

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN021424\  
 Data File : BN029749.D  
 Acq On : 14 Feb 2024 01:28  
 Operator : MA/JU  
 Sample : P1352-09  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 C0FP0

Quant Time: Feb 14 01:54:50 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN020724.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 13 22:50:45 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.869  | 152  | 182851   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.669 | 136  | 731185   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.510 | 164  | 358831   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.263 | 188  | 640511   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.463 | 240  | 519378   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.833 | 264  | 606572   | 20.000 | ng/u1 | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.323  | 96  | 23946   | 4.307  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.728  | 84  | 318227  | 21.465 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.034  | 99  | 443089  | 25.907 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.211  | 67  | 293880  | 25.746 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.399  | 132 | 343582  | 27.722 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.581  | 113 | 345554  | 27.091 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.034  | 128 | 164753  | 29.032 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.758  | 143 | 158583  | 32.086 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.287 | 165 | 278745  | 28.573 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.810 | 131 | 443535  | 26.650 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.934 | 166 | 822999  | 30.806 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.204 | 160 | 970877  | 30.602 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.722 | 143 | 108309  | 21.360 | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.510 | 176 | 695930  | 30.621 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.628 | 200 | 72746   | 23.309 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.363 | 188 | 940296  | 31.324 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.663 | 212 | 1050457 | 33.703 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.680 | 264 | 990705  | 32.640 | ng/u1 | 0.00 |

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

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

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