

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092521\  
 Data File : BM032154.D  
 Acq On : 25 Sep 2021 15:09  
 Operator : CG/JU  
 Sample : M3919-05  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GB6Y4

Quant Time: Sep 26 00:20:43 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM092421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Sep 25 01:15:33 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.666  | 152  | 407612   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.460 | 136  | 1720514  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.313 | 164  | 1090138  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.036 | 188  | 2189872  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.195 | 240  | 1888692  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.365 | 264  | 1685849  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 2.990  | 96   | 52066    | 5.393  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.413  | 84   | 271634   | 9.912  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.843  | 99   | 318456   | 9.842  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.990  | 67   | 729928   | 37.519 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.201  | 132  | 789978   | 31.449 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.390  | 113  | 566426   | 22.388 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.819  | 128  | 455780   | 44.353 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.542  | 143  | 485111   | 44.645 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.090 | 165  | 888831   | 35.782 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.589 | 131  | 1148922  | 34.222 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.713 | 166  | 2994795  | 42.282 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.001 | 160  | 3586351  | 39.853 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.507 | 143  | 96077    | 8.339  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.301 | 176  | 2537308  | 41.983 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.413 | 200  | 358491   | 34.640 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.136 | 188  | 3929190  | 43.002 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.418 | 212  | 4252049  | 42.789 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.230 | 264  | 3371927  | 42.619 | ng/u1 | 0.00     |

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

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