

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112822\  
 Data File : BP012830.D  
 Acq On : 28 Nov 2022 23:21  
 Operator : CG/JU  
 Sample : PB149200BL  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK200

Quant Time: Nov 29 00:46:57 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Nov 27 22:55:08 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.016  | 152  | 697453   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.834 | 136  | 3104724  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.651 | 164  | 2089899  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.404 | 188  | 4164876  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.486 | 240  | 2594856  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 24.004 | 264  | 2222323  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.399  | 96   | 82819    | 5.088  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 1126602  | 23.130 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.163  | 99   | 1388862  | 24.233 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.340  | 67   | 909785   | 25.762 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.540  | 132  | 1123481  | 25.726 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.716  | 113  | 1131230  | 24.498 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.181  | 128  | 536098   | 28.212 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.904  | 143  | 599995   | 29.736 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.440 | 165  | 1136989  | 24.151 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.963 | 131  | 1627893  | 25.251 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.057 | 166  | 3817981  | 26.541 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.345 | 160  | 4273067  | 25.458 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.828 | 143  | 492911   | 19.219 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.639 | 176  | 3330887  | 26.440 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.751 | 200  | 401607   | 20.175 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.504 | 188  | 4995718  | 27.249 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.727 | 212  | 4939357  | 32.760 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.839 | 264  | 2956549  | 27.085 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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