

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG091525\  
 Data File : BG064370.D  
 Acq On : 15 Sep 2025 11:25  
 Operator : RC/JU  
 Sample : PB169520BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK520

Quant Time: Sep 15 12:53:06 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091125.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 12 11:06:20 2025  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.848  | 152  | 18351    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.627 | 136  | 80138    | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.470 | 164  | 62411    | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.208 | 188  | 165750   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.438 | 240  | 186552   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.429 | 264  | 206151   | 20.000 | ng/u1 | -0.02    |

|                               |        |     |        |        |       |       |
|-------------------------------|--------|-----|--------|--------|-------|-------|
| System Monitoring Compounds   |        |     |        |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.312  | 96  | 2683   | 7.680  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.717  | 84  | 33642  | 33.008 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.014  | 99  | 46776  | 31.184 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.172  | 67  | 24765  | 31.638 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.378  | 132 | 39510  | 32.934 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.547  | 113 | 41791  | 30.622 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.994  | 128 | 20708  | 32.932 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.716  | 143 | 23418  | 30.531 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.251 | 165 | 43404  | 30.432 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.762 | 131 | 57866  | 30.012 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.876 | 166 | 187359 | 33.684 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.158 | 160 | 172106 | 32.025 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.663 | 143 | 24655  | 28.290 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.463 | 176 | 147682 | 32.725 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.580 | 200 | 30141  | 25.494 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.307 | 188 | 266725 | 32.259 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.593 | 212 | 333606 | 31.748 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 24.229 | 264 | 345839 | 32.550 | ng/u1 | -0.01 |

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

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

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