

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP071922\  
 Data File : BP011042.D  
 Acq On : 19 Jul 2022 20:44  
 Operator : CG/JU  
 Sample : N3753-12  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 GBHQ3

Quant Time: Jul 19 23:41:55 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071422.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 19 03:25:54 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.675  | 152  | 329350   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.469 | 136  | 1337125  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.340 | 164  | 760138   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.104 | 188  | 1513738  | 20.000 | ng/ul | # 0.00   |
| 79) Chrysene-d12              | 21.227 | 240  | 1469588  | 20.000 | ng/ul | # 0.00   |
| 88) Perylene-d12              | 23.574 | 264  | 1490398  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.181  | 96   | 43010    | 4.708  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.593  | 84   | 265100   | 11.227 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.852  | 99   | 175315   | 6.485  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.016  | 67   | 639587   | 36.791 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.211  | 132  | 581721   | 26.629 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.393  | 113  | 342693   | 15.834 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.840  | 128  | 371689   | 38.144 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.557  | 143  | 328908   | 34.112 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.093 | 165  | 548332   | 31.097 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.616 | 131  | 864727   | 29.747 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.757 | 166  | 2010236  | 38.694 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.028 | 160  | 2453218  | 40.021 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.557 | 143  | 32318    | 3.074  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.340 | 176  | 1732741  | 39.238 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.457 | 200  | 211074   | 28.179 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.204 | 188  | 2886993  | 44.862 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.463 | 212  | 3316555  | 42.970 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.427 | 264  | 3225721  | 45.063 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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