

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020123\  
 Data File : BP013698.D  
 Acq On : 01 Feb 2023 17:51  
 Operator : CG/JU  
 Sample : 01277-03  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EX9H3

Quant Time: Feb 01 18:27:27 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP020123.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Feb 01 14:19:27 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.140  | 152  | 289464   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.963 | 136  | 1257829  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.775 | 164  | 895070   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.528 | 188  | 2063414  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.604 | 240  | 1935465  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.180 | 264  | 1745241  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.481  | 96   | 35954    | 5.266  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.917  | 84   | 347900   | 17.759 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.281  | 99   | 296932   | 11.672 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.452  | 67   | 569084   | 38.140 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.664  | 132  | 602040   | 32.283 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.840  | 113  | 539561   | 26.604 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.310  | 128  | 318217   | 39.658 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.040 | 143  | 374983   | 40.100 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.575 | 165  | 801187   | 36.638 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.099 | 131  | 614609   | 21.281 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.175 | 166  | 2727432  | 38.657 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.469 | 160  | 3016287  | 38.694 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.951 | 143  | 70354    | 5.916  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.763 | 176  | 2379523  | 39.127 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.875 | 200  | 325025   | 28.109 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.628 | 188  | 3747935  | 38.967 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.845 | 212  | 4496465  | 41.654 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.016 | 264  | 3600564  | 40.414 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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