

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092223\  
 Data File : BP017292.D  
 Acq On : 22 Sep 2023 18:17  
 Operator : MA/JU  
 Sample : 04410-07  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BBBK1

Quant Time: Sep 22 22:29:45 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  | 133847   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.757 | 136  | 536892   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.592 | 164  | 315313   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.363 | 188  | 723093   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.469 | 240  | 709918   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.986 | 264  | 792385   | 20.000 | ng/ul | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.305  | 96   | 17848    | 5.477  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.740  | 84   | 135411   | 14.864 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.093  | 99   | 103202   | 9.388  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 268063   | 40.407 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.458  | 132  | 288000   | 33.032 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.646  | 113  | 187336   | 21.920 | ng/ul | -0.01    |
| 21) Nitrobenzene-d5           | 9.134  | 128  | 169073   | 43.560 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.846  | 143  | 174747   | 41.463 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.369 | 165  | 318296   | 39.975 | ng/ul | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.922 | 131  | 285848   | 24.783 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.010 | 166  | 1130921  | 50.067 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.293 | 160  | 1221509  | 47.019 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.834 | 143  | 31196    | 7.927  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.592 | 176  | 973828   | 50.078 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.722 | 200  | 147063   | 36.085 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.469 | 188  | 1718560  | 53.588 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.710 | 212  | 2103485  | 54.792 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.827 | 264  | 2098798  | 55.013 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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