

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042823\  
 Data File : BG057217.D  
 Acq On : 28 Apr 2023 14:02  
 Operator : CG/JU  
 Sample : SSTD16085  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD160433

Quant Time: Apr 28 23:11:44 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG042823.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 28 14:45:43 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Christian Giraldo 05/01/2023  
 Supervised By :Jagrut Upadhyay 05/01/2023

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.379  | 152  | 16574    | 20.000  | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 11.223 | 136  | 70678    | 20.000  | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.994 | 164  | 52856    | 20.000  | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.731 | 188  | 125204   | 20.000  | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 22.049 | 240  | 111062   | 20.000  | ng/u1  | # 0.01   |
| 88) Perylene-d12              | 25.556 | 264  | 117274   | 20.000  | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000   | ng/uL  |          |
| 4) Pyridine-d5                | 4.097  | 84   | 181985   | 153.262 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.522  | 99   | 250945   | 160.101 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.692  | 67   | 138113   | 149.812 | ng/u1  | 0.01     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/u1  |          |
| 15) 4-Methylphenol-d8         | 9.096  | 113  | 187281   | 157.208 | ng/u1  | 0.02     |
| 21) Nitrobenzene-d5           | 0.000  | 128  | 0d       | 0.000   | ng/u1  |          |
| 24) 2-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000   | ng/u1  |          |
| 28) 2,4-Dichlorophenol-d3     | 0.000  | 165  | 0d       | 0.000   | ng/u1  |          |
| 31) 4-Chloroaniline-d4        | 11.352 | 131  | 248299   | 147.827 | ng/u1  | 0.01     |
| 46) Dimethylphthalate-d6      | 0.000  | 166  | 0d       | 0.000   | ng/u1  |          |
| 49) Acenaphthylene-d8         | 0.000  | 160  | 0d       | 0.000   | ng/u1  |          |
| 54) 4-Nitrophenol-d4          | 15.194 | 143  | 95578    | 158.011 | ng/u1  | 0.03     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/u1  |          |
| 65) 4,6-Dinitro-2-methylph... | 16.104 | 200  | 124941   | 168.426 | ng/u1  | 0.02     |
| 73) Anthracene-d10            | 0.000  | 188  | 0d       | 0.000   | ng/u1  |          |
| 81) Pyrene-d10                | 0.000  | 212  | 0d       | 0.000   | ng/u1  |          |
| 92) Benzo(a)pyrene-d12        | 0.000  | 264  | 0d       | 0.000   | ng/u1  |          |
| Target Compounds              |        |      |          |         |        |          |
|                               |        |      |          |         |        | Qvalue   |
| 5) Pyridine                   | 4.115  | 79   | 188953   | 150.958 | ng/u1  | 96       |
| 6) Benzaldehyde               | 7.504  | 77   | 63213m   | 132.045 | ng/u1  |          |
| 8) Phenol                     | 7.551  | 94   | 248254   | 157.143 | ng/u1  | 95       |
| 10) Bis(2-Chloroethyl)ether   | 7.786  | 93   | 175713   | 148.611 | ng/u1  | 97       |
| 13) 2-Methylphenol            | 8.820  | 108  | 189137   | 154.481 | ng/u1  | 90       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.902  | 45   | 223580   | 146.450 | ng/u1  | 97       |
| 16) Acetophenone              | 9.219  | 105  | 297193   | 144.978 | ng/u1  | 100      |
| 18) 4-Methylphenol            | 9.166  | 108  | 202235   | 152.603 | ng/u1  | 95       |
| 32) 4-Chloroaniline           | 11.381 | 127  | 254716   | 145.541 | ng/u1  | 97       |
| 34) Caprolactam               | 12.174 | 113  | 59501m   | 147.545 | ng/u1  |          |
| 40) Hexachlorocyclopentadiene | 13.185 | 237  | 195697   | 204.163 | ng/u1  | 98       |
| 51) 3-Nitroaniline            | 14.900 | 138  | 95050    | 154.986 | ng/u1  | 100      |
| 53) 2,4-Dinitrophenol         | 15.106 | 184  | 88490    | 191.980 | ng/u1  | 89       |
| 55) 4-Nitrophenol             | 15.211 | 109  | 110782   | 160.303 | ng/u1  | 94       |
| 63) 4-Nitroaniline            | 16.069 | 138  | 90681    | 156.139 | ng/u1  | 89       |
| 66) 4,6-Dinitro-2-methylph... | 16.122 | 198  | 126753   | 167.683 | ng/u1# | 99       |
| 70) Atrazine                  | 17.173 | 200  | 218497   | 150.317 | ng/u1  | 99       |
| 71) Pentachlorophenol         | 17.379 | 266  | 134912   | 187.268 | ng/u1  | 97       |
| 77) Carbazole                 | 18.137 | 167  | 800761   | 146.304 | ng/u1  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.920 | 252  | 334581   | 135.648 | ng/u1  | 97       |
| 89) Di-n-octyl phthalate      | 23.171 | 149  | 1008234  | 155.685 | ng/u1  | 100      |

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