

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM081222\  
 Data File : BM036239.D  
 Acq On : 12 Aug 2022 20:35  
 Operator : CG/JU  
 Sample : N4069-16  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GBJM0

Quant Time: Aug 12 21:05:17 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM072522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Aug 12 11:23:11 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.816  | 152  | 159130   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.616 | 136  | 644395   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 363488   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.203 | 188  | 754285   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.380 | 240  | 850272   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.721 | 264  | 943612   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.216  | 96   | 19699    | 4.658  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.657  | 84   | 62533    | 5.398  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 7.004  | 99   | 90288    | 6.766  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.151  | 67   | 255775   | 28.752 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.351  | 132  | 266707   | 24.078 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.545  | 113  | 160347   | 14.832 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.986  | 128  | 164399   | 32.196 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.704  | 143  | 158741   | 31.909 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.251 | 165  | 259083   | 28.794 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.775 | 131  | 247939   | 16.277 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.880 | 166  | 881503   | 35.602 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.151 | 160  | 1076516  | 34.076 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.710 | 143  | 27944    | 5.496  | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.451 | 176  | 795639   | 35.561 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.580 | 200  | 125879   | 31.047 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.298 | 188  | 1390007  | 40.761 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.592 | 212  | 1758585  | 37.318 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.568 | 264  | 1966015  | 41.039 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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