

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112223\  
 Data File : BP018311.D  
 Acq On : 23 Nov 2023 12:10  
 Operator : MA/JU  
 Sample : 05411-01  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0QF0

Quant Time: Nov 23 23:51:12 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     | 7.875  | 152  | 351850   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.675 | 136  | 1502379  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.504 | 164  | 998336   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.251 | 188  | 2098874  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.357 | 240  | 1348631  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.780 | 264  | 1257783  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.275  | 96   | 42829    | 4.559  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 223503   | 8.657  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.028  | 99   | 184538   | 5.739  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.205  | 67   | 626018   | 32.557 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.399  | 132  | 589860   | 21.848 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.575  | 113  | 367400   | 14.158 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.034  | 128  | 406053   | 33.275 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.752  | 143  | 337541   | 29.951 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.287 | 165  | 698179   | 25.678 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.810 | 131  | 973374   | 26.544 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.922 | 166  | 3089039  | 34.774 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.198 | 160  | 3243685  | 32.046 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.687 | 143  | 45495    | 3.710  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.498 | 176  | 2676594  | 36.024 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.610 | 200  | 233856   | 26.174 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.351 | 188  | 4467295  | 39.821 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.592 | 212  | 4598671  | 42.844 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.621 | 264  | 3032130  | 41.581 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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