

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112223\  
 Data File : BP018316.D  
 Acq On : 23 Nov 2023 14:59  
 Operator : MA/JU  
 Sample : PB157189BL  
 Misc :  
 ALS Vial : 43 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK189

Quant Time: Nov 24 00:10:51 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 22 14:09:55 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.875  | 152  | 451832   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.669 | 136  | 1868142  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10      | 14.504 | 164  | 1173522  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.251 | 188  | 2155005  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12          | 21.357 | 240  | 1402779  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.780 | 264  | 1752270  | 20.000 | ng/u1 | 0.00     |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.275  | 96  | 77612   | 6.434  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.687  | 84  | 1034217 | 31.196 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.034  | 99  | 1269570 | 30.745 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.205  | 67  | 766058  | 31.024 | ng/u1 | -0.01 |
| 11) 2-Chlorophenol-d4         | 7.399  | 132 | 1058153 | 30.521 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.575  | 113 | 1035337 | 31.069 | ng/u1 | -0.01 |
| 21) Nitrobenzene-d5           | 9.034  | 128 | 514609  | 33.915 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.752  | 143 | 502949  | 35.890 | ng/u1 | -0.01 |
| 28) 2,4-Dichlorophenol-d3     | 10.287 | 165 | 1029138 | 30.440 | ng/u1 | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.810 | 131 | 1542012 | 33.817 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.922 | 166 | 3270334 | 31.319 | ng/u1 | -0.01 |
| 49) Acenaphthylene-d8         | 14.198 | 160 | 3753184 | 31.544 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.687 | 143 | 376739  | 26.135 | ng/u1 | -0.01 |
| 60) Fluorene-d10              | 15.498 | 176 | 2689737 | 30.796 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.604 | 200 | 240615  | 26.229 | ng/u1 | -0.01 |
| 73) Anthracene-d10            | 17.351 | 188 | 3669464 | 31.857 | ng/u1 | -0.01 |
| 81) Pyrene-d10                | 19.592 | 212 | 3388540 | 30.351 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.621 | 264 | 3264519 | 32.135 | ng/u1 | 0.00  |

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

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

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