

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120123\  
 Data File : BP018517.D  
 Acq On : 02 Dec 2023 16:02  
 Operator : MA/JU  
 Sample : PB157321BL  
 Misc :  
 ALS Vial : 47 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK321

Quant Time: Dec 02 22:17:34 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 01 21:56:43 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.863  | 152  | 401208   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.657 | 136  | 1696149  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.486 | 164  | 1041333  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.233 | 188  | 2135516  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.316 | 240  | 2024084  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 23.692 | 264  | 2412607  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.269  | 96  | 68860   | 5.849  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.687  | 84  | 853271  | 27.057 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.028  | 99  | 1155222 | 28.692 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.199  | 67  | 706805  | 29.619 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.393  | 132 | 891263  | 28.369 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.569  | 113 | 911617  | 27.978 | ng/u1 | -0.01 |
| 21) Nitrobenzene-d5           | 9.022  | 128 | 435590  | 28.854 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.740  | 143 | 462921  | 29.577 | ng/u1 | -0.01 |
| 28) 2,4-Dichlorophenol-d3     | 10.275 | 165 | 873897  | 27.306 | ng/u1 | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.798 | 131 | 994740  | 23.783 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.904 | 166 | 2645435 | 28.464 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.181 | 160 | 3024891 | 29.557 | ng/u1 | -0.01 |
| 54) 4-Nitrophenol-d4          | 14.686 | 143 | 433864  | 28.439 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.481 | 176 | 2231598 | 28.605 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.598 | 200 | 307273  | 23.997 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.333 | 188 | 3383841 | 30.141 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.569 | 212 | 3897713 | 24.738 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.539 | 264 | 4211991 | 29.942 | ng/u1 | -0.01 |

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

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

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