

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012421.D  
 Acq On : 01 Nov 2022 11:33  
 Operator : CG/JU  
 Sample : N5270-14  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHA70

Quant Time: Nov 01 22:56:44 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  | 309689   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.593 | 136  | 1393248  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.445 | 164  | 901002   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.204 | 188  | 1973249  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.298 | 240  | 2108176  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.686 | 264  | 2069549  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.234  | 96   | 38772    | 5.041  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.664  | 84   | 185578   | 8.399  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.963  | 99   | 255969   | 9.215  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67   | 692707   | 41.775 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.322  | 132  | 644953   | 31.796 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.505  | 113  | 476255   | 21.679 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.963  | 128  | 467730   | 51.769 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.681  | 143  | 457405   | 48.776 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.216 | 165  | 748218   | 36.816 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.740 | 131  | 1080978  | 35.723 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.857 | 166  | 2980648  | 48.326 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.134 | 160  | 3354609  | 47.593 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.669 | 143  | 94250    | 8.074  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.439 | 176  | 2594409  | 49.129 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.575 | 200  | 431232   | 41.975 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.304 | 188  | 4224453  | 49.686 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.545 | 212  | 4977091  | 47.023 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.539 | 264  | 5089323  | 50.329 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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