

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP031623\  
 Data File : BP014100.D  
 Acq On : 16 Mar 2023 13:36  
 Operator : CG/JU  
 Sample : PB151453BL  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK453

Quant Time: Mar 17 00:59:30 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030723.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 17 00:57:25 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.063  | 152  | 111883   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.887 | 136  | 472677   | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.704 | 164  | 343753   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.463 | 188  | 822228   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.545 | 240  | 908312   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 24.080 | 264  | 1033746  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.422  | 96   | 15350    | 5.184  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.858  | 84   | 131563   | 16.046 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.216  | 99   | 250641   | 25.386 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.381  | 67   | 159646   | 27.725 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.587  | 132  | 186625   | 24.809 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.775  | 113  | 203246   | 25.201 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.234  | 128  | 92525    | 24.533 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.963  | 143  | 105405   | 23.347 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.504 | 165  | 189422   | 22.299 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.028 | 131  | 285736   | 25.418 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.110 | 166  | 725444   | 24.867 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.398 | 160  | 752806   | 24.359 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.904 | 143  | 106156   | 20.649 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.698 | 176  | 614009   | 24.834 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.816 | 200  | 116120   | 18.686 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.563 | 188  | 1028483  | 25.767 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.786 | 212  | 1305506  | 26.060 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.915 | 264  | 1439901  | 26.186 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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