

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092223\  
 Data File : BM042001.D  
 Acq On : 22 Sep 2023 23:43  
 Operator : MA/JU  
 Sample : 04386-04  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0AF6

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 09/25/2023  
 Supervised By :mohammad ahmed 09/26/2023

Quant Time: Sep 23 02:40:44 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM092223.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Sep 22 13:50:58 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.954  | 152  | 5353     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.768 | 136  | 14179    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.597 | 164  | 7300     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.350 | 188  | 14516    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.521 | 240  | 5696     | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.930 | 264  | 4976     | 0.400 | ng/ul  | 0.00     |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.372  | 96   | 20970    | 2.875 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.352 | 152  | 2988     | 0.168 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.371 | 212  | 4854     | 0.207 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 14) Pentachlorophenol       | 16.978 | 266  | 218      | 0.028 | ng/ul# | 90       |
| 19) Fluoranthene            | 19.403 | 202  | 4153     | 0.065 | ng/ul# | 91       |
| 20) Pyrene                  | 19.771 | 202  | 3676     | 0.058 | ng/ul  | 95       |
| 21) Benzo(a)anthracene      | 21.507 | 228  | 1440     | 0.033 | ng/ul# | 89       |
| 22) Chrysene                | 21.559 | 228  | 1826m    | 0.041 | ng/ul  |          |
| 24) Benzo(b)fluoranthene    | 23.184 | 252  | 2334     | 0.061 | ng/ul# | 1        |
| 25) Benzo(k)fluoranthene    | 23.231 | 252  | 1146m    | 0.031 | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.825 | 252  | 1027     | 0.033 | ng/ul# | 1        |
| 27) Indeno(1,2,3-cd)pyrene  | 26.424 | 276  | 1120     | 0.026 | ng/ul# | 35       |
| 29) Benzo(g,h,i)perylene    | 27.206 | 276  | 1043     | 0.029 | ng/ul# | 8        |
| -----                       |        |      |          |       |        |          |

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

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