

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN112421\  
 Data File : BN017570.D  
 Acq On : 23 Nov 2021 17:21  
 Operator : CG/JU  
 Sample : PB140905BL  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK905

Quant Time: Nov 24 00:58:13 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN112221.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 22 16:16:36 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.846  | 152  | 240266   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.640 | 136  | 1048894  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.475 | 164  | 650004   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.216 | 188  | 1174400  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.410 | 240  | 813609   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.774 | 264  | 772172   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.305  | 96   | 38030    | 5.657  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.705  | 84   | 504348   | 28.589 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.005  | 99   | 671629   | 30.706 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.169  | 67   | 431464   | 31.478 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.375  | 132  | 521767   | 31.504 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.546  | 113  | 537646   | 31.249 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.993  | 128  | 265409   | 33.400 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.716  | 143  | 276201   | 35.131 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.257 | 165  | 486187   | 31.411 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.769 | 131  | 751189   | 33.798 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.887 | 166  | 1585851  | 33.556 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.169 | 160  | 1972155  | 32.502 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.663 | 143  | 216029   | 27.013 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.469 | 176  | 1264180  | 31.714 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.581 | 200  | 167343   | 27.434 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.316 | 188  | 1867588  | 34.214 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.616 | 212  | 1811903  | 35.081 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.621 | 264  | 1366423  | 33.659 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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