

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM080324\  
 Data File : BM046972.D  
 Acq On : 03 Aug 2024 20:09  
 Operator : MA/JU  
 Sample : P3276-15  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 E20J4

Quant Time: Aug 05 01:30:49 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM080224.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Aug 03 02:51:11 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.399  | 152  | 179800   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.157 | 136  | 700388   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.057 | 164  | 434789   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.816 | 188  | 897977   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.051 | 240  | 869371   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.762 | 264  | 1124929  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.022  | 96   | 11066    | 2.438  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1 |          |
| 7) Phenol-d5                  | 6.604  | 99   | 167658   | 12.275 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.752  | 67   | 135303   | 15.055 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 6.940  | 132  | 127804   | 11.046 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.116  | 113  | 125099   | 11.562 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.546  | 128  | 77107    | 13.720 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.257  | 143  | 52776    | 8.599  | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.793  | 165  | 104741   | 8.889  | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.304 | 131  | 178725   | 11.540 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.480 | 166  | 463500   | 13.942 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.739 | 160  | 530906   | 14.307 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.310 | 143  | 1219     | 0.190  | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.063 | 176  | 404927   | 14.239 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.198 | 200  | 1150     | 0.203  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 16.916 | 188  | 621547   | 14.601 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.227 | 212  | 627557   | 12.610 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.580 | 264  | 562783   | 9.501  | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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