

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM082222\  
 Data File : BM036369.D  
 Acq On : 22 Aug 2022 16:51  
 Operator : CG/JU  
 Sample : PB147152BL  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK152

Quant Time: Aug 23 01:03:45 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 23 00:55:01 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.792  | 152  | 223290   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.592 | 136  | 977717   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.439 | 164  | 528143   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.180 | 188  | 909085   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.362 | 240  | 742881   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.691 | 264  | 760512   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.205  | 96   | 35876    | 6.123  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.628  | 84   | 523172   | 32.380 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.975  | 99   | 644895   | 33.836 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.134  | 67   | 441294   | 36.783 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.328  | 132  | 520480   | 34.756 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.516  | 113  | 503980   | 35.414 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.963  | 128  | 265075   | 35.759 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.686  | 143  | 258682   | 34.403 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.222 | 165  | 439180   | 32.688 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.745 | 131  | 709228   | 34.051 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.857 | 166  | 1269160  | 37.984 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.133 | 160  | 1529772  | 36.708 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.669 | 143  | 141485   | 22.348 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.427 | 176  | 1091099  | 36.138 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.563 | 200  | 105702   | 22.616 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.280 | 188  | 1518331  | 37.562 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.574 | 212  | 1564151  | 40.589 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.539 | 264  | 1443166  | 38.704 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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