

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092322\  
 Data File : BM036685.D  
 Acq On : 23 Sep 2022 20:00  
 Operator : CG/JU  
 Sample : N4708-07  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 CB8K7

Quant Time: Sep 23 23:37:03 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM091522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 21 22:42:41 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.016  | 152  | 445125   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.828 | 136  | 2011692  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.639 | 164  | 1290107  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.374 | 188  | 2576720  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.544 | 240  | 2074799  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.980 | 264  | 1780802  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.393  | 96  | 45035   | 3.793  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.828  | 84  | 131105  | 3.786  | ng/u1 | 0.01  |
| 7) Phenol-d5                  | 7.169  | 99  | 215381  | 5.415  | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.340  | 67  | 557722  | 23.087 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.540  | 132 | 578220  | 20.190 | ng/u1 | -0.01 |
| 15) 4-Methylphenol-d8         | 8.722  | 113 | 403232  | 13.719 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 9.181  | 128 | 366052  | 26.884 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.904  | 143 | 354842  | 29.094 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.445 | 165 | 668085  | 23.637 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.963 | 131 | 405337  | 8.625  | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 14.051 | 166 | 2543155 | 29.228 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.333 | 160 | 2751493 | 26.659 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.833 | 143 | 80859   | 4.791  | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.627 | 176 | 2230490 | 28.605 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.739 | 200 | 261425  | 25.643 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.474 | 188 | 3497528 | 29.687 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.756 | 212 | 3915511 | 38.360 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.827 | 264 | 2813032 | 32.333 | ng/u1 | 0.00  |

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

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

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