

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032223\  
 Data File : BP014264.D  
 Acq On : 23 Mar 2023 19:17  
 Operator : CG/JU  
 Sample : PB151603BL  
 Misc : GCMS CONFIRMATION  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK603

Quant Time: Mar 24 00:53:25 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030723.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Mar 23 04:56:06 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.034  | 152  | 155341   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.857 | 136  | 686280   | 20.000 | ng/ul | # 0.00   |
| 38) Acenaphthene-d10          | 14.681 | 164  | 511137   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.439 | 188  | 1190108  | 20.000 | ng/ul | # 0.00   |
| 79) Chrysene-d12              | 21.521 | 240  | 1150781  | 20.000 | ng/ul | # 0.00   |
| 88) Perylene-d12              | 24.039 | 264  | 1059219  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.405  | 96   | 25080    | 6.100  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.834  | 84   | 347009   | 30.483 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.193  | 99   | 451013   | 32.900 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.357  | 67   | 291777   | 36.496 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.563  | 132  | 337335   | 32.299 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.746  | 113  | 374883   | 33.479 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.210  | 128  | 177184   | 32.358 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.934  | 143  | 208188   | 31.761 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.481 | 165  | 359360   | 29.137 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.998 | 131  | 525492   | 32.196 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.087 | 166  | 1388009  | 31.997 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.375 | 160  | 1455255  | 31.668 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.887 | 143  | 186148   | 24.351 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.675 | 176  | 1168697  | 31.789 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.798 | 200  | 228657   | 25.421 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.539 | 188  | 1959706  | 33.920 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.763 | 212  | 2356062  | 37.121 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.874 | 264  | 2015649  | 35.775 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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