

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM072023\  
 Data File : BM040999.D  
 Acq On : 21 Jul 2023 02:39  
 Operator : MA/JU  
 Sample : 03444-11  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 A46R8

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 07/21/2023  
 Supervised By :mohammad ahmed 07/21/2023

Quant Time: Jul 21 04:54:00 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   | 7.861  | 152  | 4192     | 0.400  | ng/ul  | -0.02    |
| 4) Naphthalene-d8           | 10.669 | 136  | 17012    | 0.400  | ng/ul  | #-0.01   |
| 9) Acenaphthene-d10         | 14.518 | 164  | 5515     | 0.400  | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.270 | 188  | 11674    | 0.400  | ng/ul  | -0.01    |
| 17) Chrysene-d12            | 21.463 | 240  | 6862     | 0.400  | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.851 | 264  | 6711     | 0.400  | ng/ul  | #-0.02   |
|                             |        |      |          |        |        |          |
| System Monitoring Compounds |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8           | 3.280  | 96   | 6624     | 0.961  | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.264 | 152  | 5826     | 0.257  | ng/ul  | -0.02    |
| 18) Fluoranthene-d10        | 19.301 | 212  | 6178     | 0.214  | ng/ul  | -0.01    |
|                             |        |      |          |        |        |          |
| Target Compounds            |        |      |          |        |        |          |
|                             |        |      |          |        | Qvalue |          |
| 5) Naphthalene              | 10.719 | 128  | 32716    | 0.703  | ng/ul  | 98       |
| 7) 2-Methylnaphthalene      | 12.335 | 142  | 4166     | 0.163  | ng/ul  | 98       |
| 8) 1-Methylnaphthalene      | 12.555 | 142  | 3737     | 0.138  | ng/ul  | 99       |
| 10) Acenaphthylene          | 14.236 | 152  | 9306     | 0.403  | ng/ul  | 96       |
| 11) Acenaphthene            | 14.578 | 153  | 11337    | 0.557  | ng/ul  | 99       |
| 12) Fluorene                | 15.568 | 166  | 12920    | 0.613  | ng/ul  | 99       |
| 15) Phenanthrene            | 17.312 | 178  | 243723   | 6.985  | ng/ul  | 98       |
| 16) Anthracene              | 17.405 | 178  | 38651    | 1.422  | ng/ul  | 98       |
| 19) Fluoranthene            | 19.329 | 202  | 400913   | 11.316 | ng/ul  | 98       |
| 20) Pyrene                  | 19.696 | 202  | 357431   | 9.875  | ng/ul  | 97       |
| 21) Benzo(a)anthracene      | 21.445 | 228  | 133059   | 6.365  | ng/ul  | 99       |
| 22) Chrysene                | 21.498 | 228  | 133903   | 5.603  | ng/ul  | 98       |
| 24) Benzo(b)fluoranthene    | 23.123 | 252  | 158873   | 6.168  | ng/ul  | 86       |
| 25) Benzo(k)fluoranthene    | 23.167 | 252  | 62457m   | 2.552  | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.743 | 252  | 99371    | 4.466  | ng/ul# | 75       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.317 | 276  | 79597    | 2.696  | ng/ul# | 91       |
| 28) Dibenzo(a,h)anthracene  | 26.327 | 278  | 18281    | 0.819  | ng/ul  | 95       |
| 29) Benzo(g,h,i)perylene    | 27.078 | 276  | 80261    | 2.942  | ng/ul  | 95       |
| -----                       |        |      |          |        |        |          |

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

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