

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081221\  
 Data File : BP006659.D  
 Acq On : 12 Aug 2021 10:21  
 Operator : CG/JU  
 Sample : PB138304BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK304

Quant Time: Aug 12 11:04:34 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP081021.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 10 15:09:30 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.710  | 152  | 134590   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.499 | 136  | 563256   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.351 | 164  | 336436   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.110 | 188  | 692983   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.210 | 240  | 696295   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.527 | 264  | 728677   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.222  | 96   | 28901    | 7.263  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.628  | 84   | 397559   | 34.755 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.893  | 99   | 432422   | 31.960 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.058  | 67   | 274594   | 33.023 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.246  | 132  | 347636   | 34.498 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.428  | 113  | 346698   | 32.074 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.869  | 128  | 170006   | 35.557 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.587  | 143  | 180484   | 36.075 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.122 | 165  | 280439   | 33.177 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.640 | 131  | 463679   | 33.750 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.775 | 166  | 1003675  | 38.901 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.045 | 160  | 1161449  | 36.699 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.569 | 143  | 185113   | 34.348 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.351 | 176  | 788074   | 37.556 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.481 | 200  | 132536   | 28.934 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.210 | 188  | 1293922  | 39.463 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.457 | 212  | 1442938  | 39.462 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.380 | 264  | 1473020  | 37.487 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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