

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121123\  
 Data File : BP018786.D  
 Acq On : 13 Dec 2023 01:56  
 Operator : MA/JU  
 Sample : 05646-09  
 Misc :  
 ALS Vial : 65 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AY5

Quant Time: Dec 13 02:26:00 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 11 22:02:13 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min)  |
|-------------------------------|--------|------|----------|--------|-------|-----------|
| Internal Standards            |        |      |          |        |       |           |
| 1) 1,4-Dichlorobenzene-d4     | 7.840  | 152  | 171725   | 20.000 | ng/u1 | 0.00      |
| 20) Naphthalene-d8            | 10.628 | 136  | 647205   | 20.000 | ng/u1 | 0.00      |
| 38) Acenaphthene-d10          | 14.463 | 164  | 422948   | 20.000 | ng/u1 | 0.00      |
| 64) Phenanthrene-d10          | 17.216 | 188  | 1009907  | 20.000 | ng/u1 | # 0.00    |
| 79) Chrysene-d12              | 21.304 | 240  | 1161918  | 20.000 | ng/u1 | # 0.00    |
| 88) Perylene-d12              | 23.663 | 264  | 1387036  | 20.000 | ng/u1 | -0.02     |
| System Monitoring Compounds   |        |      |          |        |       |           |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 24356    | 4.833  | ng/uL | 0.00      |
| 4) Pyridine-d5                | 3.681  | 84   | 296467   | 21.964 | ng/u1 | 0.00      |
| 7) Phenol-d5                  | 7.011  | 99   | 413087   | 23.970 | ng/u1 | 0.00      |
| 9) Bis-(2-Chloroethyl)eth...  | 7.175  | 67   | 237674   | 23.270 | ng/u1 | 0.00      |
| 11) 2-Chlorophenol-d4         | 7.369  | 132  | 328788   | 24.451 | ng/u1 | 0.00      |
| 15) 4-Methylphenol-d8         | 8.546  | 113  | 348157   | 24.964 | ng/u1 | 0.00      |
| 21) Nitrobenzene-d5           | 8.999  | 128  | 160598   | 27.880 | ng/u1 | 0.00      |
| 24) 2-Nitrophenol-d4          | 9.716  | 143  | 180652   | 30.249 | ng/u1 | 0.00      |
| 28) 2,4-Dichlorophenol-d3     | 10.252 | 165  | 361812   | 29.628 | ng/u1 | 0.00      |
| 31) 4-Chloroaniline-d4        | 10.775 | 131  | 450345   | 28.218 | ng/u1 | 0.00      |
| 46) Dimethylphthalate-d6      | 13.881 | 166  | 1241754  | 32.895 | ng/u1 | 0.00      |
| 49) Acenaphthylene-d8         | 14.157 | 160  | 1310848  | 31.536 | ng/u1 | 0.00      |
| 54) 4-Nitrophenol-d4          | 14.669 | 143  | 194965   | 31.465 | ng/u1 | 0.00      |
| 60) Fluorene-d10              | 15.457 | 176  | 1079971  | 34.084 | ng/u1 | 0.00      |
| 65) 4,6-Dinitro-2-methylph... | 15.581 | 200  | 160953   | 26.580 | ng/u1 | 0.00      |
| 73) Anthracene-d10            | 17.316 | 188  | 1824559  | 34.366 | ng/u1 | 0.00      |
| 81) Pyrene-d10                | 19.551 | 212  | 2361150  | 26.106 | ng/u1 | 0.00      |
| 92) Benzo(a)pyrene-d12        | 23.516 | 264  | 2764445  | 34.182 | ng/u1 | -0.01     |
| Target Compounds              |        |      |          |        |       |           |
| 12) 2-Chlorophenol            | 7.399  | 128  | 15703    | 1.154  | ng/u1 | Qvalue 92 |

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

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