

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM052824\  
 Data File : BM045771.D  
 Acq On : 28 May 2024 21:39  
 Operator : MA/JU  
 Sample : SSTD1.643  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD1.6005

Quant Time: May 28 23:52:11 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM052824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue May 28 23:49:11 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.489  | 152  | 6693     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.259 | 136  | 20022    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.126 | 164  | 10426    | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 16.870 | 188  | 20869    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.061 | 240  | 14079    | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.174 | 264  | 15023    | 0.400 | ng/ul  | 0.00     |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.063  | 96   | 13443    | 1.596 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.848 | 152  | 43836    | 1.588 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 18.898 | 212  | 77127    | 2.158 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.096  | 88   | 13549    | 1.678 | ng/ul# | 84       |
| 5) Naphthalene              | 10.303 | 128  | 82181    | 1.549 | ng/ul  | 100      |
| 7) 2-Methylnaphthalene      | 11.925 | 142  | 51694    | 1.586 | ng/ul  | 98       |
| 8) 1-Methylnaphthalene      | 12.150 | 142  | 54328    | 1.576 | ng/ul  | 100      |
| 10) Acenaphthylene          | 13.844 | 152  | 77104    | 1.554 | ng/ul  | 99       |
| 11) Acenaphthene            | 14.186 | 153  | 55106    | 1.579 | ng/ul  | 100      |
| 12) Fluorene                | 15.181 | 166  | 62159    | 1.566 | ng/ul  | 98       |
| 14) Pentachlorophenol       | 16.566 | 266  | 1778     | 0.388 | ng/ul  | 95       |
| 15) Phenanthrene            | 16.912 | 178  | 93434    | 1.595 | ng/ul  | 100      |
| 16) Anthracene              | 17.001 | 178  | 88841    | 1.518 | ng/ul  | 99       |
| 19) Fluoranthene            | 18.931 | 202  | 103959   | 2.069 | ng/ul  | 99       |
| 20) Pyrene                  | 19.294 | 202  | 106775   | 1.988 | ng/ul  | 98       |
| 21) Benzo(a)anthracene      | 21.044 | 228  | 89401    | 1.604 | ng/ul  | 99       |
| 22) Chrysene                | 21.096 | 228  | 89273    | 1.360 | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 22.546 | 252  | 87917    | 1.428 | ng/ul  | 80       |
| 25) Benzo(k)fluoranthene    | 22.587 | 252  | 87028    | 1.370 | ng/ul# | 83       |
| 26) Benzo(a)pyrene          | 23.081 | 252  | 75082    | 1.503 | ng/ul# | 73       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.246 | 276  | 80189    | 1.304 | ng/ul# | 98       |
| 28) Dibenzo(a,h)anthracene  | 25.263 | 278  | 58837    | 1.349 | ng/ul# | 84       |
| 29) Benzo(g,h,i)perylene    | 25.880 | 276  | 77813    | 1.354 | ng/ul  | 94       |
| -----                       |        |      |          |       |        |          |

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

Quant Time: May 28 23:52:11 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM052824.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Tue May 28 23:49:11 2024  
Response via : Initial Calibration

