

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN012323\  
 Data File : BN023757.D  
 Acq On : 23 Jan 2023 18:07  
 Operator : CG/JU  
 Sample : 01226-01  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 EX9F1

Quant Time: Jan 24 00:01:48 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN010523.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 20 23:02:18 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.887  | 152  | 167330   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.698 | 136  | 710980   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.539 | 164  | 439153   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.280 | 188  | 848271   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.480 | 240  | 1001030  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.892 | 264  | 1095366  | 20.000 | ng/ul | -0.02    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.269  | 96  | 16918   | 4.351  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.699  | 84  | 107259  | 9.194  | ng/ul | 0.00  |
| 7) Phenol-d5                  | 7.052  | 99  | 90682   | 6.115  | ng/ul | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.222  | 67  | 261232  | 28.714 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.416  | 132 | 258989  | 22.803 | ng/ul | 0.00  |
| 15) 4-Methylphenol-d8         | 8.599  | 113 | 164737  | 13.917 | ng/ul | -0.01 |
| 21) Nitrobenzene-d5           | 9.057  | 128 | 163126  | 30.427 | ng/ul | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.781  | 143 | 167676  | 29.544 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.316 | 165 | 278914  | 25.765 | ng/ul | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.840 | 131 | 408321  | 23.979 | ng/ul | 0.00  |
| 46) Dimethylphthalate-d6      | 13.951 | 166 | 1103049 | 33.765 | ng/ul | 0.00  |
| 49) Acenaphthylene-d8         | 14.234 | 160 | 1136292 | 31.393 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.734 | 143 | 36005   | 5.414  | ng/ul | 0.00  |
| 60) Fluorene-d10              | 15.528 | 176 | 929871  | 33.312 | ng/ul | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.651 | 200 | 155596  | 30.407 | ng/ul | 0.00  |
| 73) Anthracene-d10            | 17.380 | 188 | 1393380 | 36.707 | ng/ul | 0.00  |
| 81) Pyrene-d10                | 19.680 | 212 | 1980676 | 34.597 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.727 | 264 | 2041196 | 37.682 | ng/ul | -0.03 |

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

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

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