









Data Path : Z:\svoasr Data File : BG050895. Acq On : 9 Nov 20 Operator : CG/JU Sample : PB1404238 Misc : ALS Vial : 21 Samp Quant Time: Nov 09 01 Quant Method : Z:\svo Quant Title : SVOA C QLast Update : Tue No Response via : Initia	D 221 1:49 25 26 Multiplier: 1 2:30:18 2021 2035rv\HPCHEM1\BNA_G 24LIBRATION 20 02 14:49:05 2021	Instrument : BNA_G ClientSampleId : SLCS423 Manual IntegrationsAPPROVED Reviewed By :Jagrut Upadhyay 11/09/2021 Supervised By :mohammad ahmed 11/09/2021				
Compound	R.T.		•	Conc Units Dev		
Internal Standards						
1) 1,4-Dichloroben	zene-d4 8.227	152	38060	20.000 ng/ul	0.00	
20) Naphthalene-d8	11.053		168875	20.000 ng/ul	0.00	
38) Acenaphthene-d1			111610	20.000 ng/ul	-0.01	
64) Phenanthrene-d1			245836	20.000 ng/ul	0.00	
79) Chrysene-d12	21.893		205611	20.000 ng/ul	-0.01	
88) Perylene-d12	25.295		206459	20.000 ng/ul	-0.01	
bb) fel yiene uiz	25.255	204	200455	20.000 ng/ui	-0.01	
System Monitoring Co	mnounds					
3) 1,4-Dioxane-d8	3.586	96	7110	6.029 ng/uL	0.00	
4) Pyridine-d5	4.009	84	107030	30.339 ng/ul	0.00	
7) Phenol-d5	7.369	99	127181	31.323 ng/ul	0.00	
9) Bis-(2-Chloroet		67	82496	31.453 ng/ul	0.00	
11) 2-Chlorophenol-			89494	31.804 ng/ul	0.00	
15) 4-Methylphenol-			97437	30.482 ng/ul	0.00	
21) Nitrobenzene-d5			47398	33.027 ng/ul	0.00	
24) 2-Nitrophenol-d			51520	32.286 ng/ul	-0.01	
28) 2,4-Dichlorophe			87151	32.422 ng/ul	0.00	
31) 4-Chloroaniline			122316	30.048 ng/ul	0.00	
46) Dimethylphthala			281320	32.946 ng/ul	0.00	
49) Acenaphthylene-			352271	33.113 ng/ul	0.00	
54) 4-Nitrophenol-d			50473	32.600 ng/ul	0.00	
60) Fluorene-d10	15.842		242899	32.111 ng/ul	0.00	
65) 4,6-Dinitro-2-m			48039	32.231 ng/ul	0.00	
73) Anthracene-d10	17.698		370605	31.886 ng/ul	0.00	
81) Pyrene-d10	19.972	212	440142	33.143 ng/ul	0.00	
92) Benzo(a)pyrene-	d12 25.054	264	368764	32.310 ng/ul	-0.02	
				-		
Target Compounds				•	ralue	
2) 1,4-Dioxane	3.627	88	17451	13.473 ng/uL	97	
5) Pyridine	4.032	79	119749	32.792 ng/ul	97	
6) Benzaldehyde	7.363	77	95449	37.271 ng/ul	96	
8) Phenol	7.399	94	139247	33.152 ng/ul	98	
10) Bis(2-Chloroeth		93	108409	34.482 ng/ul	99	
12) 2-Chlorophenol	7.786	128	98140	34.350 ng/ul	97	
13) 2-Methylphenol	8.662	108	103179	33.235 ng/ul	96	
14) 2,2'-oxybis(1-C		45	166599	33.643 ng/ul	98	
16) Acetophenone	9.056	105	162689	32.763 ng/ul	97	
17) N-Nitroso-di-n-		70	96674	32.268 ng/ul	99	
18) 4-Methylphenol	8.991	108	109359	33.083 ng/ul	96	
19) Hexachloroethane		117	40679	34.054 ng/ul	95	
22) Nitrobenzene	9.443	77	141435	35.337 ng/ul	98	
23) Isophorone	9.960	82 130	266442	34.300 ng/ul	100	
25) 2-Nitrophenol	10.160	139	57094	35.664 ng/ul	99	
26) 2,4-Dimethylpher		107	110003	31.222 ng/ul	99	
27) Bis(2-Chloroetho		93 160	146708	35.052 ng/ul	98	
29) 2,4-Dichloropher		162	91863	35.036 ng/ul	96	
30) Naphthalene	11.106	128	320063	34.660 ng/ul	98	
32) 4-Chloroaniline	11.206	127	133981	33.147 ng/ul	100	
33) Hexachlorobutadi 34) Caprolactam		225 113	59728 36659m >	34.708 ng/ul	> 11/11/2/JJ	
35) 4-Chloro-3-methy	11.970 /lphenol 12.311	107	36659m≯ 114810	32.934 ng/ul ; 34.278 ng/ul	97	
		107	TT-0T0	24.270 ng/ui	57	

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\ Data File : BG050895.D Acq On : 9 Nov 2021 1:49 BNA_G	
Δcg. On 19 Nov 2021 1:49 BNA_G	
Operator : CG/JU ClientSampleId : SLCS423	
Misc : ALS Vial : 21 Sample Multiplier: 1 Manual IntegrationsAPPROVED	
Quant Time: Nov 09 01:30:18 2021 Reviewed By :Jagrut Upadhyay 11/09/2021	
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M Quant Title : SVOA CALIBRATION	
QLast Update : Tue Nov 02 14:49:05 2021	
Response via : Initial Calibration	
Compound R.T. QIon Response Conc Units Dev(Min)	
36) 2-Methylnaphthalene 12.692 142 212244 33.737 ng/ul 97	
37) 1-Methylnaphthalene 12.910 142 217363 34.098 ng/ul 96	
39) 1,2,4,5-Tetrachloroben 13.057 216 116623 35.861 ng/ul 97	
40) Hexachlorocyclopentadiene 13.027 237 46089 29.513 ng/ul 95 41) 2,4,6-Trichlorophenol 13.292 196 76529 35.963 ng/ul 98	
41) 2,4,6-Trichlorophenol 13.292 196 76529 35.963 ng/ul 98 42) 2,4,5-Trichlorophenol 13.362 196 83317 36.469 ng/ul 99	
43) 1,1'-Biphenyl 13.685 154 287123 35.196 ng/ul 99	
44) 2-Chloronaphthalene 13.738 162 225443 35.265 ng/ul 96	
45) 2-Nitroaniline 13.938 65 90391 35.588 ng/ul 95	
47) Dimethylphthalate 14.291 163 303068 35.497 ng/ul 99	
48) 2,6-Dinitrotoluene 14.426 165 65008 36.379 ng/ul 92	
50) Acenaphthylene 14.579 152 375526 35.220 ng/ul 99 51) 3-Nitroaniline 14.755 138 67397 36.466 ng/ul 95	
52) Acenaphthene 14.913 153 247392 35.287 ng/ul 96	
53) 2,4-Dinitrophenol 14.966 184 30787 31.230 ng/ul 94	
55) 4-Nitrophenol 15.054 109 49783 35.061 ng/ul 96	
56) Dibenzofuran 15.248 168 351348 35.011 ng/ul 100	
57) 2,4-Dinitrotoluene 15.207 165 94221 36.946 ng/ul 98	
58) 2, 3, 4,6-Tetrachlorophenol 15.472 232 63222 35.213 ng/ul 98 59) Diethylphthalate 15.648 149 321668 35.197 ng/ul 99	
59) Diethylphthalate 15.648 149 321668 35.197 ng/ul 99 61) Fluorene 15.895 166 277410 34.926 ng/ul 99	
62) 4-Chlorophenyl-phenyle 15.883 204 142509 34.455 ng/ul 97	
63) 4-Nitroaniline 15.918 138 68443 37.329 ng/ul 96	
66) 4,6-Dinitro-2-methylph 15.971 198 50761 34.924 ng/ul 98	
67) N-Nitrosodiphenylamine 16.094 169 246144 35.820 ng/ul 99	
68) 4-Bromophenyl-phenylether 16.776 248 88358 36.132 ng/ul 96	
69) Hexachlorobenzene 16.899 284 89846 35.739 ng/ul 97 70) Atrazine 17.034 200 99888 34.278 ng/ul 98	
71) Pentachlorophenol 17.246 266 41949 36.345 ng/ul 94	
72) Phenanthrene 17.640 178 473850 36.109 ng/ul 99	
74) Anthracene 17.734 178 454908 34.550 ng/ul 100	
75) 1,2,3,4-Tetrachloroben 13.656 216 119664 35.749 ng/uL 96	
76) Pentachlorobenzene 15.166 250 106911 34.471 ng/uL 100 77) Carbazole 17.998 167 433765 36.756 ng/ul 98	
77) Carbazole 17.998 167 433765 36.756 ng/ul 98 78) Di-n-butylphthalate 18.533 149 554592 35.764 ng/ul 99	
80) Fluoranthene 19.637 202 571278 35.844 ng/ul 99	
82) Pyrene 20.002 202 550187 35.330 ng/ul 99	
83) Butylbenzylphthalate 20.865 149 244079 36.448 ng/ul 99	
84) 3,3'-Dichlorobenzidine 21.782 252 171141 34.178 ng/ul 99	
85) Benzo(a)anthracene 21.876 228 511138 35.909 ng/ul 98	
86) Bis(2-ethylhexyl)phtha 21.741 149 345598 35.953 ng/ul 98 87) Chrysene 21.940 228 479288 35.248 ng/ul 100	
89) Di-n-octyl phthalate 23.010 149 593668 35.320 ng/ul 100	
90) Benzo(b)fluoranthene 24.208 252 524444 35.657 ng/ul 99	
91) Benzo(k)fluoranthene 24.279 252 466886 33.828 ng/ul 99	
93) Benzo(a)pyrene 25.131 252 488394 34.865 ng/ul 99	
94) Indeno(1,2,3-cd)pyrene 29.191 276 545998 35.014 ng/ul 97 95) Dibenzo(2,b)anthracene 29.273 278 458777 34.771 ng/ul 97	
95) Dibenzo(a,h)anthracene 29.273 278 458777 34.771 ng/ul 97 96) Benzo(g,h,i)perylene 30.425 276 453122 34.713 ng/ul 97	
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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