

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
 Data File : BN034347.D  
 Acq On : 02 Oct 2024 22:50  
 Operator : RC/JU  
 Sample : PB163513BL  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK513

Quant Time: Oct 03 12:46:20 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 30 14:14:38 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.387  | 152  | 261184   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.140 | 136  | 1039657  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.039 | 164  | 623368   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 16.798 | 188  | 1207724  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.027 | 240  | 1167747  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.739 | 264  | 1415215  | 20.000 | ng/ul | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.005  | 96  | 33646   | 5.257  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.387  | 84  | 422752  | 22.221 | ng/ul | 0.00 |
| 7) Phenol-d5                  | 6.587  | 99  | 539651  | 23.281 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 6.740  | 67  | 325903  | 22.660 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4         | 6.928  | 132 | 484743  | 26.567 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8         | 8.099  | 113 | 437291  | 22.927 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5           | 8.528  | 128 | 228289  | 27.549 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.240  | 143 | 245759  | 28.841 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 9.775  | 165 | 421310  | 26.253 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.287 | 131 | 597135  | 24.578 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6      | 13.463 | 166 | 1222131 | 26.234 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8         | 13.722 | 160 | 1446802 | 27.606 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.269 | 143 | 164676  | 17.224 | ng/ul | 0.00 |
| 60) Fluorene-d10              | 15.045 | 176 | 1054419 | 26.706 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.175 | 200 | 149962  | 20.825 | ng/ul | 0.00 |
| 73) Anthracene-d10            | 16.892 | 188 | 1618063 | 28.497 | ng/ul | 0.00 |
| 81) Pyrene-d10                | 19.204 | 212 | 1869216 | 28.729 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.557 | 264 | 2087360 | 28.713 | ng/ul | 0.00 |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

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

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