

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM072023\  
 Data File : BM040970.D  
 Acq On : 20 Jul 2023 07:58  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.4186

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 07/20/2023  
 Supervised By :Sohil Jodhani 07/20/2023

Quant Time: Jul 20 08:46:35 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM071923.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jul 20 05:08:42 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.878  | 152  | 2720     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.680 | 136  | 12841    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.523 | 164  | 5114     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.282 | 188  | 9414     | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.472 | 240  | 6137     | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.868 | 264  | 5210     | 0.400 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.284  | 96   | 1969     | 0.440 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.280 | 152  | 5240     | 0.306 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.311 | 212  | 9210     | 0.357 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.318  | 88   | 1987     | 0.422 | ng/ul  | 90       |
| 5) Naphthalene              | 10.730 | 128  | 13432    | 0.382 | ng/ul  | 99       |
| 7) 2-Methylnaphthalene      | 12.352 | 142  | 5905     | 0.306 | ng/ul  | 99       |
| 8) 1-Methylnaphthalene      | 12.566 | 142  | 6360     | 0.311 | ng/ul  | 99       |
| 10) Acenaphthylene          | 14.250 | 152  | 8206     | 0.383 | ng/ul  | 100      |
| 11) Acenaphthene            | 14.587 | 153  | 7504     | 0.397 | ng/ul  | 99       |
| 12) Fluorene                | 15.582 | 166  | 7776     | 0.398 | ng/ul  | 99       |
| 14) Pentachlorophenol       | 16.970 | 266  | 669      | 0.353 | ng/ul  | 99       |
| 15) Phenanthrene            | 17.324 | 178  | 11289    | 0.401 | ng/ul  | 100      |
| 16) Anthracene              | 17.422 | 178  | 8351     | 0.381 | ng/ul  | 99       |
| 19) Fluoranthene            | 19.343 | 202  | 11795    | 0.372 | ng/ul  | 100      |
| 20) Pyrene                  | 19.706 | 202  | 12685    | 0.392 | ng/ul  | 100      |
| 21) Benzo(a)anthracene      | 21.457 | 228  | 7195     | 0.385 | ng/ul  | 99       |
| 22) Chrysene                | 21.510 | 228  | 8541     | 0.400 | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 23.138 | 252  | 8218m    | 0.411 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 23.187 | 252  | 7716     | 0.406 | ng/ul  | 99       |
| 26) Benzo(a)pyrene          | 23.766 | 252  | 6742     | 0.390 | ng/ul  | 99       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.347 | 276  | 9260     | 0.404 | ng/ul# | 99       |
| 28) Dibenzo(a,h)anthracene  | 26.367 | 278  | 6988     | 0.403 | ng/ul  | 99       |
| 29) Benzo(g,h,i)perylene    | 27.105 | 276  | 8851     | 0.418 | ng/ul  | 99       |
| -----                       |        |      |          |       |        |          |

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

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