

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM062024\  
 Data File : BM046140.D  
 Acq On : 21 Jun 2024 02:05  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.4136

Quant Time: Jun 21 03:58:05 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM061824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 18 13:58:29 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.459  | 152  | 2720     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.242 | 136  | 8397     | 0.400 | ng/ul  | -0.01    |
| 9) Acenaphthene-d10         | 14.094 | 164  | 4768     | 0.400 | ng/ul  | -0.02    |
| 13) Phenanthrene-d10        | 16.849 | 188  | 9308     | 0.400 | ng/ul  | -0.01    |
| 17) Chrysene-d12            | 21.029 | 240  | 7508     | 0.400 | ng/ul  | -0.01    |
| 23) Perylene-d12            | 23.119 | 264  | 10340    | 0.400 | ng/ul  | #-0.02   |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.042  | 96   | 1529     | 0.410 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.854 | 152  | 4679     | 0.415 | ng/ul  | -0.01    |
| 18) Fluoranthene-d10        | 18.875 | 212  | 9363     | 0.416 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       |        | Qvalue   |
| 2) 1,4-Dioxane              | 3.067  | 88   | 1637     | 0.391 | ng/ul  | 95       |
| 5) Naphthalene              | 10.292 | 128  | 9644     | 0.412 | ng/ul  | 99       |
| 7) 2-Methylnaphthalene      | 11.925 | 142  | 5499     | 0.417 | ng/ul  | 99       |
| 8) 1-Methylnaphthalene      | 12.129 | 142  | 6158     | 0.426 | ng/ul  | 100      |
| 10) Acenaphthylene          | 13.821 | 152  | 9580     | 0.424 | ng/ul  | 99       |
| 11) Acenaphthene            | 14.159 | 153  | 6894     | 0.422 | ng/ul  | 99       |
| 12) Fluorene                | 15.167 | 166  | 7368     | 0.421 | ng/ul  | 98       |
| 14) Pentachlorophenol       | 16.536 | 266  | 1150     | 0.412 | ng/ul  | 95       |
| 15) Phenanthrene            | 16.887 | 178  | 10821    | 0.414 | ng/ul  | 99       |
| 16) Anthracene              | 16.992 | 178  | 8378     | 0.430 | ng/ul  | 98       |
| 19) Fluoranthene            | 18.903 | 202  | 14123    | 0.434 | ng/ul  | 100      |
| 20) Pyrene                  | 19.261 | 202  | 15036    | 0.428 | ng/ul  | 100      |
| 21) Benzo(a)anthracene      | 21.015 | 228  | 12110    | 0.418 | ng/ul  | 99       |
| 22) Chrysene                | 21.067 | 228  | 14916    | 0.403 | ng/ul  | 98       |
| 24) Benzo(b)fluoranthene    | 22.502 | 252  | 15626    | 0.419 | ng/ul  | 94       |
| 25) Benzo(k)fluoranthene    | 22.543 | 252  | 17153    | 0.387 | ng/ul  | 92       |
| 26) Benzo(a)pyrene          | 23.031 | 252  | 14561    | 0.422 | ng/ul# | 88       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.172 | 276  | 21088    | 0.373 | ng/ul# | 100      |
| 28) Dibenzo(a,h)anthracene  | 25.182 | 278  | 16223    | 0.386 | ng/ul  | 93       |
| 29) Benzo(g,h,i)perylene    | 25.799 | 276  | 17944    | 0.370 | ng/ul  | 97       |
| -----                       |        |      |          |       |        |          |

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

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