

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN120224\  
 Data File : BN035389.D  
 Acq On : 02 Dec 2024 16:56  
 Operator : RC/JU  
 Sample : SSTD16097  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD160236

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 12/03/2024  
 Supervised By :mohammad ahmed 12/05/2024

Quant Time: Dec 02 18:24:20 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN120224.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 02 18:09:44 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.305  | 152  | 116089   | 20.000  | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.057 | 136  | 489921   | 20.000  | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 13.969 | 164  | 353904   | 20.000  | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 16.739 | 188  | 764279   | 20.000  | ng/u1  | 0.01     |
| 79) Chrysene-d12              | 20.998 | 240  | 628887   | 20.000  | ng/u1  | 0.02     |
| 88) Perylene-d12              | 23.680 | 264  | 709743   | 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                | 3.340  | 84   | 967078   | 149.333 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.528  | 99   | 1300214  | 157.219 | ng/u1  | 0.02     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.675  | 67   | 677033   | 146.797 | ng/u1  | 0.01     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.000   | ng/u1  |          |
| 15) 4-Methylphenol-d8         | 8.046  | 113  | 1112976  | 154.426 | ng/u1  | 0.03     |
| 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        | 10.222 | 131  | 1466762  | 147.522 | 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          | 14.251 | 143  | 677614   | 161.935 | ng/u1  | 0.05     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.000   | ng/u1  |          |
| 65) 4,6-Dinitro-2-methylph... | 15.145 | 200  | 694364   | 172.269 | ng/u1  | 0.04     |
| 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                   | 3.358  | 79   | 960466   | 147.900 | ng/u1  | 95       |
| 6) Benzaldehyde               | 6.469  | 77   | 439302   | 98.482  | ng/u1  | 98       |
| 8) Phenol                     | 6.558  | 94   | 1296859  | 151.408 | ng/u1  | 94       |
| 10) Bis(2-Chloroethyl)ether   | 6.769  | 93   | 976410   | 145.613 | ng/u1  | 98       |
| 13) 2-Methylphenol            | 7.769  | 108  | 1024635  | 154.070 | ng/u1  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 7.840  | 45   | 773419   | 145.932 | ng/u1  | 99       |
| 16) Acetophenone              | 8.140  | 105  | 1583941  | 139.161 | ng/u1# | 86       |
| 18) 4-Methylphenol            | 8.122  | 108  | 1073381  | 145.254 | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 10.246 | 127  | 1375148  | 144.379 | ng/u1  | 100      |
| 34) Caprolactam               | 11.104 | 113  | 356981m  | 151.667 | ng/u1  |          |
| 40) Hexachlorocyclopentadiene | 12.104 | 237  | 848465   | 161.933 | ng/u1  | 95       |
| 51) 3-Nitroaniline            | 13.916 | 138  | 688897   | 156.984 | ng/u1  | 100      |
| 53) 2,4-Dinitrophenol         | 14.122 | 184  | 565733   | 191.149 | ng/u1  | 95       |
| 55) 4-Nitrophenol             | 14.269 | 109  | 642925   | 162.678 | ng/u1  | 89       |
| 63) 4-Nitroaniline            | 15.116 | 138  | 588098m  | 138.120 | ng/u1  |          |
| 66) 4,6-Dinitro-2-methylph... | 15.163 | 198  | 704999   | 161.431 | ng/u1# | 97       |
| 70) Atrazine                  | 16.263 | 200  | 1022211  | 140.958 | ng/u1  | 97       |
| 71) Pentachlorophenol         | 16.398 | 266  | 866114   | 148.689 | ng/u1  | 96       |
| 77) Carbazole                 | 17.163 | 167  | 4405477  | 133.255 | ng/u1  | 98       |
| 84) 3,3'-Dichlorobenzidine    | 20.927 | 252  | 1580762  | 140.981 | ng/u1  | 99       |
| 89) Di-n-octyl phthalate      | 21.986 | 149  | 6040723  | 187.238 | ng/u1  | 100      |

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