

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122023\  
 Data File : BP018977.D  
 Acq On : 21 Dec 2023 00:32  
 Operator : MA/JU  
 Sample : 05894-10  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH3L6

Quant Time: Dec 21 01:04:41 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 19 05:28:03 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.822  | 152  | 113115   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.616 | 136  | 411611   | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.457 | 164  | 252183   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.210 | 188  | 573297   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.298 | 240  | 663196   | 20.000 | ng/u1 | #-0.02   |
| 88) Perylene-d12              | 23.662 | 264  | 836640   | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.252  | 96   | 12134    | 3.655  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.675  | 84   | 76070    | 8.556  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.998  | 99   | 57315    | 5.049  | ng/u1 | -0.02    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.163  | 67   | 193372   | 28.742 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.357  | 132  | 190000   | 21.451 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.539  | 113  | 115665   | 12.591 | ng/u1 | -0.02    |
| 21) Nitrobenzene-d5           | 8.986  | 128  | 130025   | 35.492 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.704  | 143  | 138242   | 36.397 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.245 | 165  | 240447   | 30.960 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.763 | 131  | 273058   | 26.903 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.874 | 166  | 910246   | 40.442 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.151 | 160  | 972954   | 39.258 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.680 | 143  | 21972    | 5.947  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.451 | 176  | 784621   | 41.530 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.580 | 200  | 140980   | 41.012 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.310 | 188  | 1303370  | 43.245 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.545 | 212  | 1694753  | 32.828 | ng/u1 | -0.01    |
| 92) Benzo(a)pyrene-d12        | 23.509 | 264  | 2112510  | 43.305 | ng/u1 | -0.02    |

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

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

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