

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072721\  
 Data File : BP006357.D  
 Acq On : 27 Jul 2021 10:17  
 Operator : CG/JU  
 Sample : PB137867BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK867

Quant Time: Jul 27 11:38:34 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jul 24 04:29:12 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.787  | 152  | 290583   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.569 | 136  | 1217920  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.410 | 164  | 683747   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.163 | 188  | 1324875  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.263 | 240  | 1248800  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.610 | 264  | 1281049  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.305  | 96   | 49131    | 6.311  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.705  | 84   | 696276   | 30.129 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.958  | 99   | 795294   | 29.141 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 518364   | 30.526 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.316  | 132  | 614698   | 30.317 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.493  | 113  | 622419   | 29.001 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.940  | 128  | 285669   | 31.501 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.657  | 143  | 281024   | 32.455 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.187 | 165  | 487334   | 28.721 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.710 | 131  | 819872   | 29.350 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.834 | 166  | 1544294  | 31.734 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.104 | 160  | 1882568  | 31.487 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.610 | 143  | 272552   | 28.207 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.404 | 176  | 1279344  | 31.272 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200  | 177004   | 23.965 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.263 | 188  | 2000850  | 33.455 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.504 | 212  | 2176924  | 33.892 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.457 | 264  | 2181214  | 33.506 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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