

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM060821\  
 Data File : BM030319.D  
 Acq On : 08 Jun 2021 12:14  
 Operator : CG/JU  
 Sample : PB136605BL  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK605

Quant Time: Jun 08 12:59:59 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM060621.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jun 08 09:51:19 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.851  | 152  | 223699   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.639 | 136  | 948738   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.474 | 164  | 612128   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.209 | 188  | 1243918  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.368 | 240  | 1054640  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.650 | 264  | 984940   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.316  | 96   | 35647    | 6.085  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.728  | 84   | 430749   | 27.982 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.004  | 99   | 563756   | 28.996 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.181  | 67   | 375351   | 29.913 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.381  | 132  | 463583   | 32.970 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.545  | 113  | 450767   | 29.030 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.004  | 128  | 212506   | 32.722 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.728  | 143  | 213017   | 33.026 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.257 | 165  | 433650   | 30.865 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.769 | 131  | 635183   | 29.574 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.880 | 166  | 1443631  | 34.366 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.168 | 160  | 1766799  | 33.667 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.651 | 143  | 215860   | 27.453 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.463 | 176  | 1269690  | 33.761 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.574 | 200  | 178510   | 24.861 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.309 | 188  | 1939159  | 34.274 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.592 | 212  | 2078715  | 37.658 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.509 | 264  | 1770026  | 36.040 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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