

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP032223\  
 Data File : BP014258.D  
 Acq On : 23 Mar 2023 15:16  
 Operator : CG/JU  
 Sample : 02010-03  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 CB949

Quant Time: Mar 23 16:35:17 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  | 147073   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.858 | 136  | 675018   | 20.000 | ng/ul | # 0.00   |
| 38) Acenaphthene-d10          | 14.681 | 164  | 492330   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.440 | 188  | 1150504  | 20.000 | ng/ul | # 0.00   |
| 79) Chrysene-d12              | 21.522 | 240  | 1140556  | 20.000 | ng/ul | # 0.00   |
| 88) Perylene-d12              | 24.039 | 264  | 1179639  | 20.000 | ng/ul | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.405  | 96   | 14972    | 3.846  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.846  | 84   | 80240    | 7.445  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.199  | 99   | 80689    | 6.217  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.358  | 67   | 235795   | 31.152 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.564  | 132  | 215082   | 21.751 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.752  | 113  | 172423   | 16.264 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.211  | 128  | 149257   | 27.712 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.940  | 143  | 162903   | 25.267 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.481 | 165  | 315400   | 25.999 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.999 | 131  | 372850   | 23.225 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 14.087 | 166  | 1301451  | 31.148 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.375 | 160  | 1340517  | 30.286 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.910 | 143  | 27142    | 3.686  | ng/ul | 0.02     |
| 60) Fluorene-d10              | 15.675 | 176  | 1094412  | 30.906 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.798 | 200  | 111359   | 12.806 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.540 | 188  | 1868958  | 33.463 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.763 | 212  | 2284738  | 36.320 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.874 | 264  | 2178077  | 34.712 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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