

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP071822\  
 Data File : BP011005.D  
 Acq On : 18 Jul 2022 22:11  
 Operator : CG/JU  
 Sample : N3710-02  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 H0B45

Quant Time: Jul 19 00:47:33 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071422.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 14 15:34:34 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.681  | 152  | 386927   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.475 | 136  | 1589905  | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.340 | 164  | 868541   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.104 | 188  | 1739968  | 20.000 | ng/u1 | #-0.01   |
| 79) Chrysene-d12              | 21.222 | 240  | 1723938  | 20.000 | ng/u1 | #-0.02   |
| 88) Perylene-d12              | 23.574 | 264  | 1814784  | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.181  | 96   | 53313    | 4.967  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.593  | 84   | 324465   | 11.696 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.858  | 99   | 233028   | 7.337  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.022  | 67   | 747151   | 36.583 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.216  | 132  | 671317   | 26.158 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.393  | 113  | 431396   | 16.966 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.840  | 128  | 430481   | 37.154 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.563  | 143  | 368354   | 32.129 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.099 | 165  | 629553   | 30.026 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.622 | 131  | 908425   | 26.281 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.763 | 166  | 2473547  | 41.669 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.028 | 160  | 2593353  | 37.027 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.557 | 143  | 57574    | 4.792  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.340 | 176  | 1887739  | 37.412 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.463 | 200  | 236119   | 27.424 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.204 | 188  | 3154974  | 42.652 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.463 | 212  | 3654362  | 40.361 | ng/u1 | -0.01    |
| 92) Benzo(a)pyrene-d12        | 23.427 | 264  | 3786882  | 43.446 | ng/u1 | -0.02    |

Target Compounds Qvalue

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

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