

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092723\  
 Data File : BP017404.D  
 Acq On : 27 Sep 2023 18:51  
 Operator : MA/JU  
 Sample : 04359-08 10X  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 JLAD6

Quant Time: Sep 27 23:13:35 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 04:52:29 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.922  | 152  | 143066   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.746 | 136  | 623255   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.581 | 164  | 378246   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.351 | 188  | 844798   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.457 | 240  | 766262   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.969 | 264  | 831757   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.293  | 96   | 10256    | 2.945  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.728  | 84   | 141289   | 14.510 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.075  | 99   | 104011   | 8.852  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.264  | 67   | 361565   | 50.989 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.440  | 132  | 336662   | 36.125 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.634  | 113  | 206837   | 22.642 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.116  | 128  | 225449   | 50.036 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.834  | 143  | 232177   | 47.456 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.358 | 165  | 397450   | 42.999 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.910 | 131  | 466295   | 34.825 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.999 | 166  | 1492255  | 55.072 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.281 | 160  | 1586529  | 50.909 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.816 | 143  | 25339    | 5.368  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 1293575  | 55.453 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.710 | 200  | 173740   | 36.489 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.451 | 188  | 2097255  | 55.975 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.698 | 212  | 2468649  | 59.575 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.804 | 264  | 2338774  | 58.401 | ng/u1 | 0.00     |

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

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

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