

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032023\  
 Data File : BP014181.D  
 Acq On : 21 Mar 2023 14:58  
 Operator : CG/JU  
 Sample : PB151552BL  
 Misc :  
 ALS Vial : 51 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK552

Quant Time: Mar 22 01:20:26 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030723.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 21 04:47:07 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.046  | 152  | 192039   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.869 | 136  | 814790   | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.686 | 164  | 575737   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.445 | 188  | 1346352  | 20.000 ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.521 | 240  | 1182468  | 20.000 ng/u1 | # 0.00   |
| 88) Perylene-d12              | 24.039 | 264  | 1016463  | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.411  | 96   | 27098    | 5.331 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.846  | 84   | 380192   | 27.016 ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.205  | 99   | 509451   | 30.062 ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.369  | 67   | 318975   | 32.273 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.575  | 132  | 377951   | 29.272 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.763  | 113  | 417499   | 30.160 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.222  | 128  | 196754   | 30.264 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.951  | 143  | 231331   | 29.726 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.493 | 165  | 400321   | 27.339 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.010 | 131  | 574013   | 29.622 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.092 | 166  | 1426341  | 29.192 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.381 | 160  | 1545468  | 29.858 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.892 | 143  | 214605   | 24.923 ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.681 | 176  | 1209753  | 29.214 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.798 | 200  | 249867   | 24.555 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.545 | 188  | 2004485  | 30.669 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.769 | 212  | 2355285  | 36.115 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.874 | 264  | 1694272  | 31.336 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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