

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112522\  
 Data File : BP012776.D  
 Acq On : 26 Nov 2022 19:52  
 Operator : CG/JU  
 Sample : N5602-05  
 Misc :  
 ALS Vial : 49 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AF7

Quant Time: Nov 27 21:53:20 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 25 15:59:22 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.022  | 152  | 561923   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.833 | 136  | 2515006  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.657 | 164  | 1657200  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.410 | 188  | 3468037  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.492 | 240  | 2531193  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 24.009 | 264  | 1873350  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.399  | 96   | 68854    | 5.251  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 507986   | 12.945 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.163  | 99   | 533707   | 11.558 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.340  | 67   | 932933   | 32.789 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.545  | 132  | 1077832  | 30.634 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.716  | 113  | 925499   | 24.877 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.186  | 128  | 599788   | 38.964 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.910  | 143  | 666212   | 40.760 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.445 | 165  | 1435498  | 37.641 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.969 | 131  | 1410054  | 27.000 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.063 | 166  | 4858950  | 42.597 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.351 | 160  | 5555878  | 41.744 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.833 | 143  | 180888   | 8.895  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.645 | 176  | 4246741  | 42.511 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.757 | 200  | 612185   | 36.933 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.510 | 188  | 6553389  | 42.928 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.739 | 212  | 6966596  | 47.367 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.850 | 264  | 4138290  | 44.973 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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