

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092424\  
 Data File : BM047615.D  
 Acq On : 24 Sep 2024 17:54  
 Operator : RC/JU  
 Sample : SSTD0.489  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.4010

Quant Time: Sep 25 02:50:37 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM092424.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 25 02:35:09 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.830  | 152  | 7574     | 0.400 | ng/u1  | 0.00     |
| 4) Naphthalene-d8           | 10.623 | 136  | 22985    | 0.400 | ng/u1  | 0.00     |
| 9) Acenaphthene-d10         | 14.460 | 164  | 10768    | 0.400 | ng/u1  | 0.00     |
| 13) Phenanthrene-d10        | 17.196 | 188  | 22026    | 0.400 | ng/u1  | 0.00     |
| 17) Chrysene-d12            | 21.362 | 240  | 18126    | 0.400 | ng/u1  | 0.00     |
| 23) Perylene-d12            | 23.642 | 264  | 19359    | 0.400 | ng/u1  | 0.00     |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.273  | 96   | 4111     | 0.538 | ng/u1  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.212 | 152  | 12468    | 0.363 | ng/u1  | 0.00     |
| 18) Fluoranthene-d10        | 19.216 | 212  | 20829    | 0.398 | ng/u1  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.307  | 88   | 4907     | 0.575 | ng/u1  | 100      |
| 5) Naphthalene              | 10.672 | 128  | 24868    | 0.414 | ng/u1  | 100      |
| 7) 2-Methylnaphthalene      | 12.289 | 142  | 14920    | 0.381 | ng/u1  | 100      |
| 8) 1-Methylnaphthalene      | 12.509 | 142  | 15587    | 0.380 | ng/u1  | 100      |
| 10) Acenaphthylene          | 14.183 | 152  | 21725    | 0.435 | ng/u1  | 100      |
| 11) Acenaphthene            | 14.525 | 153  | 15106    | 0.444 | ng/u1  | 100      |
| 12) Fluorene                | 15.506 | 166  | 16585    | 0.419 | ng/u1  | 100      |
| 14) Pentachlorophenol       | 16.850 | 266  | 2414     | 0.344 | ng/u1  | 100      |
| 15) Phenanthrene            | 17.238 | 178  | 25620    | 0.404 | ng/u1  | 100      |
| 16) Anthracene              | 17.331 | 178  | 22753    | 0.402 | ng/u1  | 100      |
| 19) Fluoranthene            | 19.248 | 202  | 28007    | 0.381 | ng/u1  | 100      |
| 20) Pyrene                  | 19.606 | 202  | 29827    | 0.388 | ng/u1  | 100      |
| 21) Benzo(a)anthracene      | 21.344 | 228  | 26699    | 0.401 | ng/u1  | 100      |
| 22) Chrysene                | 21.397 | 228  | 28140    | 0.398 | ng/u1  | 100      |
| 24) Benzo(b)fluoranthene    | 22.955 | 252  | 31373    | 0.408 | ng/u1  | 100      |
| 25) Benzo(k)fluoranthene    | 22.998 | 252  | 31750    | 0.395 | ng/u1  | 100      |
| 26) Benzo(a)pyrene          | 23.542 | 252  | 24593    | 0.405 | ng/u1  | 100      |
| 27) Indeno(1,2,3-cd)pyrene  | 25.970 | 276  | 39027    | 0.423 | ng/u1  | 100      |
| 28) Dibenzo(a,h)anthracene  | 25.980 | 278  | 29985    | 0.418 | ng/u1  | 100      |
| 29) Benzo(g,h,i)perylene    | 26.674 | 276  | 31601    | 0.413 | ng/u1  | 100      |
| -----                       |        |      |          |       |        |          |

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

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