

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN021324\  
 Data File : BN029717.D  
 Acq On : 12 Feb 2024 17:11  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD0.4345

Quant Time: Feb 12 18:17:09 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-SIM-BN020624.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sun Feb 11 23:18:17 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.880  | 152  | 5479     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.672 | 136  | 14765    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.516 | 164  | 6226     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.272 | 188  | 10712    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.473 | 240  | 7741     | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.843 | 264  | 7370     | 0.400 | ng/ul  | # 0.00   |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.332  | 96   | 3178     | 0.465 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.267 | 152  | 8167     | 0.435 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.304 | 212  | 10584    | 0.416 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       |        | Qvalue   |
| 2) 1,4-Dioxane              | 3.365  | 88   | 3264     | 0.439 | ng/ul  | 100      |
| 5) Naphthalene              | 10.722 | 128  | 18366    | 0.442 | ng/ul  | 99       |
| 7) 2-Methylnaphthalene      | 12.338 | 142  | 10284    | 0.432 | ng/ul  | 100      |
| 8) 1-Methylnaphthalene      | 12.558 | 142  | 10457    | 0.431 | ng/ul  | 100      |
| 10) Acenaphthylene          | 14.238 | 152  | 13091    | 0.431 | ng/ul  | 99       |
| 11) Acenaphthene            | 14.580 | 153  | 10118    | 0.438 | ng/ul  | 98       |
| 12) Fluorene                | 15.570 | 166  | 10102    | 0.418 | ng/ul  | 99       |
| 14) Pentachlorophenol       | 16.917 | 266  | 934      | 0.391 | ng/ul# | 50       |
| 15) Phenanthrene            | 17.314 | 178  | 14307    | 0.428 | ng/ul  | 100      |
| 16) Anthracene              | 17.407 | 178  | 12023    | 0.427 | ng/ul  | 99       |
| 19) Fluoranthene            | 19.336 | 202  | 14491    | 0.408 | ng/ul  | 98       |
| 20) Pyrene                  | 19.699 | 202  | 14983    | 0.411 | ng/ul  | 98       |
| 21) Benzo(a)anthracene      | 21.458 | 228  | 11309    | 0.405 | ng/ul  | 100      |
| 22) Chrysene                | 21.508 | 228  | 12936    | 0.424 | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 23.121 | 252  | 12965    | 0.468 | ng/ul  | 99       |
| 25) Benzo(k)fluoranthene    | 23.168 | 252  | 13663    | 0.444 | ng/ul  | 96       |
| 26) Benzo(a)pyrene          | 23.738 | 252  | 11519    | 0.442 | ng/ul  | 98       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.292 | 276  | 14625    | 0.462 | ng/ul# | 100      |
| 28) Dibenzo(a,h)anthracene  | 26.329 | 278  | 11409    | 0.460 | ng/ul  | 98       |
| 29) Benzo(g,h,i)perylene    | 27.043 | 276  | 13551    | 0.472 | ng/ul  | 99       |
| -----                       |        |      |          |       |        |          |

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

