

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012397.D  
 Acq On : 31 Oct 2022 21:16  
 Operator : CG/JU  
 Sample : N5216-05  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHA83

Quant Time: Oct 31 23:52:43 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Oct 28 06:05:59 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.793  | 152  | 330385   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.593 | 136  | 1470150  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.445 | 164  | 958871   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.210 | 188  | 2104711  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.298 | 240  | 2285888  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.692 | 264  | 2293554  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.240  | 96   | 31446    | 3.832  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.669  | 84   | 117915   | 5.003  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.969  | 99   | 162554   | 5.486  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.134  | 67   | 515806   | 29.158 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.328  | 132  | 492367   | 22.753 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.510  | 113  | 320261   | 13.665 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.963  | 128  | 337804   | 35.433 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.687  | 143  | 343402   | 34.704 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.222 | 165  | 582368   | 27.157 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.746 | 131  | 757414   | 23.721 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.863 | 166  | 2257838  | 34.397 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.139 | 160  | 2478683  | 33.044 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 63728    | 5.130  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.445 | 176  | 1984896  | 35.319 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.575 | 200  | 346004   | 31.575 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.304 | 188  | 3338944  | 36.818 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.551 | 212  | 4020475  | 35.032 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.539 | 264  | 4159288  | 37.115 | ng/u1 | -0.01    |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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