

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP063022\  
 Data File : BP010916.D  
 Acq On : 30 Jun 2022 12:37  
 Operator : CG/JU  
 Sample : N3523-05  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0JW9

Quant Time: Jul 01 02:32:38 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 27 23:12:49 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.705  | 152  | 402963   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.499 | 136  | 1599469  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.357 | 164  | 795734   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.122 | 188  | 1474436  | 20.000 | ng/u1 | #-0.01   |
| 79) Chrysene-d12              | 21.239 | 240  | 1495523  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.598 | 264  | 1644041  | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.205  | 96   | 50914    | 4.780  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.617  | 84   | 147808   | 5.213  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.881  | 99   | 218025   | 6.835  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.046  | 67   | 642116   | 30.780 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.240  | 132  | 653170   | 24.680 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.416  | 113  | 392955   | 15.583 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.863  | 128  | 368014   | 36.143 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.587  | 143  | 327733   | 38.905 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.122 | 165  | 562516   | 27.025 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.646 | 131  | 613715   | 18.063 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.781 | 166  | 1701494  | 31.215 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.051 | 160  | 2140405  | 30.767 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.575 | 143  | 49495    | 5.813  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.357 | 176  | 1520438  | 32.504 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.481 | 200  | 151037   | 30.919 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.222 | 188  | 2534648  | 38.638 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.475 | 212  | 2898149  | 35.570 | ng/u1 | -0.01    |
| 92) Benzo(a)pyrene-d12        | 23.445 | 264  | 3293926  | 40.545 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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