

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072622\  
 Data File : BP011106.D  
 Acq On : 26 Jul 2022 12:45  
 Operator : CG/JU  
 Sample : N3801-17  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 GBHY2

Quant Time: Jul 26 14:44:27 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071422.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 26 14:31:36 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.664  | 152  | 505427   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.458 | 136  | 2319778  | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.328 | 164  | 1364351  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.092 | 188  | 2441843  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.210 | 240  | 1478385  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.551 | 264  | 1439496  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.164  | 96   | 59426    | 4.239  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.576  | 84   | 285417   | 7.876  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.840  | 99   | 290869   | 7.011  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.005  | 67   | 924496   | 34.653 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.193  | 132  | 869798   | 25.945 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.375  | 113  | 595573   | 17.931 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.828  | 128  | 580760   | 34.354 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.546  | 143  | 580709   | 34.715 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.081 | 165  | 944880   | 30.887 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.605 | 131  | 919651   | 18.235 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.751 | 166  | 3474669  | 37.262 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.016 | 160  | 3999434  | 36.351 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.546 | 143  | 62534    | 3.314  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.328 | 176  | 2748243  | 34.673 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.451 | 200  | 396511   | 32.816 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.192 | 188  | 3765034  | 36.269 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.451 | 212  | 3535208  | 45.530 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.404 | 264  | 2588589  | 37.441 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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