

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM021524\  
 Data File : BM044141.D  
 Acq On : 16 Feb 2024 16:22  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 51 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.4188

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 02/17/2024  
 Supervised By :mohammad ahmed 02/19/2024

Quant Time: Feb 16 23:41:57 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM021524.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 15 13:40:24 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.900  | 152  | 12701    | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.699 | 136  | 42038    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.543 | 164  | 23461    | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.297 | 188  | 49123    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.499 | 240  | 29651    | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.890 | 264  | 27959    | 0.400 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.356  | 96   | 7341     | 0.408 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.294 | 152  | 26124    | 0.444 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.327 | 212  | 49695    | 0.562 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.390  | 88   | 7195     | 0.400 | ng/ul# | 88       |
| 5) Naphthalene              | 10.748 | 128  | 53164    | 0.426 | ng/ul  | 98       |
| 7) 2-Methylnaphthalene      | 12.365 | 142  | 34401    | 0.446 | ng/ul  | 100      |
| 8) 1-Methylnaphthalene      | 12.585 | 142  | 34730    | 0.445 | ng/ul  | 100      |
| 10) Acenaphthylene          | 14.261 | 152  | 56998    | 0.442 | ng/ul  | 99       |
| 11) Acenaphthene            | 14.604 | 153  | 39353    | 0.421 | ng/ul  | 99       |
| 12) Fluorene                | 15.594 | 166  | 43852    | 0.416 | ng/ul  | 99       |
| 14) Pentachlorophenol       | 16.942 | 266  | 6430     | 0.366 | ng/ul  | 98       |
| 15) Phenanthrene            | 17.339 | 178  | 67362    | 0.433 | ng/ul  | 99       |
| 16) Anthracene              | 17.432 | 178  | 67122    | 0.436 | ng/ul  | 99       |
| 19) Fluoranthene            | 19.360 | 202  | 74475    | 0.551 | ng/ul  | 98       |
| 20) Pyrene                  | 19.722 | 202  | 77041    | 0.542 | ng/ul  | 96       |
| 21) Benzo(a)anthracene      | 21.481 | 228  | 60452    | 0.445 | ng/ul  | 100      |
| 22) Chrysene                | 21.537 | 228  | 59190    | 0.404 | ng/ul  | 100      |
| 24) Benzo(b)fluoranthene    | 23.162 | 252  | 50723    | 0.440 | ng/ul  | 96       |
| 25) Benzo(k)fluoranthene    | 23.212 | 252  | 51659m   | 0.412 | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.785 | 252  | 45612    | 0.425 | ng/ul# | 94       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.369 | 276  | 48804    | 0.357 | ng/ul# | 93       |
| 28) Dibenzo(a,h)anthracene  | 26.402 | 278  | 38480    | 0.358 | ng/ul  | 95       |
| 29) Benzo(g,h,i)perylene    | 27.127 | 276  | 40512    | 0.335 | ng/ul  | 97       |
| -----                       |        |      |          |       |        |          |

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

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