

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041423\  
 Data File : BN024898.D  
 Acq On : 14 Apr 2023 15:17  
 Operator : CG/JU  
 Sample : 02252-06  
 Misc :  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 COME2

Quant Time: Apr 14 22:47:50 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041023.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 14 22:34:56 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.940  | 152  | 303308   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.752 | 136  | 1320064  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.587 | 164  | 838168   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.334 | 188  | 1799371  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.522 | 240  | 1696622  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.980 | 264  | 1864483  | 20.000 | ng/u1 | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.311  | 96  | 27900   | 3.501  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.746  | 84  | 120469  | 5.048  | ng/u1 | 0.01 |
| 7) Phenol-d5                  | 7.099  | 99  | 157223  | 5.317  | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.269  | 67  | 505123  | 25.539 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.469  | 132 | 437504  | 20.132 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.652  | 113 | 303847  | 12.943 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.111  | 128 | 284632  | 26.630 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.834  | 143 | 300525  | 25.958 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.369 | 165 | 495495  | 23.945 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.893 | 131 | 334368  | 10.648 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.993 | 166 | 1852614 | 28.512 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.281 | 160 | 1998814 | 26.845 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.787 | 143 | 57905   | 4.347  | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.575 | 176 | 1590852 | 28.998 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.693 | 200 | 287466  | 24.548 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.428 | 188 | 2705034 | 32.008 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.728 | 212 | 3195078 | 32.431 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.827 | 264 | 3196675 | 32.791 | ng/u1 | 0.00 |

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

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

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