

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091525\  
 Data File : BN037720.D  
 Acq On : 15 Sep 2025 14:13  
 Operator : RC/JU  
 Sample : PB169502BL  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SBLK502

Quant Time: Sep 15 15:28:14 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091225.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 12 15:43:05 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.852  | 152  | 241380   | 20.000 | ng/u1 | 0.02     |
| 20) Naphthalene-d8            | 10.657 | 136  | 914866   | 20.000 | ng/u1 | 0.01     |
| 38) Acenaphthene-d10          | 14.498 | 164  | 512998   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.245 | 188  | 907144   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.492 | 240  | 842910   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.586 | 264  | 918649   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.270  | 96   | 35009    | 6.143  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.676  | 84   | 477654   | 30.142 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.999  | 99   | 558491   | 31.012 | ng/u1 | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.169  | 67   | 350619   | 31.576 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.375  | 132  | 477057   | 32.100 | ng/u1 | 0.01     |
| 15) 4-Methylphenol-d8         | 8.546  | 113  | 435922   | 30.750 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.005  | 128  | 223831   | 32.442 | ng/u1 | 0.01     |
| 24) 2-Nitrophenol-d4          | 9.728  | 143  | 187373   | 33.194 | ng/u1 | 0.01     |
| 28) 2,4-Dichlorophenol-d3     | 10.269 | 165  | 401695   | 31.163 | ng/u1 | 0.01     |
| 31) 4-Chloroaniline-d4        | 10.781 | 131  | 611383   | 31.269 | ng/u1 | 0.01     |
| 46) Dimethylphthalate-d6      | 13.910 | 166  | 1126253  | 32.881 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.193 | 160  | 1390246  | 32.481 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 159007   | 24.771 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.493 | 176  | 993097   | 32.416 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.598 | 200  | 76869    | 18.795 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.345 | 188  | 1458405  | 32.335 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.633 | 212  | 1570801  | 35.650 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.380 | 264  | 1598418  | 34.462 | ng/u1 | 0.01     |

Target Compounds Qvalue

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

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