

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012405.D  
 Acq On : 01 Nov 2022 01:51  
 Operator : CG/JU  
 Sample : N5216-03  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHA60

Quant Time: Nov 01 02:18:24 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 01 01:03:58 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.793  | 152  | 249642   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.593 | 136  | 1115210  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.446 | 164  | 730549   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.204 | 188  | 1624563  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.298 | 240  | 1798812  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.686 | 264  | 1839092  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.240  | 96   | 27211    | 4.389  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.670  | 84   | 99825    | 5.605  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 6.969  | 99   | 137376   | 6.135  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 470133   | 35.172 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.322  | 132  | 457038   | 27.951 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.511  | 113  | 277245   | 15.655 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.963  | 128  | 322918   | 44.652 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.681  | 143  | 331485   | 44.161 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.216 | 165  | 542511   | 33.350 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.746 | 131  | 774953   | 31.995 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.857 | 166  | 2037971  | 40.751 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.140 | 160  | 2301991  | 40.279 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 50498    | 5.335  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.440 | 176  | 1800979  | 42.062 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.575 | 200  | 314987   | 37.241 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.304 | 188  | 3076665  | 43.953 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.545 | 212  | 3690139  | 40.860 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.533 | 264  | 3964329  | 44.116 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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