

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070125\  
 Data File : BP025057.D  
 Acq On : 01 Jul 2025 15:50  
 Operator : RC/JU  
 Sample : PB168432BL  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK432

Quant Time: Jul 01 16:39:14 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 | 7.437  | 152  | 403559   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.184 | 136  | 1561806  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10      | 14.095 | 164  | 1000847  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 16.925 | 188  | 2127513  | 20.000 | ng/u1 | 0.01     |
| 79) Chrysene-d12          | 21.354 | 240  | 2254342  | 20.000 | ng/u1 | 0.01     |
| 88) Perylene-d12          | 24.489 | 264  | 2420110  | 20.000 | ng/u1 | 0.01     |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.090  | 96  | 47862   | 5.308  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.472  | 84  | 833533  | 32.653 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 6.631  | 99  | 996595  | 32.079 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 6.790  | 67  | 545901  | 30.525 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 6.984  | 132 | 842486  | 33.514 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.131  | 113 | 797799  | 31.704 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.566  | 128 | 342895  | 32.706 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.278  | 143 | 299354  | 32.755 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 9.807  | 165 | 750550  | 32.319 | ng/u1 | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.313 | 131 | 1274685 | 35.644 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.507 | 166 | 2302412 | 32.024 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 13.778 | 160 | 2567652 | 30.497 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.289 | 143 | 341998  | 25.183 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.113 | 176 | 1952047 | 31.625 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.231 | 200 | 182562  | 19.749 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.019 | 188 | 3149754 | 30.795 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.419 | 212 | 3672974 | 31.767 | ng/u1 | 0.02  |
| 92) Benzo(a)pyrene-d12        | 24.265 | 264 | 3720050 | 30.563 | ng/u1 | 0.00  |

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

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

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