

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM121024\  
 Data File : BM048951.D  
 Acq On : 11 Dec 2024 11:41  
 Operator : RC/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.4107

Quant Time: Dec 11 12:10:38 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM120924.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Dec 09 17:52:26 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.582  | 152  | 6570     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.341 | 136  | 19434    | 0.400 | ng/ul  | -0.01    |
| 9) Acenaphthene-d10         | 14.214 | 164  | 10552    | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 16.958 | 188  | 20500    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.155 | 240  | 15300    | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.321 | 264  | 16208    | 0.400 | ng/ul  | #-0.01   |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.160  | 96   | 3091     | 0.403 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.942 | 152  | 10347    | 0.417 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 18.992 | 212  | 19007    | 0.416 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.193  | 88   | 3353     | 0.377 | ng/ul  | 97       |
| 5) Naphthalene              | 10.391 | 128  | 19878    | 0.399 | ng/ul  | 99       |
| 7) 2-Methylnaphthalene      | 12.019 | 142  | 12499    | 0.411 | ng/ul  | 99       |
| 8) 1-Methylnaphthalene      | 12.239 | 142  | 12870    | 0.415 | ng/ul  | 99       |
| 10) Acenaphthylene          | 13.932 | 152  | 17823    | 0.400 | ng/ul  | 99       |
| 11) Acenaphthene            | 14.274 | 153  | 13657    | 0.405 | ng/ul  | 98       |
| 12) Fluorene                | 15.269 | 166  | 14894    | 0.398 | ng/ul  | 98       |
| 14) Pentachlorophenol       | 16.617 | 266  | 1410     | 0.334 | ng/ul  | 95       |
| 15) Phenanthrene            | 17.001 | 178  | 22921    | 0.408 | ng/ul  | 99       |
| 16) Anthracene              | 17.094 | 178  | 20644    | 0.412 | ng/ul  | 99       |
| 19) Fluoranthene            | 19.024 | 202  | 25835    | 0.407 | ng/ul  | 98       |
| 20) Pyrene                  | 19.387 | 202  | 27677    | 0.405 | ng/ul  | 99       |
| 21) Benzo(a)anthracene      | 21.140 | 228  | 22120    | 0.411 | ng/ul  | 99       |
| 22) Chrysene                | 21.193 | 228  | 24395    | 0.400 | ng/ul  | 100      |
| 24) Benzo(b)fluoranthene    | 22.674 | 252  | 22538    | 0.402 | ng/ul  | 98       |
| 25) Benzo(k)fluoranthene    | 22.718 | 252  | 25113    | 0.388 | ng/ul  | 99       |
| 26) Benzo(a)pyrene          | 23.224 | 252  | 20795    | 0.401 | ng/ul  | 96       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.471 | 276  | 28353    | 0.390 | ng/ul# | 99       |
| 28) Dibenzo(a,h)anthracene  | 25.488 | 278  | 21630    | 0.382 | ng/ul  | 97       |
| 29) Benzo(g,h,i)perylene    | 26.121 | 276  | 22036    | 0.376 | ng/ul  | 99       |
| -----                       |        |      |          |       |        |          |

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

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