

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110322\  
 Data File : BM037356.D  
 Acq On : 03 Nov 2022 20:51  
 Operator : CG/JU  
 Sample : N5276-16  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 EW4X6

Quant Time: Nov 04 01:06:28 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 | 7.851  | 152  | 402630   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.651 | 136  | 1784547  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.486 | 164  | 1152071  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.227 | 188  | 2312295  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.397 | 240  | 2067205  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.744 | 264  | 1761911  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.275  | 96  | 43633   | 4.777  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.699  | 84  | 155751  | 5.712  | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.022  | 99  | 240165  | 6.854  | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.187  | 67  | 678627  | 31.561 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.381  | 132 | 664235  | 24.811 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.569  | 113 | 446044  | 15.589 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 9.022  | 128 | 436792  | 32.767 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.739  | 143 | 434910  | 29.951 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.275 | 165 | 729712  | 27.258 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.798 | 131 | 855520  | 20.912 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.904 | 166 | 2651798 | 33.290 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.180 | 160 | 3160059 | 32.998 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.710 | 143 | 85943   | 5.398  | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.474 | 176 | 2498716 | 35.176 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.604 | 200 | 400821  | 27.368 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.327 | 188 | 4087974 | 38.388 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.615 | 212 | 4537319 | 42.003 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.591 | 264 | 3505655 | 38.653 | ng/u1 | -0.01 |

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

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

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