

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061722\  
 Data File : BM035507.D  
 Acq On : 17 Jun 2022 23:12  
 Operator : CG/JU  
 Sample : N3226-02  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 EW4L3

Quant Time: Jun 18 00:12:48 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM061622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 17 01:58:49 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.810  | 152  | 220953   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.610 | 136  | 940559   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 560664   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.204 | 188  | 1133270  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.386 | 240  | 1142438  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.727 | 264  | 1248414  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.269  | 96   | 36479    | 6.591  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 394910   | 28.968 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.998  | 99   | 583990   | 34.487 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.146  | 67   | 383397   | 36.408 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.351  | 132  | 549274   | 37.449 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.540  | 113  | 493895   | 34.841 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.975  | 128  | 256101   | 36.579 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.698  | 143  | 281144   | 37.240 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.251 | 165  | 444789   | 34.700 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.757 | 131  | 682309   | 33.830 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.874 | 166  | 1496316  | 39.235 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.145 | 160  | 1810065  | 38.300 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.716 | 143  | 179295   | 30.605 | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.451 | 176  | 1312898  | 39.759 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.592 | 200  | 194301   | 28.759 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.304 | 188  | 2007041  | 39.400 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.598 | 212  | 2306940  | 41.074 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.574 | 264  | 2460606  | 40.864 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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