

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN111522\  
 Data File : BN022703.D  
 Acq On : 17 Nov 2022 01:33  
 Operator : CG/JU  
 Sample : N5577-04  
 Misc :  
 ALS Vial : 63 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 C9816

Quant Time: Nov 17 03:23:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN111522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 15 22:53:43 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.869  | 152  | 85676    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.699 | 136  | 372126   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.557 | 164  | 236776   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.316 | 188  | 474342   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.522 | 240  | 432755   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.957 | 264  | 384047   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.193  | 96   | 9447     | 3.765  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.634  | 84   | 40603    | 5.721  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.040  | 99   | 55103    | 6.810  | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.205  | 67   | 144633   | 29.811 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.399  | 132  | 135425   | 22.853 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.599  | 113  | 98303    | 15.388 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.052  | 128  | 89605    | 31.108 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.781  | 143  | 92660    | 28.830 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.322 | 165  | 150087   | 27.392 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.846 | 131  | 108541   | 13.005 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.969 | 166  | 578857   | 33.611 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.245 | 160  | 608036   | 32.042 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.787 | 143  | 20608    | 6.002  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.551 | 176  | 482803   | 34.181 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.687 | 200  | 84990    | 28.189 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.416 | 188  | 776055   | 37.363 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.722 | 212  | 849462   | 37.877 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.798 | 264  | 722109   | 37.404 | ng/u1 | 0.00     |

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

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

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