

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN033124\  
 Data File : BN030557.D  
 Acq On : 30 Mar 2024 20:02  
 Operator : MA/JU  
 Sample : PB159886BL  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK886

Quant Time: Mar 31 13:10:51 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN032824.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Mar 31 12:24:07 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.698  | 152  | 232073   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.486 | 136  | 1010352  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.339 | 164  | 655853   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.086 | 188  | 1388646  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.327 | 240  | 1298874  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.268 | 264  | 1472958  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.199  | 96   | 34505    | 6.971  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.599  | 84   | 341809   | 23.031 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.863  | 99   | 572580   | 30.499 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.040  | 67   | 365934   | 31.558 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.228  | 132  | 468573   | 32.901 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.398  | 113  | 456618   | 29.021 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.851  | 128  | 231711   | 33.060 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.569  | 143  | 217552   | 33.931 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.104 | 165  | 429153   | 30.574 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.622 | 131  | 205718   | 8.902  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.757 | 166  | 1485580  | 33.926 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.033 | 160  | 1675159  | 33.273 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.539 | 143  | 243794   | 29.107 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.339 | 176  | 1262536  | 32.949 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.451 | 200  | 148870   | 24.288 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.186 | 188  | 1980066  | 33.094 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.486 | 212  | 2327729  | 33.333 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.074 | 264  | 2356248  | 34.717 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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