

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072522\  
 Data File : BP011083.D  
 Acq On : 25 Jul 2022 09:50  
 Operator : CG/JU  
 Sample : PB146370BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK370

Quant Time: Jul 25 11:42:36 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     | 7.669  | 152  | 409816   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.463 | 136  | 1910072  | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.334 | 164  | 1349746  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.104 | 188  | 2773855  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.222 | 240  | 2178272  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.568 | 264  | 1751254  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.164  | 96   | 78813    | 6.933  | ng/uL | -0.02    |
| 4) Pyridine-d5                | 3.575  | 84   | 1090223  | 37.105 | ng/u1 | -0.02    |
| 7) Phenol-d5                  | 6.852  | 99   | 1281995  | 38.109 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.016  | 67   | 875092   | 40.454 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.205  | 132  | 1058537  | 38.942 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.387  | 113  | 1093733  | 40.612 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.834  | 128  | 564019   | 40.520 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.557  | 143  | 592160   | 42.993 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.093 | 165  | 941122   | 37.363 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.616 | 131  | 1587619  | 38.232 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.757 | 166  | 3872219  | 41.975 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.028 | 160  | 4507994  | 41.417 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.557 | 143  | 607968   | 32.563 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.339 | 176  | 3150399  | 40.177 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.469 | 200  | 522148   | 38.041 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.204 | 188  | 4745456  | 40.242 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.457 | 212  | 5114597  | 44.707 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.421 | 264  | 3430301  | 40.783 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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