

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112723\  
 Data File : BP018416.D  
 Acq On : 28 Nov 2023 00:34  
 Operator : MA/JU  
 Sample : PB157281BL  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK281

Quant Time: Nov 28 03:18:05 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 28 02:41:10 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.857  | 152  | 502008   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.657 | 136  | 1989131  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.492 | 164  | 1136747  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.245 | 188  | 2075776  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.386 | 240  | 1954342  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.921 | 264  | 2295277  | 20.000 | ng/u1 | 0.01     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.269  | 96  | 87625   | 6.538  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.687  | 84  | 1150913 | 31.246 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.022  | 99  | 1413997 | 30.820 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.193  | 67  | 824606  | 30.057 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.387  | 132 | 1208491 | 31.373 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.569  | 113 | 1132297 | 30.582 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.022  | 128 | 570523  | 35.312 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.740  | 143 | 575228  | 38.551 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.275 | 165 | 1131877 | 31.442 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.798 | 131 | 1603734 | 33.031 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.910 | 166 | 3227032 | 31.904 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.187 | 160 | 3718564 | 32.265 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.687 | 143 | 398681  | 28.552 | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.486 | 176 | 2658022 | 31.418 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.598 | 200 | 261278  | 29.568 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.345 | 188 | 3798424 | 34.236 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.592 | 212 | 4263739 | 27.412 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.751 | 264 | 4715720 | 35.438 | ng/u1 | 0.01 |

Target Compounds Qvalue

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

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