

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111423\  
 Data File : BP018151.D  
 Acq On : 14 Nov 2023 12:13  
 Operator : MA/JU  
 Sample : PB156906TB 10X  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Nov 15 00:37:30 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP111023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 13 15:50:18 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.787  | 152  | 71240    | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.604 | 136  | 297597   | 20.000 | ng/ul | # 0.00   |
| 38) Acenaphthene-d10          | 14.469 | 164  | 154568   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.251 | 188  | 403701   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.380 | 240  | 357122   | 20.000 | ng/ul | -0.01    |
| 88) Perylene-d12              | 23.845 | 264  | 400602   | 20.000 | ng/ul | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.158  | 96   | 14309    | 7.085  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.593  | 84   | 126056   | 23.940 | ng/ul | -0.01    |
| 7) Phenol-d5                  | 6.952  | 99   | 210671   | 30.631 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.140  | 67   | 144790   | 35.823 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.310  | 132  | 169775   | 31.172 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.510  | 113  | 168887   | 29.967 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.999  | 128  | 84647    | 31.571 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.710  | 143  | 81870    | 26.953 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.234 | 165  | 145321   | 26.843 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.787 | 131  | 193647   | 26.083 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.892 | 166  | 477770   | 33.500 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.163 | 160  | 489637   | 31.521 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.739 | 143  | 36259    | 17.014 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.469 | 176  | 388792   | 33.020 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.634 | 200  | 39697    | 13.602 | ng/ul | -0.03    |
| 73) Anthracene-d10            | 17.351 | 188  | 733757   | 33.607 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.604 | 212  | 838528   | 35.152 | ng/ul | -0.01    |
| 92) Benzo(a)pyrene-d12        | 23.686 | 264  | 789595   | 33.897 | ng/ul | -0.02    |

Target Compounds Qvalue

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

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