

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101624\  
 Data File : BN034611.D  
 Acq On : 16 Oct 2024 15:16  
 Operator : RC/JU  
 Sample : SSTD16091  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD160229

Quant Time: Oct 16 15:48:14 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101624.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 16 14:56:03 2024  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/17/2024  
 Supervised By :mohammad ahmed 10/17/2024

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.669  | 152  | 232473   | 20.000  | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.445 | 136  | 1022000  | 20.000  | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.304 | 164  | 613166   | 20.000  | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.051 | 188  | 1194046  | 20.000  | ng/u1  | 0.01     |
| 79) Chrysene-d12              | 21.280 | 240  | 1015956  | 20.000  | ng/u1  | 0.01     |
| 88) Perylene-d12              | 24.174 | 264  | 1240091  | 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                | 3.599  | 84   | 2613602  | 151.635 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.846  | 99   | 3063932  | 148.752 | ng/u1  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.016  | 67   | 1803090  | 141.262 | ng/u1  | 0.01     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/u1  |          |
| 15) 4-Methylphenol-d8         | 8.381  | 113  | 2407677  | 143.460 | 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        | 10.587 | 131  | 3415857  | 141.896 | 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          | 14.516 | 143  | 1286753  | 141.908 | ng/u1  | 0.03     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/u1  |          |
| 65) 4,6-Dinitro-2-methylph... | 15.433 | 200  | 882768   | 135.409 | 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                   | 3.616  | 79   | 2664145  | 151.209 | ng/u1  | 98       |
| 6) Benzaldehyde               | 6.810  | 77   | 943294   | 86.086  | ng/u1  | 100      |
| 8) Phenol                     | 6.875  | 94   | 3161805  | 146.848 | ng/u1  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 7.110  | 93   | 2416725  | 144.316 | ng/u1  | 97       |
| 13) 2-Methylphenol            | 8.110  | 108  | 2355304  | 146.895 | ng/u1  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.199  | 45   | 3533791  | 143.152 | ng/u1  | 98       |
| 16) Acetophenone              | 8.493  | 105  | 3631917  | 136.105 | ng/u1  | 97       |
| 18) 4-Methylphenol            | 8.457  | 108  | 2549875  | 143.703 | ng/u1  | 97       |
| 32) 4-Chloroaniline           | 10.610 | 127  | 3204805  | 135.871 | ng/u1  | 100      |
| 34) Caprolactam               | 11.410 | 113  | 828851m  | 141.876 | ng/u1  |          |
| 40) Hexachlorocyclopentadiene | 12.475 | 237  | 1314749  | 132.488 | ng/u1  | 96       |
| 51) 3-Nitroaniline            | 14.216 | 138  | 1359576  | 137.491 | ng/u1# | 85       |
| 53) 2,4-Dinitrophenol         | 14.416 | 184  | 627653   | 127.007 | ng/u1  | 93       |
| 55) 4-Nitrophenol             | 14.539 | 109  | 859322   | 113.828 | ng/u1  | 93       |
| 63) 4-Nitroaniline            | 15.386 | 138  | 1098364  | 107.855 | ng/u1  | 95       |
| 66) 4,6-Dinitro-2-methylph... | 15.445 | 198  | 986731   | 135.029 | ng/u1  | 99       |
| 70) Atrazine                  | 16.545 | 200  | 1740387  | 136.173 | ng/u1  | 98       |
| 71) Pentachlorophenol         | 16.704 | 266  | 1175627  | 132.098 | ng/u1  | 97       |
| 77) Carbazole                 | 17.451 | 167  | 8289834  | 133.110 | ng/u1  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.192 | 252  | 2944303  | 129.433 | ng/u1  | 93       |
| 89) Di-n-octyl phthalate      | 22.351 | 149  | 11661551 | 143.998 | ng/u1  | 100      |

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