

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121022\  
 Data File : BP013057.D  
 Acq On : 11 Dec 2022 00:58  
 Operator : CG/JU  
 Sample : N5927-21  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EXZA7

Quant Time: Dec 12 03:30:02 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 12 01:32:54 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.969  | 152  | 615498   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.781 | 136  | 2691340  | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.610 | 164  | 1830335  | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.363 | 188  | 4164213  | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.451 | 240  | 3850569  | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.939 | 264  | 3850669  | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.352  | 96   | 70982    | 4.931 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 357400   | 8.740 ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.122  | 99   | 392688   | 7.833 ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.293  | 67   | 1177690  | 38.941 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.499  | 132  | 1178957  | 29.873 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.675  | 113  | 780604   | 19.003 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.140  | 128  | 779520   | 39.421 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.863  | 143  | 880471   | 39.551 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.399 | 165  | 1469409  | 33.982 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.922 | 131  | 1827748  | 29.942 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.016 | 166  | 5618369  | 39.940 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.304 | 160  | 6080839  | 39.340 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.798 | 143  | 152173   | 6.014 ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.604 | 176  | 4813856  | 40.450 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.722 | 200  | 943901   | 35.223 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.463 | 188  | 7815021  | 41.599 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.698 | 212  | 9170167  | 41.490 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.780 | 264  | 8121516  | 42.314 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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