

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110222\  
 Data File : BM037316.D  
 Acq On : 02 Nov 2022 15:02  
 Operator : CG/JU  
 Sample : N5301-09  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBWU6

Quant Time: Nov 02 22:53:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM110122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 02 03:33:37 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.857  | 152  | 311512   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.657 | 136  | 1370333  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.492 | 164  | 871843   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.233 | 188  | 1774508  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.403 | 240  | 1522096  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.750 | 264  | 1354331  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.275  | 96   | 30811    | 4.360  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.699  | 84   | 248154   | 11.764 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.028  | 99   | 173716   | 6.408  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.193  | 67   | 480214   | 28.866 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.387  | 132  | 468487   | 22.618 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.575  | 113  | 330784   | 14.942 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.028  | 128  | 298937   | 29.204 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.745  | 143  | 301229   | 27.016 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.281 | 165  | 554610   | 26.979 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.804 | 131  | 735007   | 23.397 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.910 | 166  | 2055732  | 34.102 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.186 | 160  | 2372120  | 32.732 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.716 | 143  | 58756    | 4.876  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.480 | 176  | 1882945  | 35.027 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.610 | 200  | 277279   | 24.670 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.327 | 188  | 2966589  | 36.300 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.621 | 212  | 3243988  | 40.785 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.603 | 264  | 2517833  | 36.116 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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