

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041024\  
 Data File : BN030832.D  
 Acq On : 11 Apr 2024 01:45  
 Operator : MA/JU  
 Sample : P2014-03  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 E0MD0

Quant Time: Apr 11 02:35:19 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN032824.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Apr 09 12:10:11 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.675  | 152  | 158290   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.457 | 136  | 739117   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.316 | 164  | 509442   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.069 | 188  | 1154077  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.304 | 240  | 1130779  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.233 | 264  | 1221768  | 20.000 | ng/u1 | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.187  | 96  | 14428   | 4.274  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.593  | 84  | 135609  | 13.397 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 6.846  | 99  | 102277  | 7.987  | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.016  | 67  | 287961  | 36.410 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.210  | 132 | 294786  | 30.347 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.381  | 113 | 204373  | 19.044 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 8.834  | 128 | 202658  | 39.526 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.546  | 143 | 198620  | 42.346 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.081 | 165 | 358273  | 34.891 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.604 | 131 | 532132  | 31.478 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.734 | 166 | 1397396 | 41.084 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.010 | 160 | 1617067 | 41.350 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.528 | 143 | 35578   | 5.468  | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.316 | 176 | 1273922 | 42.800 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.434 | 200 | 164190  | 32.232 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.169 | 188 | 2169755 | 43.635 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.469 | 212 | 2664734 | 43.831 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 24.033 | 264 | 2547399 | 45.250 | ng/u1 | 0.00 |

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

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

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