

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110222\  
 Data File : BP012478.D  
 Acq On : 03 Nov 2022 07:27  
 Operator : CG/JU  
 Sample : N5321-23  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DBWW5

Quant Time: Nov 03 07:55:15 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Nov 03 00:09:04 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.781  | 152  | 283755   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.581 | 136  | 1337514  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.434 | 164  | 900684   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.198 | 188  | 1951399  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.292 | 240  | 1842034  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.674 | 264  | 1639388  | 20.000 | ng/u1 | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.234  | 96   | 26175    | 3.714  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.664  | 84   | 99624    | 4.921  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.958  | 99   | 213828   | 8.402  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67   | 457878   | 30.137 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.311  | 132  | 511873   | 27.541 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.499  | 113  | 410725   | 20.405 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.952  | 128  | 320901   | 36.998 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.669  | 143  | 365732   | 40.626 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.210 | 165  | 685854   | 35.154 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.734 | 131  | 482064   | 16.594 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.851 | 166  | 2379377  | 38.591 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.128 | 160  | 2433153  | 34.532 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.669 | 143  | 96352    | 8.257  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.434 | 176  | 2029698  | 38.449 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.563 | 200  | 406911   | 40.051 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.298 | 188  | 3306436  | 39.324 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.539 | 212  | 3751858  | 40.568 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.522 | 264  | 3260504  | 40.704 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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