

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN030822\  
 Data File : BN018875.D  
 Acq On : 08 Mar 2022 17:43  
 Operator : CG/JU  
 Sample : N1733-26  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DBRU4

Quant Time: Mar 09 01:56:07 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN030822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 08 14:15:20 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.905  | 152  | 114127   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.716 | 136  | 537218   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.540 | 164  | 377121   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.281 | 188  | 837737   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.457 | 240  | 935896   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.845 | 264  | 1052139  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.270  | 96   | 19557    | 6.854  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.693  | 84   | 225606   | 25.839 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.058  | 99   | 376682   | 33.807 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.228  | 67   | 242711   | 36.165 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.434  | 132  | 285897   | 35.804 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.610  | 113  | 315457   | 34.818 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.063  | 128  | 144818   | 35.094 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.793  | 143  | 166865   | 36.805 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.328 | 165  | 298879   | 34.463 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.840 | 131  | 458834   | 35.136 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.945 | 166  | 1075955  | 37.683 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.234 | 160  | 1309583  | 36.847 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.716 | 143  | 179675   | 31.073 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.528 | 176  | 952565   | 38.380 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.639 | 200  | 146922   | 28.299 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.375 | 188  | 1543243  | 38.514 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.669 | 212  | 1869552  | 36.913 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.698 | 264  | 2115542  | 38.580 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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