

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN111522\  
 Data File : BN022700.D  
 Acq On : 16 Nov 2022 23:43  
 Operator : CG/JU  
 Sample : N5577-01  
 Misc :  
 ALS Vial : 60 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 C9811

Quant Time: Nov 17 01:15:44 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  | 81755    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.698 | 136  | 348439   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.557 | 164  | 219938   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.316 | 188  | 459513   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.527 | 240  | 430424   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.957 | 264  | 410315   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.193  | 96   | 10046    | 4.195  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.634  | 84   | 45671    | 6.744  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.046  | 99   | 40456    | 5.239  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.205  | 67   | 131028   | 28.302 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.399  | 132  | 115038   | 20.343 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.599  | 113  | 75032    | 12.309 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.051  | 128  | 79913    | 29.630 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.781  | 143  | 79684    | 26.478 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.322 | 165  | 128450   | 25.037 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.845 | 131  | 124548   | 15.938 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.969 | 166  | 519077   | 32.448 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.251 | 160  | 541751   | 30.735 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.792 | 143  | 15251    | 4.782  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.551 | 176  | 430157   | 32.785 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.686 | 200  | 71045    | 24.324 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.416 | 188  | 699867   | 34.783 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.721 | 212  | 784278   | 35.160 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.798 | 264  | 736201   | 35.693 | ng/u1 | 0.00     |

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

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

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