

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN042523\  
 Data File : BN025123.D  
 Acq On : 25 Apr 2023 10:31  
 Operator : CG/JU  
 Sample : PB152285BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK285

Quant Time: Apr 26 00:44:56 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN041023.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Apr 25 08:52:18 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.910  | 152  | 307295   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.722 | 136  | 1347237  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.557 | 164  | 863488   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.304 | 188  | 1865940  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.498 | 240  | 1667424  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.939 | 264  | 1709660  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.293  | 96   | 50651    | 6.273  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.717  | 84   | 700043   | 28.955 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.075  | 99   | 897960   | 29.972 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.240  | 67   | 644046   | 32.140 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.440  | 132  | 700755   | 31.828 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.622  | 113  | 727154   | 30.573 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.081  | 128  | 347394   | 31.847 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.804  | 143  | 386458   | 32.708 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.340 | 165  | 662035   | 31.348 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.863 | 131  | 1079759  | 33.690 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.969 | 166  | 2245331  | 33.543 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.251 | 160  | 2496952  | 32.551 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.757 | 143  | 389541   | 28.389 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.551 | 176  | 1914370  | 33.872 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.675 | 200  | 280473   | 23.097 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.404 | 188  | 3191954  | 36.423 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.704 | 212  | 3565299  | 36.822 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.780 | 264  | 3273129  | 36.615 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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