

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM100722\  
 Data File : BM036898.D  
 Acq On : 08 Oct 2022 16:18  
 Operator : CG/JU  
 Sample : SSTDCCC0.4EC  
 Misc :  
 ALS Vial : 53 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.4136

Quant Time: Oct 10 01:26:25 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM100622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Oct 10 00:54:15 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units   | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| -----                       |        |      |          |       |         |          |
| Internal Standards          |        |      |          |       |         |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.959  | 152  | 7721     | 0.400 | ng/ul   | 0.00     |
| 4) Naphthalene-d8           | 10.765 | 136  | 26845    | 0.400 | ng/ul # | 0.00     |
| 9) Acenaphthene-d10         | 14.584 | 164  | 15271    | 0.400 | ng/ul   | 0.00     |
| 13) Phenanthrene-d10        | 17.322 | 188  | 31519    | 0.400 | ng/ul   | 0.00     |
| 17) Chrysene-d12            | 21.488 | 240  | 23341    | 0.400 | ng/ul   | 0.00     |
| 23) Perylene-d12            | 23.888 | 264  | 16110    | 0.400 | ng/ul # | 0.00     |
| System Monitoring Compounds |        |      |          |       |         |          |
| 3) 1,4-Dioxane-d8           | 3.352  | 96   | 3761     | 0.345 | ng/ul   | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.349 | 152  | 16472    | 0.406 | ng/ul   | 0.00     |
| 18) Fluoranthene-d10        | 19.340 | 212  | 32074    | 0.459 | ng/ul   | 0.00     |
| Target Compounds            |        |      |          |       |         |          |
|                             |        |      |          |       |         | Qvalue   |
| 2) 1,4-Dioxane              | 3.390  | 88   | 3670     | 0.345 | ng/ul#  | 85       |
| 5) Naphthalene              | 10.814 | 128  | 30830    | 0.402 | ng/ul   | 96       |
| 7) 2-Methylnaphthalene      | 12.420 | 142  | 19475    | 0.415 | ng/ul   | 91       |
| 8) 1-Methylnaphthalene      | 12.640 | 142  | 19860    | 0.417 | ng/ul   | 99       |
| 10) Acenaphthylene          | 14.307 | 152  | 26226    | 0.413 | ng/ul   | 98       |
| 11) Acenaphthene            | 14.644 | 153  | 22182    | 0.403 | ng/ul   | 99       |
| 12) Fluorene                | 15.630 | 166  | 24775    | 0.397 | ng/ul   | 99       |
| 14) Pentachlorophenol       | 16.971 | 266  | 2748     | 0.360 | ng/ul   | 97       |
| 15) Phenanthrene            | 17.364 | 178  | 40117    | 0.399 | ng/ul   | 99       |
| 16) Anthracene              | 17.453 | 178  | 34370    | 0.409 | ng/ul   | 99       |
| 19) Fluoranthene            | 19.373 | 202  | 42986    | 0.463 | ng/ul   | 97       |
| 20) Pyrene                  | 19.731 | 202  | 43961    | 0.454 | ng/ul   | 99       |
| 21) Benzo(a)anthracene      | 21.474 | 228  | 34413    | 0.408 | ng/ul   | 98       |
| 22) Chrysene                | 21.526 | 228  | 36109    | 0.399 | ng/ul   | 98       |
| 24) Benzo(b)fluoranthene    | 23.154 | 252  | 30046    | 0.417 | ng/ul   | 94       |
| 25) Benzo(k)fluoranthene    | 23.204 | 252  | 29302    | 0.418 | ng/ul   | 92       |
| 26) Benzo(a)pyrene          | 23.780 | 252  | 24215    | 0.415 | ng/ul#  | 90       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.376 | 276  | 29748    | 0.360 | ng/ul#  | 45       |
| 28) Dibenzo(a,h)anthracene  | 26.400 | 278  | 23637    | 0.354 | ng/ul#  | 90       |
| 29) Benzo(g,h,i)perylene    | 27.144 | 276  | 25295    | 0.347 | ng/ul   | 97       |
| -----                       |        |      |          |       |         |          |

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

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