

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112223\  
 Data File : BP018282.D  
 Acq On : 22 Nov 2023 19:12  
 Operator : MA/JU  
 Sample : 05233-01  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH376

Quant Time: Nov 22 23:19:16 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 22 14:09:55 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.881  | 152  | 381510   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.681 | 136  | 1659080  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.510 | 164  | 1127934  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.257 | 188  | 2373170  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.362 | 240  | 1583498  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.786 | 264  | 1417518  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.275  | 96   | 44488    | 4.368  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.693  | 84   | 324043   | 11.576 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.034  | 99   | 319286   | 9.157  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.210  | 67   | 809145   | 38.809 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.404  | 132  | 803424   | 27.445 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.581  | 113  | 532942   | 18.941 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.040  | 128  | 528650   | 39.230 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.757  | 143  | 487221   | 39.149 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.292 | 165  | 960436   | 31.988 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.816 | 131  | 1443811  | 35.653 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.927 | 166  | 3813846  | 38.000 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.204 | 160  | 4233480  | 37.019 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.692 | 143  | 95042    | 6.860  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.504 | 176  | 3186867  | 37.963 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.616 | 200  | 316489   | 31.328 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.357 | 188  | 4861205  | 38.324 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.598 | 212  | 5116559  | 40.599 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.627 | 264  | 3301708  | 40.176 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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