

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG091625\  
 Data File : BG064383.D  
 Acq On : 16 Sep 2025 13:16  
 Operator : RC/JU  
 Sample : PB169689BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK689

Quant Time: Sep 16 14:38:33 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 | 7.842  | 152  | 16867    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.627 | 136  | 68817    | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10      | 14.469 | 164  | 54597    | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.207 | 188  | 137166   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.438 | 240  | 161320   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.434 | 264  | 177689   | 20.000 | ng/u1 | -0.01    |

|                               |        |     |        |        |       |       |
|-------------------------------|--------|-----|--------|--------|-------|-------|
| System Monitoring Compounds   |        |     |        |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.312  | 96  | 2681   | 8.350  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.717  | 84  | 39063  | 41.699 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.013  | 99  | 53517  | 38.817 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.172  | 67  | 25652  | 35.655 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.378  | 132 | 46340  | 42.026 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.547  | 113 | 47456  | 37.832 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.993  | 128 | 20407  | 37.792 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.716  | 143 | 24634  | 37.400 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.251 | 165 | 48287  | 39.425 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.762 | 131 | 59261  | 35.791 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.876 | 166 | 183947 | 37.804 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.158 | 160 | 176488 | 37.540 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.669 | 143 | 24850  | 32.595 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.462 | 176 | 149767 | 37.937 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.574 | 200 | 27074  | 27.671 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.307 | 188 | 266878 | 39.004 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.593 | 212 | 347217 | 38.212 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 24.228 | 264 | 366650 | 40.036 | ng/u1 | -0.01 |

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

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

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