

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032223\  
 Data File : BP014245.D  
 Acq On : 23 Mar 2023 07:52  
 Operator : CG/JU  
 Sample : 01977-15  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EYGA2

Quant Time: Mar 23 08:45:42 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  | 157693   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.857 | 136  | 707816   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.681 | 164  | 517102   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.439 | 188  | 1234820  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.521 | 240  | 1259794  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 24.045 | 264  | 1329763  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.405  | 96   | 25241    | 6.048  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.834  | 84   | 295066   | 25.534 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.193  | 99   | 488132   | 35.077 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.357  | 67   | 313419   | 38.618 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.563  | 132  | 359110   | 33.871 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.751  | 113  | 418886   | 36.851 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.210  | 128  | 198967   | 35.230 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.940  | 143  | 238007   | 35.206 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.481 | 165  | 407881   | 32.065 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.004 | 131  | 554540   | 32.942 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.086 | 166  | 1559352  | 35.533 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.375 | 160  | 1683303  | 36.209 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.892 | 143  | 239479   | 30.966 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.675 | 176  | 1339141  | 36.005 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.798 | 200  | 266933   | 28.602 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.539 | 188  | 2206488  | 36.809 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.769 | 212  | 2678687  | 38.552 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.880 | 264  | 2630920  | 37.195 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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