

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112922\  
 Data File : BP012868.D  
 Acq On : 30 Nov 2022 00:19  
 Operator : CG/JU  
 Sample : N5725-12  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHB61

Quant Time: Nov 30 00:50:14 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 29 22:53:34 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.011  | 152  | 620578   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.828 | 136  | 2846907  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.645 | 164  | 1976289  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.398 | 188  | 4212995  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.486 | 240  | 3465978  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.992 | 264  | 2900806  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 55180    | 3.810  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 368921   | 8.512  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.158  | 99   | 272345   | 5.341  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.334  | 67   | 882489   | 28.084 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.534  | 132  | 841802   | 21.664 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.710  | 113  | 564093   | 13.729 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.175  | 128  | 558741   | 32.066 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.905  | 143  | 595691   | 32.197 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.440 | 165  | 1105150  | 25.600 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 1179554  | 19.953 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.051 | 166  | 4248134  | 31.229 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.340 | 160  | 4709596  | 29.672 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.828 | 143  | 81320    | 3.353  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.640 | 176  | 3732626  | 31.332 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.745 | 200  | 505729   | 25.116 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.498 | 188  | 6001734  | 32.363 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.728 | 212  | 6619561  | 32.869 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.833 | 264  | 4641619  | 32.576 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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