

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090823\  
 Data File : BP017051.D  
 Acq On : 09 Sep 2023 14:34  
 Operator : MA/JU  
 Sample : 04227-15  
 Misc :  
 ALS Vial : 62 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHA06

Quant Time: Sep 09 23:06:45 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP090523.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Sep 09 01:41:21 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.969  | 152  | 325695   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.799 | 136  | 1402600  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.628 | 164  | 823498   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.392 | 188  | 1747457  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.492 | 240  | 1349542  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.027 | 264  | 1343987  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.322  | 96   | 49917    | 5.797  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.764  | 84   | 405670   | 16.489 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.116  | 99   | 286924   | 9.643  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.311  | 67   | 865400   | 46.497 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.493  | 132  | 766314   | 35.584 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.681  | 113  | 522749   | 23.060 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.169  | 128  | 491240   | 45.145 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.881  | 143  | 492459   | 41.749 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.404 | 165  | 843427   | 38.730 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 1111016  | 35.361 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.045 | 166  | 2932289  | 47.836 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.328 | 160  | 3327207  | 47.032 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.840 | 143  | 80664    | 6.737  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.622 | 176  | 2498066  | 47.597 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.745 | 200  | 370360   | 35.218 | ng/ul | -0.01    |
| 73) Anthracene-d10            | 17.492 | 188  | 3906383  | 48.677 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.733 | 212  | 4290264  | 52.193 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.863 | 264  | 3355006  | 49.206 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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