

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP052622\  
 Data File : BP010567.D  
 Acq On : 26 May 2022 13:54  
 Operator : CG/JU  
 Sample : PB145098BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK098

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 05/27/2022  
 Supervised By :mohammad ahmed 05/31/2022

Quant Time: May 27 02:13:37 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 24 05:50:40 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.881  | 152  | 129878   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.681 | 136  | 568689   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.510 | 164  | 398728   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.263 | 188  | 891287   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.351 | 240  | 760058   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.763 | 264  | 603997   | 20.000 | ng/u1 | -0.01    |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.334  | 96   | 18081m   | 5.934  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.752  | 84   | 270651   | 31.283 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.046  | 99   | 332601   | 30.511 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.210  | 67   | 243157   | 33.200 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.410  | 132  | 269088   | 32.260 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.587  | 113  | 284100   | 30.667 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.040  | 128  | 148435m  | 33.725 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.763  | 143  | 157541   | 33.776 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.304 | 165  | 265388   | 30.091 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.816 | 131  | 325430   | 24.816 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.922 | 166  | 1036905  | 35.296 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.204 | 160  | 1142821  | 33.726 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.716 | 143  | 125500m  | 24.001 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.504 | 176  | 878880   | 34.388 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.634 | 200  | 138862   | 26.640 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.363 | 188  | 1426747  | 35.352 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.598 | 212  | 1670093  | 41.376 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.610 | 264  | 1088824  | 35.901 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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