

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018078.D
 Acq On : 11 Nov 2023 22:33
 Operator : MA/JU
 Sample : PB156719BL
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK719

Quant Time: Nov 12 08:47:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.846 | 152 | 70406 | 20.000 | ng/u1 | 0.00 |
| 20) Naphthalene-d8 | 10.669 | 136 | 282933 | 20.000 | ng/u1 | 0.00 |
| 38) Acenaphthene-d10 | 14.522 | 164 | 178605 | 20.000 | ng/u1 | 0.00 |
| 64) Phenanthrene-d10 | 17.304 | 188 | 403980 | 20.000 | ng/u1 | 0.00 |
| 79) Chrysene-d12 | 21.427 | 240 | 399563 | 20.000 | ng/u1 | 0.00 |
| 88) Perylene-d12 | 23.933 | 264 | 463772 | 20.000 | ng/u1 | 0.00 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.228 | 96 | 10233 | 5.127 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.664 | 84 | 129448 | 24.875 | ng/u1 | 0.00 |
| 7) Phenol-d5 | 7.016 | 99 | 152458 | 22.430 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.199 | 67 | 99108 | 24.811 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4 | 7.369 | 132 | 127467 | 23.681 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8 | 8.575 | 113 | 127307 | 22.856 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5 | 9.052 | 128 | 61517 | 24.133 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4 | 9.769 | 143 | 71035 | 24.598 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.299 | 165 | 115566 | 22.453 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4 | 10.851 | 131 | 160664 | 22.762 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6 | 13.945 | 166 | 402917 | 24.450 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8 | 14.222 | 160 | 427389 | 23.811 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4 | 14.798 | 143 | 43517 | 17.672 | ng/u1 | 0.02 |
| 60) Fluorene-d10 | 15.522 | 176 | 329093 | 24.188 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.681 | 200 | 58363 | 19.984 | ng/u1 | 0.00 |
| 73) Anthracene-d10 | 17.404 | 188 | 544989 | 24.944 | ng/u1 | 0.00 |
| 81) Pyrene-d10 | 19.657 | 212 | 668496 | 25.047 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12 | 23.762 | 264 | 690140 | 25.592 | ng/u1 | -0.01 |

Target Compounds Qvalue

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

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