

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM081722\  
 Data File : BM036338.D  
 Acq On : 18 Aug 2022 04:48  
 Operator : CG/JU  
 Sample : N4176-08  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 EX7Q3

Quant Time: Aug 18 05:21:25 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 16 15:00:33 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.798  | 152  | 270306   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.598 | 136  | 1161934  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.439 | 164  | 629390   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.186 | 188  | 1155165  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.368 | 240  | 1022877  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.697 | 264  | 1084608  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.210  | 96   | 24512    | 3.456  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.640  | 84   | 112728   | 5.763  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.981  | 99   | 444587   | 19.269 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.134  | 67   | 291621   | 20.080 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.328  | 132  | 370675   | 20.447 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.522  | 113  | 353081   | 20.495 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.969  | 128  | 186326   | 21.150 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.686  | 143  | 188405   | 21.084 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.228 | 165  | 315275   | 19.746 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.745 | 131  | 482783   | 19.504 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.863 | 166  | 935487   | 23.494 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.133 | 160  | 1097693  | 22.103 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.674 | 143  | 114372   | 15.159 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.433 | 176  | 816972   | 22.706 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.568 | 200  | 66237    | 11.153 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.280 | 188  | 1343836  | 26.163 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.574 | 212  | 1490716  | 28.095 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.544 | 264  | 1484243  | 27.911 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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