

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122222\  
 Data File : BP013238.D  
 Acq On : 22 Dec 2022 10:31  
 Operator : CG/JU  
 Sample : N6130-01 10X  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Dec 22 21:44:46 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 21 23:38:37 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.946  | 152  | 473341   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 1896098  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.592 | 164  | 1045288  | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.345 | 188  | 1968762  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.433 | 240  | 1975514  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.909 | 264  | 2286042  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.334  | 96   | 24942    | 2.253  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.769  | 84   | 280275   | 8.912  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.104  | 99   | 247243   | 6.413  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 729742   | 31.376 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.475  | 132  | 621615   | 20.481 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.657  | 113  | 413633   | 13.093 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 9.116  | 128  | 423665   | 30.411 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.840  | 143  | 412466   | 26.299 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165  | 712133   | 23.377 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.898 | 131  | 948565   | 22.056 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.998 | 166  | 2471177  | 30.761 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.286 | 160  | 2818394  | 31.927 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.792 | 143  | 47287    | 3.272  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.580 | 176  | 2168177  | 31.902 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.704 | 200  | 253606   | 20.017 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.445 | 188  | 3446983  | 38.809 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.680 | 212  | 4135015  | 36.466 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.751 | 264  | 4511717  | 39.595 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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