

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN112122\  
 Data File : BN022826.D  
 Acq On : 22 Nov 2022 16:00  
 Operator : CG/JU  
 Sample : N5593-08  
 Misc :  
 ALS Vial : 48 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 H0AD7

Quant Time: Nov 22 17:38:09 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN111522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 22 05:15:51 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.846  | 152  | 59387    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.663 | 136  | 285317   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.516 | 164  | 184929   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.269 | 188  | 383773   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.474 | 240  | 371810   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.863 | 264  | 379392   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.175  | 96   | 8303     | 4.773  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.616  | 84   | 51274    | 10.423 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.016  | 99   | 46932    | 8.367  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.181  | 67   | 127885   | 38.027 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.369  | 132  | 116056   | 28.253 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.569  | 113  | 84913    | 19.176 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.028  | 128  | 79367    | 35.938 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.751  | 143  | 83508    | 33.888 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.287 | 165  | 133510   | 31.781 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.816 | 131  | 149296   | 23.331 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.934 | 166  | 556438   | 41.368 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.210 | 160  | 584883   | 39.464 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.739 | 143  | 19621    | 7.317  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.510 | 176  | 452077   | 40.979 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.645 | 200  | 81375    | 33.359 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.369 | 188  | 666874   | 39.684 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.674 | 212  | 809940   | 42.035 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.715 | 264  | 786052   | 41.216 | ng/u1 | 0.01     |

Target Compounds Qvalue

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

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