

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM062424\  
 Data File : BM046305.D  
 Acq On : 25 Jun 2024 23:24  
 Operator : MA/JU  
 Sample : P2844-15  
 Misc :  
 ALS Vial : 65 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 A4C69

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 06/26/2024  
 Supervised By :mohammad ahmed 06/27/2024

Quant Time: Jun 26 00:05:28 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM061824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 25 08:04:59 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|--------|--------|----------|
| -----                       |        |      |          |        |        |          |
| Internal Standards          |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.392  | 152  | 2297     | 0.400  | ng/ul  | -0.06    |
| 4) Naphthalene-d8           | 10.161 | 136  | 6977     | 0.400  | ng/ul  | -0.09    |
| 9) Acenaphthene-d10         | 14.040 | 164  | 3979     | 0.400  | ng/ul  | -0.05    |
| 13) Phenanthrene-d10        | 16.782 | 188  | 9006m    | 0.400  | ng/ul  | -0.07    |
| 17) Chrysene-d12            | 20.991 | 240  | 6016m    | 0.400  | ng/ul  | -0.04    |
| 23) Perylene-d12            | 23.083 | 264  | 9528m    | 0.400  | ng/ul  | -0.03    |
|                             |        |      |          |        |        |          |
| System Monitoring Compounds |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8           | 3.016  | 96   | 9873     | 3.133  | ng/ul  | -0.01    |
| 6) 2-Methylnaphthalene-d10  | 11.761 | 152  | 2669     | 0.285  | ng/ul  | -0.11    |
| 18) Fluoranthene-d10        | 18.821 | 212  | 6409m    | 0.356  | ng/ul  | -0.05    |
|                             |        |      |          |        |        |          |
| Target Compounds            |        |      |          |        |        |          |
|                             |        |      |          |        | Qvalue |          |
| 5) Naphthalene              | 10.205 | 128  | 21672    | 1.114  | ng/ul  | 94       |
| 7) 2-Methylnaphthalene      | 11.838 | 142  | 8973     | 0.819  | ng/ul  | 96       |
| 8) 1-Methylnaphthalene      | 12.053 | 142  | 6838     | 0.569  | ng/ul  | 99       |
| 10) Acenaphthylene          | 13.753 | 152  | 58561    | 3.103  | ng/ul  | 97       |
| 11) Acenaphthene            | 14.100 | 153  | 24273    | 1.782  | ng/ul  | 98       |
| 12) Fluorene                | 15.094 | 166  | 31235    | 2.136  | ng/ul  | 100      |
| 14) Pentachlorophenol       | 16.470 | 266  | 420      | 0.156  | ng/ul  | 95       |
| 15) Phenanthrene            | 16.825 | 178  | 724040m  | 28.615 | ng/ul  |          |
| 16) Anthracene              | 16.918 | 178  | 138303m  | 7.339  | ng/ul  |          |
| 19) Fluoranthene            | 18.853 | 202  | 1552040m | 59.583 | ng/ul  |          |
| 20) Pyrene                  | 19.216 | 202  | 1127854m | 40.037 | ng/ul  |          |
| 21) Benzo(a)anthracene      | 20.973 | 228  | 651347m  | 28.030 | ng/ul  |          |
| 22) Chrysene                | 21.026 | 228  | 743191m  | 25.061 | ng/ul  |          |
| 24) Benzo(b)fluoranthene    | 22.469 | 252  | 1083433m | 31.562 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 22.504 | 252  | 377536m  | 9.251  | ng/ul  |          |
| 26) Benzo(a)pyrene          | 22.993 | 252  | 427766m  | 13.448 | ng/ul  |          |
| 27) Indeno(1,2,3-cd)pyrene  | 25.125 | 276  | 416967m  | 8.009  | ng/ul  |          |
| 28) Dibenzo(a,h)anthracene  | 25.122 | 278  | 134864m  | 3.480  | ng/ul  |          |
| 29) Benzo(g,h,i)perylene    | 25.742 | 276  | 67240m   | 1.506  | ng/ul  |          |
| -----                       |        |      |          |        |        |          |

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

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