

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072621\  
 Data File : BP006333.D  
 Acq On : 26 Jul 2021 12:03  
 Operator : CG/JU  
 Sample : PB137844TB 10X  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLEB844

Quant Time: Jul 26 13:10:28 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 26 12:12:19 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.787  | 152  | 245738   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.569 | 136  | 1030179  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.416 | 164  | 592192   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.169 | 188  | 1163712  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.269 | 240  | 1103923  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.615 | 264  | 1124645  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 48021    | 7.294  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.705  | 84   | 698760   | 35.754 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.958  | 99   | 775628   | 33.607 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 500758   | 34.871 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.316  | 132  | 604122   | 35.233 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.493  | 113  | 603275   | 33.239 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.940  | 128  | 282149   | 36.783 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.657  | 143  | 279992   | 38.228 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.193 | 165  | 483956   | 33.720 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.710 | 131  | 814110   | 34.454 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.834 | 166  | 1622890  | 38.505 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.110 | 160  | 1892619  | 36.550 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.616 | 143  | 283014   | 33.818 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.410 | 176  | 1308642  | 36.934 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.528 | 200  | 185246   | 28.555 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.269 | 188  | 2079998  | 39.595 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.510 | 212  | 2273212  | 40.035 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.468 | 264  | 2206276  | 38.604 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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