

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN020623\  
 Data File : BN023954.D  
 Acq On : 06 Feb 2023 15:58  
 Operator : CG/JU  
 Sample : PB150667BL  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK667

Quant Time: Feb 07 03:00:44 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN020623.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 07 02:28:26 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.840  | 152  | 148553   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.651 | 136  | 600069   | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.492 | 164  | 392048   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.245 | 188  | 832828   | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.439 | 240  | 721589   | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.816 | 264  | 741739   | 20.000 ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.228  | 96   | 18663    | 5.080 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.652  | 84   | 237833   | 23.338 ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.011  | 99   | 305194   | 24.032 ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.175  | 67   | 202259   | 25.239 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.369  | 132  | 250167   | 25.328 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.557  | 113  | 242628   | 24.456 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.010  | 128  | 119742   | 24.996 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.734  | 143  | 134413   | 24.989 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.269 | 165  | 241076   | 24.050 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.793 | 131  | 343072   | 25.324 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.910 | 166  | 775889   | 25.606 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.187 | 160  | 839670   | 25.013 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.704 | 143  | 116371   | 22.326 ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.487 | 176  | 654372   | 25.648 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.610 | 200  | 107203   | 18.631 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.339 | 188  | 980108   | 25.834 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.639 | 212  | 1123507  | 25.847 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.668 | 264  | 996709   | 26.265 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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