

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM091922\  
 Data File : BM036598.D  
 Acq On : 20 Sep 2022 02:00  
 Operator : CG/JU  
 Sample : N4576-01  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 CB8F6

Quant Time: Sep 20 02:32:21 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 16 02:39:54 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.028  | 152  | 347029   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.839 | 136  | 1557194  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.651 | 164  | 999901   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.386 | 188  | 2030075  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.551 | 240  | 1931654  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.992 | 264  | 1684398  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 41560    | 4.489  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 174469   | 6.462  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.175  | 99   | 187767   | 6.055  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.351  | 67   | 558708   | 29.666 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.551  | 132  | 562327   | 25.186 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.728  | 113  | 364971   | 15.927 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.192  | 128  | 368835   | 34.994 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.916  | 143  | 351043   | 37.184 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.451 | 165  | 656666   | 30.014 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.969 | 131  | 652925   | 17.949 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.063 | 166  | 2447341  | 36.290 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.345 | 160  | 2783985  | 34.803 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.827 | 143  | 75481    | 5.771  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.639 | 176  | 2145851  | 35.507 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.745 | 200  | 278727   | 34.703 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.486 | 188  | 3586853  | 38.643 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.768 | 212  | 4267269  | 44.905 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.833 | 264  | 3484945  | 42.348 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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