

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081822\  
 Data File : BP011489.D  
 Acq On : 18 Aug 2022 19:14  
 Operator : CG/JU  
 Sample : N4221-17  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0KN9

Quant Time: Aug 18 23:19:07 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 17 01:58:37 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.558  | 152  | 76438    | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.346 | 136  | 312539   | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.228 | 164  | 167833   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.998 | 188  | 331539   | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.127 | 240  | 324739   | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.421 | 264  | 362149   | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.070  | 96   | 10583    | 4.296 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.487  | 84   | 37448    | 5.938 ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.740  | 99   | 33205    | 4.727 ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.905  | 67   | 120464   | 26.397 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.087  | 132  | 109327   | 21.161 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.275  | 113  | 64204    | 11.420 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.722  | 128  | 66511    | 29.560 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.440  | 143  | 65906    | 28.781 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.969  | 165  | 103538   | 24.495 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.499 | 131  | 155355   | 22.079 ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.657 | 166  | 367223   | 30.936 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.916 | 160  | 395994   | 30.015 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.469 | 143  | 5953     | 2.603 ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.234 | 176  | 307902   | 30.832 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.369 | 200  | 44555    | 24.871 ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.098 | 188  | 505287   | 34.412 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.363 | 212  | 571189   | 34.765 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.274 | 264  | 601863   | 34.165 ng/u1 | 0.00     |

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

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

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