

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061122\  
 Data File : BN020103.D  
 Acq On : 10 Jun 2022 16:43  
 Operator : CG/JU  
 Sample : N3250-15  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 EXYT0

Quant Time: Jun 10 22:46:06 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN060622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 06 15:14:59 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.822  | 152  | 137125   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.610 | 136  | 651312   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.445 | 164  | 431009   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.186 | 188  | 907464   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.374 | 240  | 825239   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.704 | 264  | 754516   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.281  | 96   | 13753    | 4.344  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.693  | 84   | 73127    | 8.451  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.987  | 99   | 71019    | 6.259  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.152  | 67   | 244026   | 34.422 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.352  | 132  | 244556   | 26.526 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.522  | 113  | 155368   | 15.827 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.969  | 128  | 168539   | 34.014 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.693  | 143  | 178104   | 33.869 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.228 | 165  | 274720   | 29.925 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.740 | 131  | 370030   | 24.890 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.857 | 166  | 1095272  | 34.925 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.139 | 160  | 1259116  | 34.621 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.639 | 143  | 29294    | 4.661  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.439 | 176  | 927206   | 35.729 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.557 | 200  | 164113   | 31.220 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.286 | 188  | 1538469  | 38.846 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.580 | 212  | 1730718  | 38.686 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.557 | 264  | 1468924  | 39.790 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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