

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121421\  
 Data File : BG051502.D  
 Acq On : 14 Dec 2021 13:30  
 Operator : CG/JU  
 Sample : SSTD16040  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Dec 14 14:04:34 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 14 13:31:18 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.287  | 152  | 33260    | 20.000  | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 11.130 | 136  | 135009   | 20.000  | ng/ul  | # 0.00   |
| 38) Acenaphthene-d10          | 14.919 | 164  | 96355    | 20.000  | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.668 | 188  | 228376   | 20.000  | ng/ul  | # 0.00   |
| 79) Chrysene-d12              | 21.998 | 240  | 202455   | 20.000  | ng/ul  | # 0.01   |
| 88) Perylene-d12              | 25.499 | 264  | 207630   | 20.000  | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000   | ng/uL  |          |
| 4) Pyridine-d5                | 4.046  | 84   | 378836m  | 160.288 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.429  | 99   | 470703   | 155.121 | ng/ul  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.600  | 67   | 260452   | 144.089 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/ul  |          |
| 15) 4-Methylphenol-d8         | 8.998  | 113  | 354690   | 144.298 | 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        | 11.254 | 131  | 472568   | 148.316 | ng/ul  | 0.01     |
| 46) Dimethylphthalate-d6      | 0.000  | 166  | 0d       | 0.000   | ng/ul  |          |
| 49) Acenaphthylene-d8         | 0.000  | 160  | 0d       | 0.000   | ng/ul  |          |
| 54) 4-Nitrophenol-d4          | 15.107 | 143  | 193767   | 168.104 | ng/ul  | 0.03     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/ul  |          |
| 65) 4,6-Dinitro-2-methylph... | 16.029 | 200  | 212914   | 204.573 | 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  |          |
| Target Compounds              |        |      |          |         |        |          |
|                               |        |      |          |         | Qvalue |          |
| 5) Pyridine                   | 4.063  | 79   | 371103   | 150.722 | ng/ul  | 94       |
| 6) Benzaldehyde               | 7.406  | 77   | 109687m  | 60.054  | ng/ul  |          |
| 8) Phenol                     | 7.459  | 94   | 451835   | 149.753 | ng/ul# | 89       |
| 10) Bis(2-Chloroethyl)ether   | 7.694  | 93   | 327052   | 142.793 | ng/ul  | 97       |
| 13) 2-Methylphenol            | 8.722  | 108  | 345663   | 148.895 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.822  | 45   | 440011   | 139.225 | ng/ul  | 95       |
| 16) Acetophenone              | 9.127  | 105  | 526607   | 137.347 | ng/ul  | 98       |
| 18) 4-Methylphenol            | 9.068  | 108  | 357569   | 142.572 | ng/ul  | 94       |
| 32) 4-Chloroaniline           | 11.283 | 127  | 473926   | 147.063 | ng/ul  | 99       |
| 34) Caprolactam               | 12.088 | 113  | 111452m  | 161.807 | ng/ul  |          |
| 40) Hexachlorocyclopentadiene | 13.116 | 237  | 385020   | 162.777 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.819 | 138  | 170935   | 136.228 | ng/ul  | 96       |
| 53) 2,4-Dinitrophenol         | 15.019 | 184  | 148723   | 217.575 | ng/ul  | 90       |
| 55) 4-Nitrophenol             | 15.125 | 109  | 207477   | 170.264 | ng/ul  | 86       |
| 63) 4-Nitroaniline            | 15.988 | 138  | 159572   | 123.667 | ng/ul  | 86       |
| 66) 4,6-Dinitro-2-methylph... | 16.041 | 198  | 204446   | 195.088 | ng/ul# | 99       |
| 70) Atrazine                  | 17.116 | 200  | 413349   | 143.946 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 17.316 | 266  | 300976   | 167.325 | ng/ul  | 97       |
| 77) Carbazole                 | 18.068 | 167  | 1487085  | 136.816 | ng/ul  | 95       |
| 84) 3,3'-Dichlorobenzidine    | 21.874 | 252  | 637537   | 127.045 | ng/ul# | 96       |
| 89) Di-n-octyl phthalate      | 23.190 | 149  | 1661993  | 143.096 | ng/ul  | 100      |
| -----                         |        |      |          |         |        |          |

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



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