

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP101123\  
 Data File : BP017652.D  
 Acq On : 11 Oct 2023 18:02  
 Operator : MA/JU  
 Sample : 04744-06RX  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH583RX

Quant Time: Oct 12 02:23:57 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP100923.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 10 04:45:41 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.922  | 152  | 117861   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.758 | 136  | 476797   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.622 | 164  | 321427   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.410 | 188  | 807313   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.551 | 240  | 855378   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.122 | 264  | 996123   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.281  | 96   | 14598    | 5.025  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.723  | 84   | 102641   | 12.551 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.099  | 99   | 269      | 0.027  | ng/ul | 0.02     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.264  | 67   | 206700   | 33.439 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.452  | 132  | 24816    | 3.165  | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.634  | 113  | 14       | 0.002  | ng/ul | -0.01    |
| 21) Nitrobenzene-d5           | 9.134  | 128  | 119334   | 32.481 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.875  | 143  | 5776     | 1.431  | ng/ul | 0.02     |
| 28) 2,4-Dichlorophenol-d3     | 10.399 | 165  | 23436    | 2.990  | ng/ul | 0.02     |
| 31) 4-Chloroaniline-d4        | 10.922 | 131  | 155      | 0.014  | ng/ul | -0.01    |
| 46) Dimethylphthalate-d6      | 14.034 | 166  | 885538   | 34.699 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.316 | 160  | 472153   | 16.993 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.846 | 143  | 15       | 0.004  | ng/ul | -0.02    |
| 60) Fluorene-d10              | 15.622 | 176  | 833966   | 37.366 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.763 | 200  | 12       | 0.003  | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.510 | 188  | 917246   | 23.674 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.763 | 212  | 1482156  | 30.377 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.951 | 264  | 1329793  | 25.750 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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