

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM091323\  
 Data File : BM041827.D  
 Acq On : 13 Sep 2023 12:22  
 Operator : MA/JU  
 Sample : SSTD0.432  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.4024

Quant Time: Sep 13 16:09:47 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM091323.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 13 14:45:01 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.975  | 152  | 4245     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.790 | 136  | 10464    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.615 | 164  | 5115     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.367 | 188  | 11111    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.539 | 240  | 7177     | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.956 | 264  | 8277     | 0.400 | ng/ul  | 0.00     |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.381  | 96   | 2287     | 0.424 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.368 | 152  | 5863     | 0.412 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.390 | 212  | 10333    | 0.420 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.419  | 88   | 2395     | 0.423 | ng/ul  | 100      |
| 5) Naphthalene              | 10.840 | 128  | 14654    | 0.416 | ng/ul  | 100      |
| 7) 2-Methylnaphthalene      | 12.445 | 142  | 8738     | 0.412 | ng/ul  | 100      |
| 8) 1-Methylnaphthalene      | 12.665 | 142  | 8912     | 0.412 | ng/ul  | 100      |
| 10) Acenaphthylene          | 14.342 | 152  | 13251    | 0.413 | ng/ul  | 100      |
| 11) Acenaphthene            | 14.680 | 153  | 10195    | 0.415 | ng/ul  | 100      |
| 12) Fluorene                | 15.665 | 166  | 11759    | 0.410 | ng/ul  | 100      |
| 14) Pentachlorophenol       | 16.991 | 266  | 2565     | 0.405 | ng/ul  | 100      |
| 15) Phenanthrene            | 17.409 | 178  | 19701    | 0.406 | ng/ul  | 100      |
| 16) Anthracene              | 17.502 | 178  | 17476    | 0.413 | ng/ul  | 100      |
| 19) Fluoranthene            | 19.417 | 202  | 22900    | 0.412 | ng/ul  | 100      |
| 20) Pyrene                  | 19.785 | 202  | 23988    | 0.412 | ng/ul  | 100      |
| 21) Benzo(a)anthracene      | 21.521 | 228  | 19321    | 0.408 | ng/ul  | 100      |
| 22) Chrysene                | 21.577 | 228  | 19811    | 0.409 | ng/ul  | 100      |
| 24) Benzo(b)fluoranthene    | 23.208 | 252  | 23322    | 0.396 | ng/ul  | 100      |
| 25) Benzo(k)fluoranthene    | 23.260 | 252  | 21898    | 0.403 | ng/ul  | 100      |
| 26) Benzo(a)pyrene          | 23.845 | 252  | 18202    | 0.405 | ng/ul  | 100      |
| 27) Indeno(1,2,3-cd)pyrene  | 26.455 | 276  | 27574    | 0.399 | ng/ul  | 100      |
| 28) Dibenzo(a,h)anthracene  | 26.475 | 278  | 20383    | 0.400 | ng/ul  | 100      |
| 29) Benzo(g,h,i)perylene    | 27.236 | 276  | 23495    | 0.402 | ng/ul  | 100      |
| -----                       |        |      |          |       |        |          |

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

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