

8.00

10.00

12.00

14.00

16.00

18.00

20.00

22.00

24.00

26.00

6.00

40000₫

20000

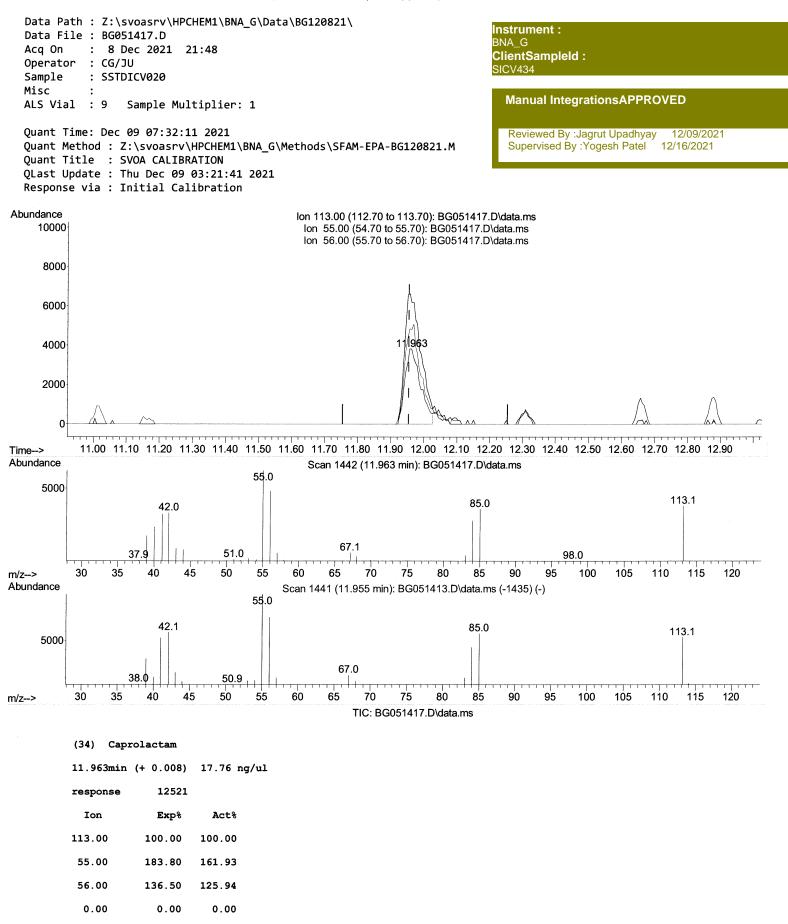
Time-->

30.00

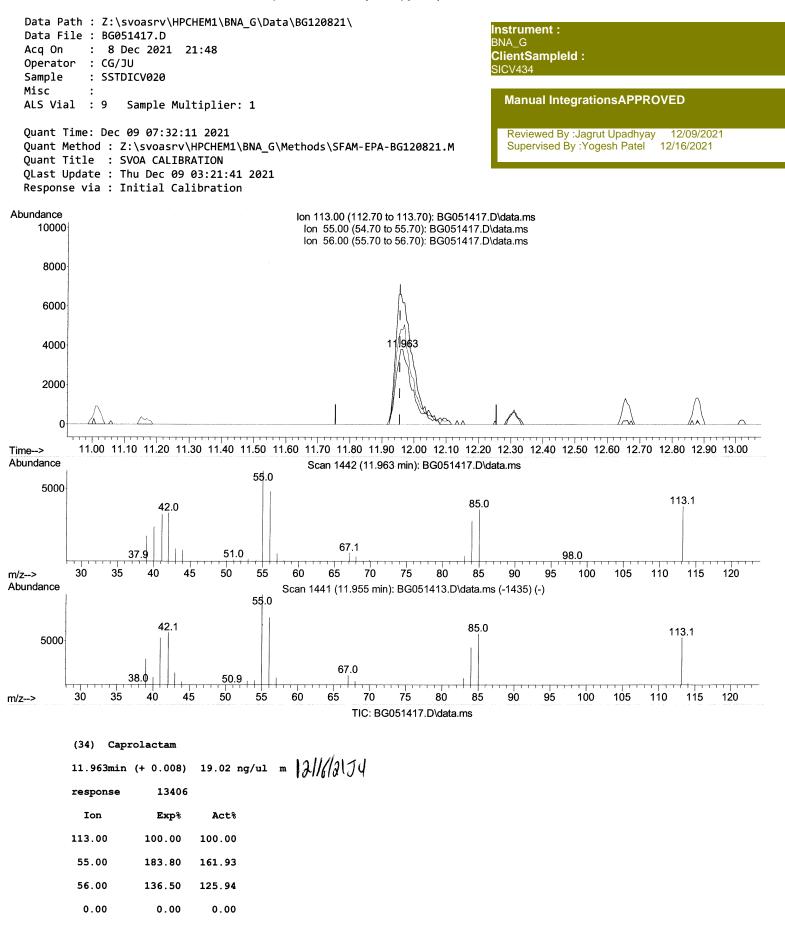
32.00

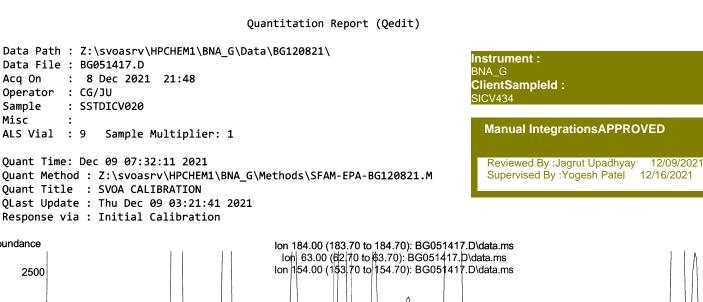
28.00

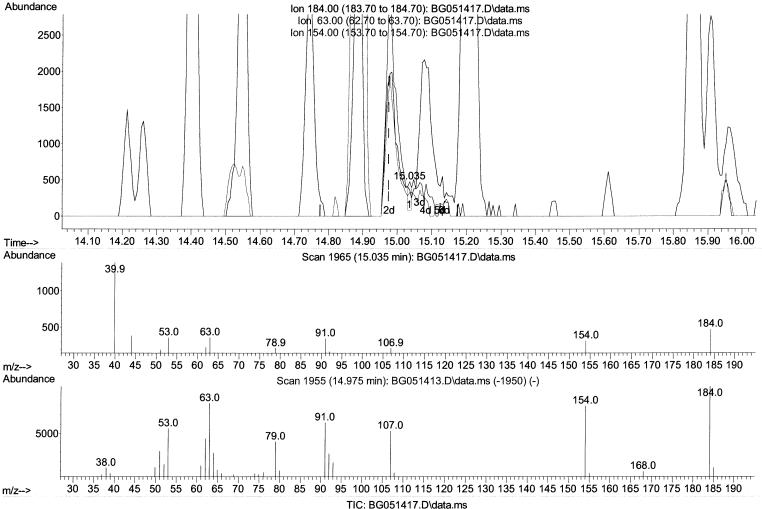












(53) 2,4-Dinitrophenol

Acq On

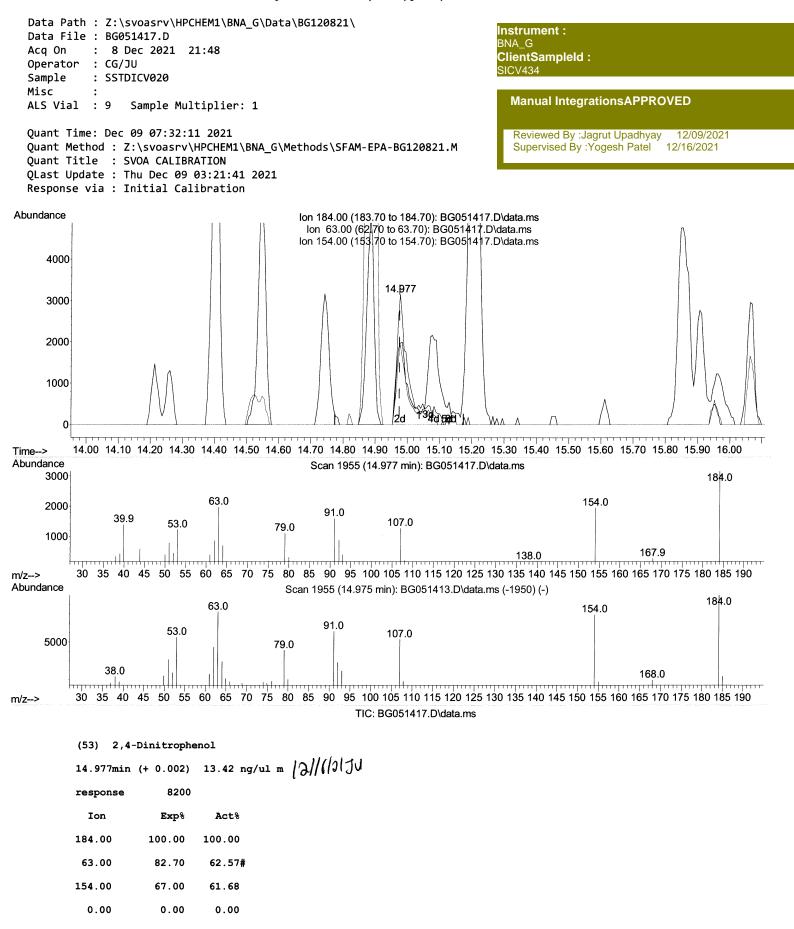
Sample Misc

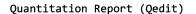
ALS Vial : 9

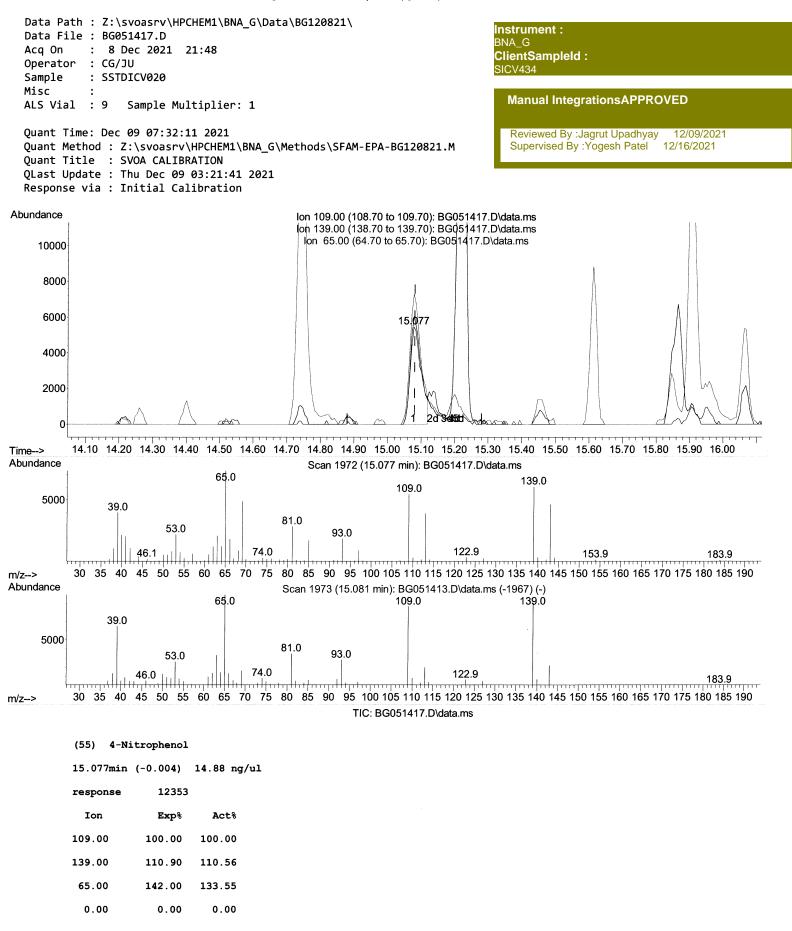
15.035min (+ 0.061) 0.41 ng/ul response 250 Ion Ехр% Act% 184.00 100.00 100.00 63.00 82.70 76.16 154.00 67.00 67.51 0.00 0.00 0.00

12/16/2021

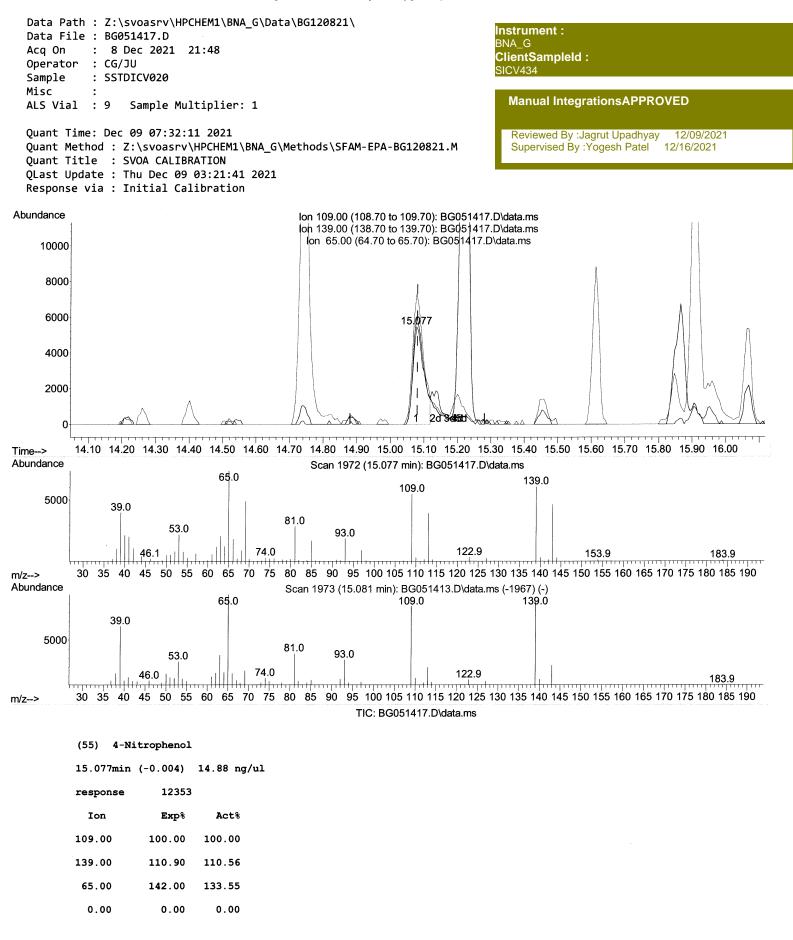




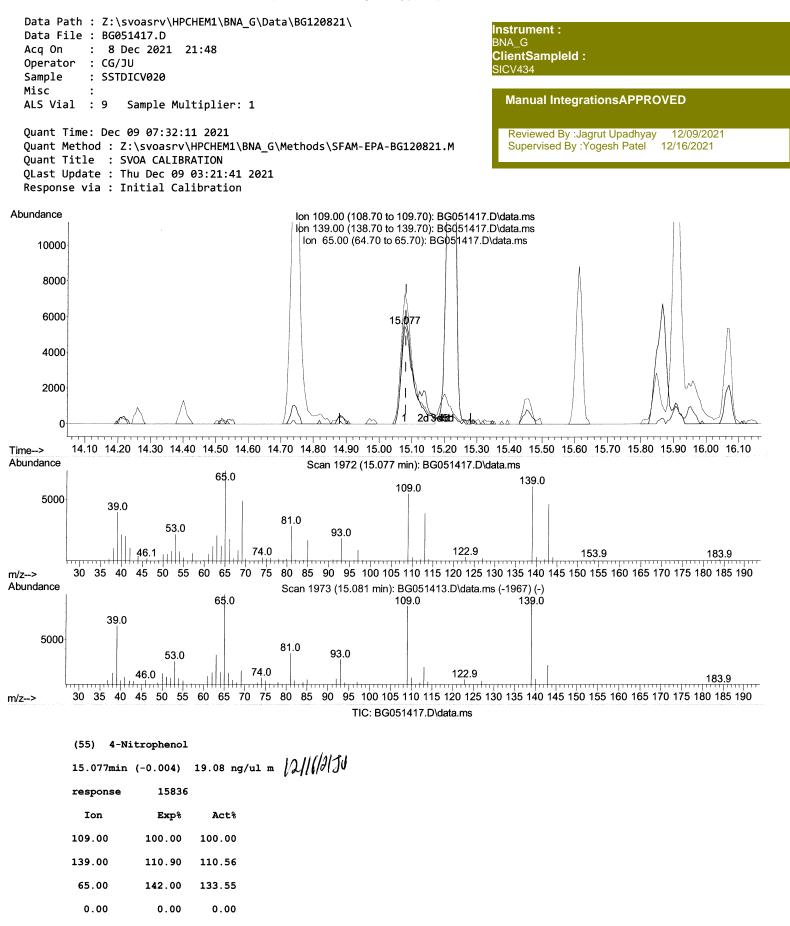


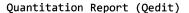


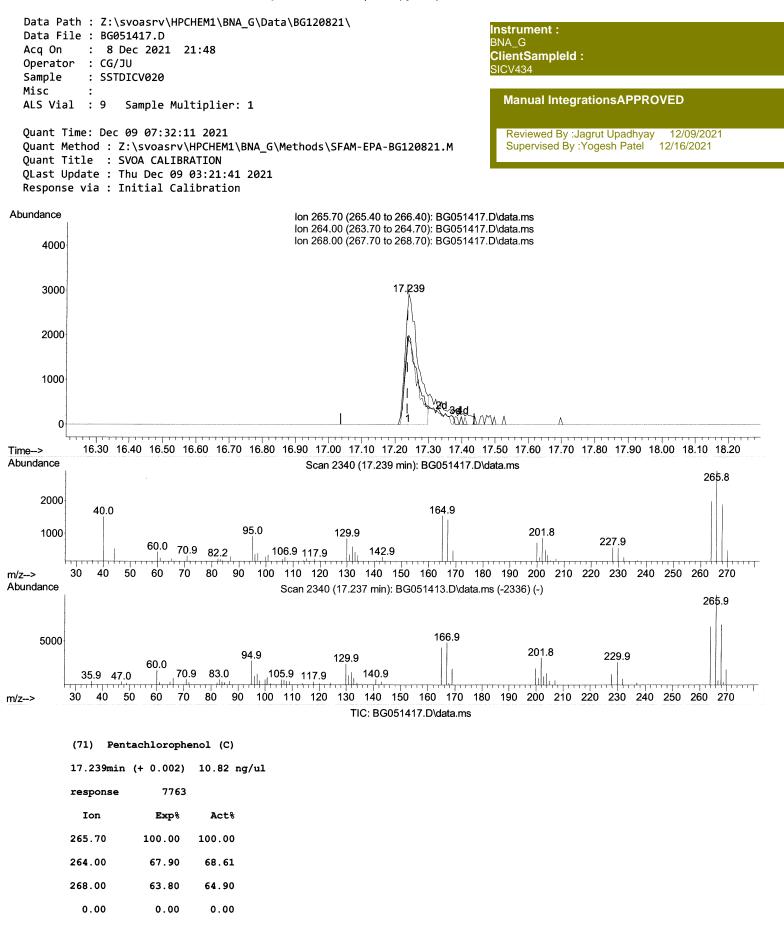


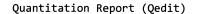


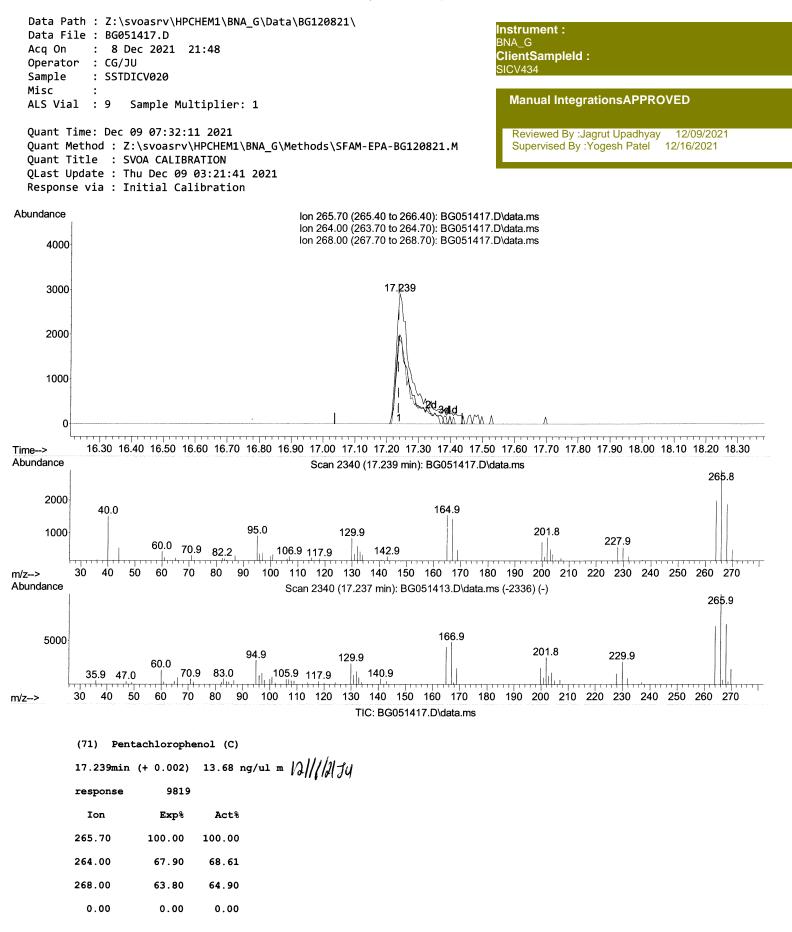












Data Path : Z:\svoasrv\HPCHEM1\ Data File : BG051417.D Acq On : 8 Dec 2021 21:48 Operator : CG/JU Sample : SSTDICV020 Misc : ALS Vial : 9 Sample Multiplic Quant Time: Dec 09 07:32:11 202: Quant Method : Z:\svoasrv\HPCHEM Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 03:21 Response via : Initial Calibrat:	_ er: 1 1 1\BNA_G` :41 2021			A-BG120821.M	Instrument : BNA_G ClientSampleId : SICV434 Manual IntegrationsAPPROVED Reviewed By :Jagrut Upadhyay 12/09/2021 Supervised By :Yogesh Patel 12/16/2021
Compound		-	•	Conc Units Dev	
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.185	152	24496	20.000 ng/ul	0.00
20) Naphthalene-d8	11.017		108668	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.818		71623	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.574		161075	20.000 ng/ul	0.00
79) Chrysene-d12	21.875		147999	20.000 ng/ul	0.00
88) Perylene-d12	25.271		146184	20.000 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.531	96	5599	7.506 ng/uL	0.00
4) Pyridine-d5	3.966	84	38829	18.127 ng/ul	0.00
7) Phenol-d5	7.356	99	45405	18.207 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth	7.503	67	29030	18.153 ng/ul	0.00
11) 2-Chlorophenol-d4	7.721	132	34177	19.263 ng/ul	0.00
15) 4-Methylphenol-d8	8.907	113	36044	18.399 ng/ul	0.00
21) Nitrobenzene-d5	9.366	128	17317	18.370 ng/ul	0.00
24) 2-Nitrophenol-d4	10.088	143	19606	18.380 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.647	165	32953	18.990 ng/ul	0.00
31) 4-Chloroaniline-d4	11.158	131	46776	18.429 ng/ul	0.00
<pre>46) Dimethylphthalate-d6</pre>	14.213	166	104525	18.860 ng/ul	0.00
49) Acenaphthylene-d8	14.518	160	135066	19.243 ng/ul	0.00
54) 4-Nitrophenol-d4	15.065	143	14141	16.947 ng/ul	0.00
60) Fluorene-d10	15.811		95134	19.283 ng/ul	0.00
65) 4,6-Dinitro-2-methylph			16942	17.700 ng/ul	0.00
73) Anthracene-d10	17.674	188 212	149077 177412	19.780 ng/ul 19.944 ng/ul	0.00 0.00
81) Pyrene-d10 92) Benzo(a)pyrene-d12	19.953		147969	19.625 ng/ul	0.00
	25.036	264	147909	19.025 lig/u1	0.00
Target Compounds				0va	alue
2) 1,4-Dioxane	3.567	88	6434	7.730 ng/uL	96
5) Pyridine	3.984	79	40790	18.243 ng/ul	97
6) Benzaldehyde	7.321	77	33390	21.062 ng/ul	97
8) Phenol	7.386	94	47542	18.626 ng/ul	97
10) Bis(2-Chloroethyl)ether	7.597	93	36323	18.579 ng/ul	99
<pre>12) 2-Chlorophenol</pre>	7.756	128	34518	18.993 ng/ul	94
13) 2-Methylphenol	8.643	108	34600	18.207 ng/ul	96
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.708	45	55739	18.947 ng/ul	98
16) Acetophenone	9.019	105	57262	18.876 ng/ul	96
17) N-Nitroso-di-n-propyla	8.990	70	34241	18.828 ng/ul	97
18) 4-Methylphenol	8.972	108	37200	18.634 ng/ul	98
19) Hexachloroethane	9.272	117 77	14457 48045	18.397 ng/ul	96 97
22) Nitrobenzene	9.407 9.924	82	92244	18.734 ng/ul	99
23) Isophorone 25) 2-Nitrophenol	9.924	82 139	20714	18.734 ng/ul 19.393 ng/ul	94
26) 2,4-Dimethylphenol	10.129	107	42724	18.888 ng/ul	98
27) Bis(2-Chloroethoxy)met	10.170	93	50498	18.937 ng/ul	97
29) 2,4-Dichlorophenol	10.400	162	32759	19.255 ng/ul	97
30) Naphthalene	11.064	128	114430	19.175 ng/ul	97
32) 4-Chloroaniline	11.004 11.181	127	47828	18.733 ng/ul	100
33) Hexachlorobutadiene	11.328	225	22080	19.026 ng/ul	96
34) Caprolactam	11.963	113	13406m >		
35) 4-Chloro-3-methylphenol	12.309	107	39405	18.646 ng/ul	97

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Data Path : Z:\svoasrv\HPCHEM1\ Data File : BG051417.D Acq On : 8 Dec 2021 21:48 Operator : CG/JU Sample : SSTDICV020 Misc : ALS Vial : 9 Sample Multipli		120821\	Instrument : BNA_G ClientSampleId : SICV434 Manual IntegrationsAPPROVED
Quant Time: Dec 09 07:32:11 202 Quant Method : Z:\svoasrv\HPCHE Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 03:21 Response via : Initial Calibrat	Reviewed By :Jagrut Upadhyay 12/09/2021 Supervised By :Yogesh Patel 12/16/2021		
Compound	R.T. QIon	Response Conc Units Dev(M	in)
<ul> <li>36) 2-Methylnaphthalene</li> <li>37) 1-Methylnaphthalene</li> <li>39) 1,2,4,5-Tetrachloroben</li> <li>40) Hexachlorocyclopentadiene</li> <li>41) 2,4,6-Trichlorophenol</li> <li>42) 2,4,5-Trichlorophenol</li> <li>43) 1,1'-Biphenyl</li> <li>44) 2-Chloronaphthalene</li> <li>45) 2-Nitroaniline</li> <li>47) Dimethylphthalate</li> <li>48) 2,6-Dinitrotoluene</li> </ul>	13.27319613.36119613.65515413.70816213.9196514.26016314.401165	7720918.849 ng/ul#4279519.189 ng/ul1608113.617 ng/ul2711818.823 ng/ul2864018.569 ng/ul10301219.238 ng/ul8228219.585 ng/ul2996418.869 ng/ul10694019.146 ng/ul2273919.239 ng/ul	99 96 98 94 92 98 96 99 99 96
<ul> <li>50) Acenaphthylene</li> <li>51) 3-Nitroaniline</li> <li>52) Acenaphthene</li> <li>53) 2,4-Dinitrophenol</li> <li>55) 4-Nitrophenol</li> <li>56) Dibenzofuran</li> <li>57) 2,4-Dinitrotoluene</li> </ul>	14.54815214.74213814.88315314.97718415.07710915.21816815.200165	133272 19.242 ng/ul 23355 20.519 ng/ul 87326 19.216 ng/ul 15836m 13.415 ng/ul 123955 19.238 ng/ul 32153 19.032 ng/ul	98 99 97 12///21 JU 97 99
<ul> <li>58) 2,3,4,6-Tetrachlorophenol</li> <li>59) Diethylphthalate</li> <li>61) Fluorene</li> <li>62) 4-Chlorophenyl-phenyle</li> <li>63) 4-Nitroaniline</li> <li>66) 4,6-Dinitro-2-methylph</li> <li>67) N-Nitrosodiphenylamine</li> </ul>	15.453 232 15.611 149 15.870 166 15.846 204 15.905 138 15.970 198 16.070 169	2126118.193 ng/ul11274818.703 ng/ul9924719.020 ng/ul5265019.220 ng/ul2189321.668 ng/ul1648217.711 ng/ul8718719.417 ng/ul	100 99 98 98 96 97 97
<ul> <li>68) 4-Bromophenyl-phenylether</li> <li>69) Hexachlorobenzene</li> <li>70) Atrazine</li> <li>71) Pentachlorophenol</li> <li>72) Phenanthrene</li> <li>74) Anthracene</li> </ul>	16.74524816.87428417.01020017.23926617.61517817.709178	3132419.260 ng/ul3208319.353 ng/ul3818819.710 ng/ul9819m13.681 ng/ul17171219.786 ng/ul17375620.000 ng/ul	93 97 18/16/21 JU 99 98
<ul> <li>75) 1,2,3,4-Tetrachloroben</li> <li>76) Pentachlorobenzene</li> <li>77) Carbazole</li> <li>78) Di-n-butylphthalate</li> <li>80) Fluoranthene</li> <li>82) Pyrene</li> </ul>	13.625         216           15.135         250           17.985         167           18.502         149           19.618         202           19.983         202	45292       20.108 ng/uL         41499       20.350 ng/uL         157431       20.357 ng/ul         201346       19.419 ng/ul         216055       19.726 ng/ul         212410       19.762 ng/ul	99 97 98 99 97 97
<ul> <li>83) Butylbenzylphthalate</li> <li>84) 3,3'-Dichlorobenzidine</li> <li>85) Benzo(a)anthracene</li> <li>86) Bis(2-ethylhexyl)phtha</li> <li>87) Chrysene</li> <li>89) Di-n-octyl phthalate</li> </ul>	20.840 149 21.757 252 21.851 228 21.704 149 21.922 228 22.967 149	90372       19.278 ng/ul         67244       21.516 ng/ul         194451       19.890 ng/ul         128376       19.716 ng/ul         184458       19.788 ng/ul         216003       20.092 ng/ul	97 97 99 98 98 100
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	24.184       252         24.254       252         25.112       252         29.195       276         29.242       278         30.429       276	187783       19.556 ng/ul         176556       19.742 ng/ul         180490       19.735 ng/ul         198154       19.522 ng/ul         165245       19.313 ng/ul         163493       19.261 ng/ul	99 96 99 99 97 93

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