

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM042522\  
 Data File : BM034813.D  
 Acq On : 25 Apr 2022 23:37  
 Operator : CG/JU  
 Sample : N2373-10  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BFFE7

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 04/26/2022  
 Supervised By :Yogesh Patel 04/29/2022

Quant Time: Apr 26 03:07:44 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM042222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Apr 22 16:54:34 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.943  | 152  | 6196     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.743 | 136  | 15911    | 0.400 | ng/ul  | # 0.00   |
| 9) Acenaphthene-d10         | 14.571 | 164  | 8330     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.309 | 188  | 15248    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.486 | 240  | 12051    | 0.400 | ng/ul  | # 0.00   |
| 23) Perylene-d12            | 23.879 | 264  | 12444    | 0.400 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.382  | 96   | 25668    | 3.746 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.327 | 152  | 5488     | 0.245 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.331 | 212  | 9778     | 0.257 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 5) Naphthalene              | 10.793 | 128  | 9116     | 0.200 | ng/ul# | 89       |
| 7) 2-Methylnaphthalene      | 12.404 | 142  | 5422     | 0.179 | ng/ul  | 99       |
| 8) 1-Methylnaphthalene      | 12.619 | 142  | 5563     | 0.187 | ng/ul  | 99       |
| 10) Acenaphthylene          | 14.289 | 152  | 26059    | 0.716 | ng/ul# | 91       |
| 11) Acenaphthene            | 14.631 | 153  | 3567     | 0.116 | ng/ul  | 97       |
| 12) Fluorene                | 15.616 | 166  | 8438     | 0.243 | ng/ul  | 99       |
| 15) Phenanthrene            | 17.352 | 178  | 187422   | 3.718 | ng/ul  | 93       |
| 16) Anthracene              | 17.444 | 178  | 24641    | 0.549 | ng/ul# | 92       |
| 19) Fluoranthene            | 19.364 | 202  | 290458   | 5.021 | ng/ul  | 97       |
| 20) Pyrene                  | 19.726 | 202  | 328909   | 5.661 | ng/ul  | 100      |
| 21) Benzo(a)anthracene      | 21.468 | 228  | 115692   | 2.468 | ng/ul  | 98       |
| 22) Chrysene                | 21.521 | 228  | 168635   | 3.374 | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 23.151 | 252  | 198371   | 3.491 | ng/ul  | 88       |
| 25) Benzo(k)fluoranthene    | 23.192 | 252  | 66387m   | 1.169 | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.771 | 252  | 135128   | 2.914 | ng/ul# | 78       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.353 | 276  | 104035   | 1.598 | ng/ul# | 90       |
| 28) Dibenzo(a,h)anthracene  | 26.367 | 278  | 25722    | 0.486 | ng/ul  | 93       |
| 29) Benzo(g,h,i)perylene    | 27.115 | 276  | 96041    | 1.709 | ng/ul  | 93       |
| -----                       |        |      |          |       |        |          |

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

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