

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092123\  
 Data File : BP017279.D  
 Acq On : 21 Sep 2023 21:46  
 Operator : MA/JU  
 Sample : 04387-12RE  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DFTPP653

Quant Time: Sep 22 00:01:20 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Sep 21 23:51:07 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.934  | 152  | 177278   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 715894   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.598 | 164  | 434499   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.369 | 188  | 979079   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.474 | 240  | 957346   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.998 | 264  | 1104697  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.305  | 96   | 37907    | 8.783  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.740  | 84   | 587028   | 48.652 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.099  | 99   | 808530   | 55.532 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.281  | 67   | 455340   | 51.821 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.463  | 132  | 632550   | 54.777 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.651  | 113  | 666535   | 58.883 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.140  | 128  | 321422   | 62.105 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.851  | 143  | 373221   | 66.414 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 665388   | 62.671 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.928 | 131  | 891090   | 57.939 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.016 | 166  | 2041792  | 65.597 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.298 | 160  | 2321187  | 64.840 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.833 | 143  | 349982   | 64.540 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.598 | 176  | 1767692  | 65.967 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.728 | 200  | 284783   | 51.607 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.469 | 188  | 2844827  | 65.514 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.716 | 212  | 3434076  | 66.332 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.833 | 264  | 3507264  | 65.941 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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