

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG091023\  
 Data File : BG058960.D  
 Acq On : 10 Sep 2023 20:55  
 Operator : MA/JU  
 Sample : SSTD16012  
 Misc :  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD160412

Quant Time: Sep 10 23:32:28 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Sep 10 22:58:52 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 09/11/2023  
 Supervised By :mohammad ahmed 09/13/2023

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.329  | 152  | 67433    | 20.000  | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 11.173 | 136  | 313053   | 20.000  | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.951 | 164  | 205195   | 20.000  | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.707 | 188  | 457616   | 20.000  | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 22.037 | 240  | 430367   | 20.000  | ng/u1  | 0.01     |
| 88) Perylene-d12              | 25.609 | 264  | 499948   | 20.000  | ng/u1  | 0.02     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000   | ng/uL  |          |
| 4) Pyridine-d5                | 4.070  | 84   | 797454   | 165.200 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.448  | 99   | 1111325  | 166.161 | ng/u1  | 0.02     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.654  | 67   | 621813   | 164.343 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/u1  |          |
| 15) 4-Methylphenol-d8         | 9.028  | 113  | 842037   | 162.442 | 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        | 11.326 | 131  | 1023763  | 136.184 | ng/u1  | 0.02     |
| 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          | 15.151 | 143  | 396512   | 189.470 | ng/u1  | 0.03     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/u1  |          |
| 65) 4,6-Dinitro-2-methylph... | 16.067 | 200  | 381865   | 183.288 | ng/u1  | 0.03     |
| 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                   | 4.093  | 79   | 858702   | 168.987 | ng/u1  | 93       |
| 6) Benzaldehyde               | 7.477  | 77   | 584730   | 148.320 | ng/u1# | 80       |
| 8) Phenol                     | 7.483  | 94   | 1157977  | 166.024 | ng/u1  | 90       |
| 10) Bis(2-Chloroethyl)ether   | 7.754  | 93   | 816247   | 157.795 | ng/u1  | 91       |
| 13) 2-Methylphenol            | 8.752  | 108  | 872930   | 167.151 | ng/u1  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.852  | 45   | 991528   | 166.548 | ng/u1  | 96       |
| 16) Acetophenone              | 9.187  | 105  | 1373031  | 148.554 | ng/u1# | 97       |
| 18) 4-Methylphenol            | 9.099  | 108  | 948930   | 166.048 | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 11.349 | 127  | 957748   | 134.325 | ng/u1  | 99       |
| 34) Caprolactam               | 12.178 | 113  | 240021m  | 130.615 | ng/u1  |          |
| 40) Hexachlorocyclopentadiene | 13.100 | 237  | 735576   | 133.488 | ng/u1  | 99       |
| 51) 3-Nitroaniline            | 14.892 | 138  | 433759   | 188.315 | ng/u1  | 93       |
| 53) 2,4-Dinitrophenol         | 15.080 | 184  | 291717   | 150.864 | ng/u1  | 89       |
| 55) 4-Nitrophenol             | 15.168 | 109  | 489967   | 187.191 | ng/u1  | 86       |
| 63) 4-Nitroaniline            | 16.067 | 138  | 373878m  | 174.449 | ng/u1  |          |
| 66) 4,6-Dinitro-2-methylph... | 16.085 | 198  | 413312   | 188.965 | ng/u1# | 91       |
| 70) Atrazine                  | 17.143 | 200  | 612387   | 135.536 | ng/u1  | 99       |
| 71) Pentachlorophenol         | 17.325 | 266  | 568918   | 144.462 | ng/u1  | 100      |
| 77) Carbazole                 | 18.118 | 167  | 2817402  | 163.649 | ng/u1  | 94       |
| 84) 3,3'-Dichlorobenzidine    | 21.925 | 252  | 1331173  | 141.609 | ng/u1# | 92       |
| 89) Di-n-octyl phthalate      | 23.194 | 149  | 3486935  | 223.333 | ng/u1  | 100      |

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