

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP011524\  
 Data File : BP019375.D  
 Acq On : 15 Jan 2024 17:16  
 Operator : MA/JU  
 Sample : P1130-09  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Jan 15 23:51:43 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP010824.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jan 09 03:05:24 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.775  | 152  | 115528   | 20.000 | ng/ul | -0.01    |
| 20) Naphthalene-d8            | 10.569 | 136  | 471249   | 20.000 | ng/ul | -0.01    |
| 38) Acenaphthene-d10          | 14.416 | 164  | 316253   | 20.000 | ng/ul | -0.01    |
| 64) Phenanthrene-d10          | 17.174 | 188  | 786124   | 20.000 | ng/ul | -0.01    |
| 79) Chrysene-d12              | 21.280 | 240  | 807719   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.633 | 264  | 795958   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.216  | 96   | 11880    | 3.888  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.640  | 84   | 77157    | 8.733  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.963  | 99   | 56614    | 5.127  | ng/ul | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.116  | 67   | 210042   | 31.803 | ng/ul | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.310  | 132  | 174473   | 23.002 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.498  | 113  | 119229   | 13.611 | ng/ul | -0.02    |
| 21) Nitrobenzene-d5           | 8.940  | 128  | 127526   | 32.166 | ng/ul | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.657  | 143  | 143608   | 31.175 | ng/ul | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.198 | 165  | 268439   | 28.844 | ng/ul | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.722 | 131  | 279708   | 23.793 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.839 | 166  | 971182   | 34.759 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.110 | 160  | 997894   | 33.318 | ng/ul | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 20223    | 4.713  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.416 | 176  | 846279   | 36.996 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.557 | 200  | 181191   | 31.219 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.274 | 188  | 1417045  | 36.142 | ng/ul | -0.01    |
| 81) Pyrene-d10                | 19.521 | 212  | 1845490  | 35.040 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.480 | 264  | 1626761  | 37.587 | ng/ul | 0.00     |

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

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

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