

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080122\  
 Data File : BP011199.D  
 Acq On : 01 Aug 2022 14:45  
 Operator : CG/JU  
 Sample : N3790-17  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Aug 01 23:00:03 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP072822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 28 16:05:22 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.646  | 152  | 246376   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.440 | 136  | 1020127  | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.310 | 164  | 511479   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.086 | 188  | 962381   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.210 | 240  | 984219   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.551 | 264  | 1066352  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.152  | 96   | 13633    | 2.534  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.569  | 84   | 61363    | 4.032  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 6.828  | 99   | 423435   | 19.690 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.993  | 67   | 293945   | 19.331 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.181  | 132  | 329924   | 19.468 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.363  | 113  | 320258   | 18.205 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.810  | 128  | 155238   | 19.581 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.528  | 143  | 147792   | 19.461 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.063 | 165  | 257418   | 19.234 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.587 | 131  | 399418   | 18.410 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.734 | 166  | 770787   | 22.119 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.004 | 160  | 960206   | 21.927 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.539 | 143  | 65496    | 10.449 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.316 | 176  | 657572   | 22.456 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.445 | 200  | 48469    | 9.991  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.180 | 188  | 1024565  | 24.161 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.445 | 212  | 1195372  | 22.603 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.404 | 264  | 1253220  | 23.661 | ng/u1 | 0.00     |

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

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

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