

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070225\  
 Data File : BP025067.D  
 Acq On : 02 Jul 2025 13:38  
 Operator : RC/JU  
 Sample : PB168468BL  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK468

Quant Time: Jul 02 14:57:56 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP063025.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 01 11:17:33 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.443  | 152  | 356393   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.184 | 136  | 1379398  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.096 | 164  | 897982   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.919 | 188  | 1899673  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.354 | 240  | 2055293  | 20.000 | ng/u1 | 0.01     |
| 88) Perylene-d12              | 24.483 | 264  | 2227295  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.090  | 96   | 51675    | 6.490  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.473  | 84   | 777204   | 34.476 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.631  | 99   | 895404   | 32.636 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.790  | 67   | 494952   | 31.338 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 6.984  | 132  | 751207   | 33.838 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.131  | 113  | 722495   | 32.511 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.572  | 128  | 297089   | 32.084 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.278  | 143  | 257169   | 31.861 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.808  | 165  | 677785   | 33.045 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.325 | 131  | 1189724  | 37.668 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.501 | 166  | 2137541  | 33.136 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 13.772 | 160  | 2406923  | 31.863 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.296 | 143  | 308020   | 25.279 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.113 | 176  | 1798166  | 32.469 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.225 | 200  | 174840   | 21.182 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.013 | 188  | 2923996  | 32.016 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.419 | 212  | 3504694  | 33.247 | ng/u1 | 0.02     |
| 92) Benzo(a)pyrene-d12        | 24.271 | 264  | 3675991  | 32.815 | ng/u1 | 0.00     |

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

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

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