

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102722\  
 Data File : BP012353.D  
 Acq On : 27 Oct 2022 21:15  
 Operator : CG/JU  
 Sample : N5015-16  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C1BG1

Quant Time: Oct 27 23:23:48 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 26 23:40:20 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.804  | 152  | 330775   | 20.000 | ng/u1  | -0.01    |
| 20) Naphthalene-d8            | 10.610 | 136  | 1492157  | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.457 | 164  | 1005390  | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.222 | 188  | 2195930  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.315 | 240  | 2185226  | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.709 | 264  | 1914417  | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.246  | 96   | 63549    | 7.736  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.675  | 84   | 289936   | 12.286 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.981  | 99   | 1317412  | 44.405 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.146  | 67   | 803550   | 45.371 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.340  | 132  | 1024561  | 47.290 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.522  | 113  | 1004293  | 42.800 | ng/u1  | -0.01    |
| 21) Nitrobenzene-d5           | 8.975  | 128  | 541369   | 55.948 | ng/u1  | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.698  | 143  | 598116   | 59.553 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.234 | 165  | 1049039  | 48.197 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.757 | 131  | 1414019  | 43.631 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.875 | 166  | 3342181  | 48.561 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.151 | 160  | 3784407  | 48.116 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 557466   | 42.797 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.457 | 176  | 2890597  | 49.055 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.586 | 200  | 431352   | 37.729 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.322 | 188  | 4465443  | 47.195 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.563 | 212  | 5321521  | 48.504 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.557 | 264  | 4696029  | 50.203 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 8) Phenol                     | 7.004  | 94   | 54451    | 1.765  | ng/u1# | 86       |
| 16) Acetophenone              | 8.640  | 105  | 37976    | 1.042  | ng/u1  | 97       |
| -----                         |        |      |          |        |        |          |

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

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