

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM051622\  
 Data File : BM035100.D  
 Acq On : 16 May 2022 12:40  
 Operator : CG/JU  
 Sample : N2707-06MSD  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 EXL87MSD

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 05/17/2022  
 Supervised By :mohammad ahmed 05/18/2022

Quant Time: May 17 00:50:57 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM051222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu May 12 18:15:39 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units   | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| -----                       |        |      |          |       |         |          |
| Internal Standards          |        |      |          |       |         |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.917  | 152  | 3814     | 0.400 | ng/ul   | 0.00     |
| 4) Naphthalene-d8           | 10.715 | 136  | 11909    | 0.400 | ng/ul # | 0.00     |
| 9) Acenaphthene-d10         | 14.552 | 164  | 7601     | 0.400 | ng/ul   | 0.00     |
| 13) Phenanthrene-d10        | 17.292 | 188  | 16281    | 0.400 | ng/ul # | 0.00     |
| 17) Chrysene-d12            | 21.473 | 240  | 12137    | 0.400 | ng/ul   | 0.00     |
| 23) Perylene-d12            | 23.859 | 264  | 9974     | 0.400 | ng/ul # | 0.00     |
|                             |        |      |          |       |         |          |
| System Monitoring Compounds |        |      |          |       |         |          |
| 3) 1,4-Dioxane-d8           | 3.361  | 96   | 1034     | 0.226 | ng/ul   | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.305 | 152  | 7424     | 0.391 | ng/ul   | 0.00     |
| 18) Fluoranthene-d10        | 19.322 | 212  | 19954    | 0.479 | ng/ul   | 0.00     |
|                             |        |      |          |       |         |          |
| Target Compounds            |        |      |          |       | Qvalue  |          |
| 2) 1,4-Dioxane              | 3.394  | 88   | 2872m    | 0.626 | ng/ul   |          |
| 5) Naphthalene              | 10.765 | 128  | 13473    | 0.343 | ng/ul#  | 89       |
| 7) 2-Methylnaphthalene      | 12.376 | 142  | 9194     | 0.356 | ng/ul   | 99       |
| 8) 1-Methylnaphthalene      | 12.596 | 142  | 9105     | 0.379 | ng/ul   | 100      |
| 10) Acenaphthylene          | 14.270 | 152  | 15198    | 0.338 | ng/ul#  | 90       |
| 11) Acenaphthene            | 14.612 | 153  | 11031    | 0.340 | ng/ul   | 96       |
| 12) Fluorene                | 15.597 | 166  | 13247    | 0.349 | ng/ul#  | 97       |
| 14) Pentachlorophenol       | 16.946 | 266  | 4750     | 0.772 | ng/ul   | 98       |
| 15) Phenanthrene            | 17.334 | 178  | 23498    | 0.379 | ng/ul   | 94       |
| 16) Anthracene              | 17.427 | 178  | 21524    | 0.381 | ng/ul#  | 91       |
| 19) Fluoranthene            | 19.350 | 202  | 28097    | 0.412 | ng/ul   | 98       |
| 20) Pyrene                  | 19.712 | 202  | 28831    | 0.416 | ng/ul   | 100      |
| 21) Benzo(a)anthracene      | 21.459 | 228  | 22728    | 0.405 | ng/ul   | 97       |
| 22) Chrysene                | 21.511 | 228  | 22981    | 0.393 | ng/ul   | 98       |
| 24) Benzo(b)fluoranthene    | 23.136 | 252  | 20655    | 0.395 | ng/ul   | 90       |
| 25) Benzo(k)fluoranthene    | 23.183 | 252  | 19660    | 0.394 | ng/ul#  | 90       |
| 26) Benzo(a)pyrene          | 23.753 | 252  | 18665    | 0.392 | ng/ul#  | 84       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.336 | 276  | 20439    | 0.434 | ng/ul#  | 82       |
| 28) Dibenzo(a,h)anthracene  | 26.350 | 278  | 16533    | 0.441 | ng/ul#  | 89       |
| 29) Benzo(g,h,i)perylene    | 27.091 | 276  | 17406    | 0.430 | ng/ul   | 94       |
| -----                       |        |      |          |       |         |          |

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

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