392	41.78 ng/ul 43.20 ng/ul 42.54 ng/ul 43.73 ng/ul 41.25 ng/ul 41.40 ng/ul 43.17 ng/ul 40.93 ng/ul	99 98 96 99 99 99 99	
<pre>89) Di-n-octyl phthalate 90) Benzo(b)fluoranthene 91) Benzo(k)fluoranthene 93) Benzo(a)pyrene 94) Indeno(1,2,3-cd)pyrene 95) Dibenzo(a,h)anthracene 96) Benzo(g,h,i)perylene</pre>	24.190 24.260 25.124 29.207 29.254	149 252 252 252 276 278 276	443384 393804 357946 369162 414272 349244 344931	43.33 ng/ul 41.32 ng/ul 40.02 ng/ul 40.60 ng/ul 40.71 ng/ul 40.46 ng/ul 40.29 ng/ul	100 98 99 98 98 97 98	

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