

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092622\  
 Data File : BM036754.D  
 Acq On : 26 Sep 2022 20:35  
 Operator : CG/JU  
 Sample : PB147882BL  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK882

Quant Time: Sep 26 23:16:10 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 26 23:10:22 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.004  | 152  | 447162   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.816 | 136  | 1958048  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.627 | 164  | 1248849  | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.368 | 188  | 2397747  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.533 | 240  | 1703302  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.968 | 264  | 1476545  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.381  | 96   | 55701    | 4.670  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.810  | 84   | 742444   | 21.342 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.157  | 99   | 1062885  | 26.598 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.328  | 67   | 670971   | 27.649 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.534  | 132  | 824394   | 28.655 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.710  | 113  | 856545   | 29.009 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.169  | 128  | 424151   | 32.004 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.892  | 143  | 410214   | 34.556 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.428 | 165  | 773365   | 28.111 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.951 | 131  | 1178759  | 25.771 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.045 | 166  | 2649218  | 31.453 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.321 | 160  | 2982404  | 29.851 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.821 | 143  | 393202   | 24.069 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.615 | 176  | 2271278  | 30.091 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.727 | 200  | 235309   | 24.805 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.462 | 188  | 3478712  | 31.731 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.750 | 212  | 3559063  | 42.473 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.815 | 264  | 2402890  | 33.310 | ng/ul | 0.01     |

Target Compounds Qvalue

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

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