

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122222\  
 Data File : BP013242.D  
 Acq On : 22 Dec 2022 12:48  
 Operator : CG/JU  
 Sample : N6015-01  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DBXT5

Quant Time: Dec 22 21:46:29 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 21 23:38:37 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.946  | 152  | 572234   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 2440890  | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.587 | 164  | 1542935  | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.345 | 188  | 3203826  | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.433 | 240  | 2131808  | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.910 | 264  | 2332442  | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.334  | 96   | 56691    | 4.236 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.764  | 84   | 352536   | 9.273 ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.105  | 99   | 1172298  | 25.151 ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 763413   | 27.151 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.475  | 132  | 952688   | 25.965 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.657  | 113  | 866870   | 22.698 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.116  | 128  | 480795   | 26.809 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.840  | 143  | 530764   | 26.288 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 1023105  | 26.089 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.898 | 131  | 968746   | 17.498 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.998 | 166  | 3096800  | 26.115 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.281 | 160  | 3578889  | 27.466 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.787 | 143  | 297582   | 13.951 ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 2767413  | 27.586 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.704 | 200  | 323407   | 15.686 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.445 | 188  | 3970390  | 27.469 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.674 | 212  | 4067988  | 33.245 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.745 | 264  | 3221423  | 27.709 ng/u1 | 0.00     |

Target Compounds Qvalue

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

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