

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062422\  
 Data File : BP010787.D  
 Acq On : 24 Jun 2022 11:38  
 Operator : CG/JU  
 Sample : N3395-10  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0157

Quant Time: Jun 24 22:44:57 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 24 22:42:26 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.722  | 152  | 360994   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.516 | 136  | 1465349  | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.375 | 164  | 780968   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.139 | 188  | 1404854  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.251 | 240  | 1168450  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.627 | 264  | 1306355  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.211  | 96   | 57536    | 6.029  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.611  | 84   | 331604   | 13.054 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.887  | 99   | 368537   | 12.896 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.063  | 67   | 760577   | 40.697 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.252  | 132  | 771136   | 32.525 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.428  | 113  | 533530   | 23.618 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.881  | 128  | 421021   | 45.133 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.604  | 143  | 345734   | 44.798 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.140 | 165  | 704940   | 36.968 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.657 | 131  | 1052104  | 33.800 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.798 | 166  | 2352788  | 43.980 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.069 | 160  | 2878654  | 42.162 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.575 | 143  | 77187    | 9.237  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.381 | 176  | 1979153  | 43.111 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.492 | 200  | 150933   | 32.428 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.239 | 188  | 2949878  | 47.195 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.492 | 212  | 3048812  | 47.894 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.474 | 264  | 3108808  | 48.158 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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