

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM012825\  
 Data File : BM049473.D  
 Acq On : 28 Jan 2025 18:04  
 Operator : RC/JU  
 Sample : PB166287BL  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK287

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 01/29/2025  
 Supervised By :mohammad ahmed 02/05/2025

Quant Time: Jan 29 07:38:09 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM012825.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jan 28 14:45:18 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.510  | 152  | 152805   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.269 | 136  | 536151   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.151 | 164  | 328572   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.904 | 188  | 685874   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.133 | 240  | 661398   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.915 | 264  | 720748   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.116  | 96   | 23653    | 5.735  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.510  | 84   | 342481   | 28.264 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.710  | 99   | 346143   | 27.915 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.863  | 67   | 237272   | 27.983 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.051  | 132  | 272474   | 28.397 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.228  | 113  | 246921   | 27.179 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.657  | 128  | 110595   | 26.963 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.369  | 143  | 120927   | 27.029 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.904  | 165  | 245456   | 26.885 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.422 | 131  | 298563   | 25.494 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.575 | 166  | 699157   | 29.021 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.839 | 160  | 807669   | 28.534 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.386 | 143  | 70952m   | 17.953 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.151 | 176  | 627356   | 29.612 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.280 | 200  | 79053    | 18.291 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.004 | 188  | 994381   | 29.033 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.309 | 212  | 1182786  | 30.591 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.727 | 264  | 1109240  | 28.855 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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