

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN070823\  
 Data File : BN026401.D  
 Acq On : 09 Jul 2023 00:46  
 Operator : MA/JU  
 Sample : 03351-03MSD  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 YBW10MSD

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 07/10/2023  
 Supervised By :mohammad ahmed 07/10/2023

Quant Time: Jul 09 02:44:41 2023  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-SIM-BN062223.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Jul 08 09:20:18 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.925  | 152  | 3900     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.730 | 136  | 16444    | 0.400 | ng/ul  | #-0.01   |
| 9) Acenaphthene-d10         | 14.570 | 164  | 9587     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.321 | 188  | 17824    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.519 | 240  | 10174    | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.969 | 264  | 11362    | 0.400 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.301  | 96   | 22911    | 5.216 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.331 | 152  | 6641     | 0.263 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.353 | 212  | 11585    | 0.299 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.335  | 88   | 6119     | 1.312 | ng/ul# | 89       |
| 5) Naphthalene              | 10.780 | 128  | 13322    | 0.283 | ng/ul  | 98       |
| 7) 2-Methylnaphthalene      | 12.402 | 142  | 8048     | 0.268 | ng/ul  | 99       |
| 8) 1-Methylnaphthalene      | 12.617 | 142  | 8460     | 0.272 | ng/ul  | 99       |
| 10) Acenaphthylene          | 14.288 | 152  | 14930    | 0.333 | ng/ul  | 100      |
| 11) Acenaphthene            | 14.630 | 153  | 10057    | 0.274 | ng/ul  | 100      |
| 12) Fluorene                | 15.615 | 166  | 11525    | 0.274 | ng/ul  | 99       |
| 14) Pentachlorophenol       | 16.983 | 266  | 1573     | 0.246 | ng/ul  | 98       |
| 15) Phenanthrene            | 17.359 | 178  | 16759    | 0.285 | ng/ul  | 99       |
| 16) Anthracene              | 17.452 | 178  | 15495    | 0.296 | ng/ul  | 97       |
| 19) Fluoranthene            | 19.381 | 202  | 16405    | 0.302 | ng/ul  | 99       |
| 20) Pyrene                  | 19.744 | 202  | 16743m   | 0.296 | ng/ul  |          |
| 21) Benzo(a)anthracene      | 21.505 | 228  | 12065    | 0.287 | ng/ul  | 99       |
| 22) Chrysene                | 21.554 | 228  | 12936    | 0.296 | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 23.220 | 252  | 12859    | 0.265 | ng/ul  | 92       |
| 25) Benzo(k)fluoranthene    | 23.264 | 252  | 12910    | 0.265 | ng/ul# | 90       |
| 26) Benzo(a)pyrene          | 23.864 | 252  | 10637    | 0.253 | ng/ul  | 93       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.550 | 276  | 12323    | 0.268 | ng/ul# | 92       |
| 28) Dibenzo(a,h)anthracene  | 26.563 | 278  | 9518     | 0.273 | ng/ul# | 92       |
| 29) Benzo(g,h,i)perylene    | 27.348 | 276  | 9817     | 0.240 | ng/ul# | 91       |
| -----                       |        |      |          |       |        |          |

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

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