

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121922\  
 Data File : BP013214.D  
 Acq On : 21 Dec 2022 18:37  
 Operator : CG/JU  
 Sample : N6032-01  
 Misc :  
 ALS Vial : 96 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 COL84

Quant Time: Dec 21 21:54:41 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 19 21:43:11 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.952  | 152  | 555320   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 2382521  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.593 | 164  | 1539928  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.345 | 188  | 3137279  | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.433 | 240  | 2099140  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.910 | 264  | 2398476  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.334  | 96   | 57946    | 4.462  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 52903    | 1.434  | ng/u1 | 0.02     |
| 7) Phenol-d5                  | 7.111  | 99   | 233699   | 5.167  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 531715   | 19.487 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.481  | 132  | 514628   | 14.453 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 368750   | 9.950  | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.116  | 128  | 325871   | 18.616 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.840  | 143  | 328276   | 16.658 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165  | 610245   | 15.942 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.904 | 131  | 307233   | 5.685  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.998 | 166  | 2173361  | 18.364 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.287 | 160  | 2251375  | 17.312 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.793 | 143  | 51185    | 2.404  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.587 | 176  | 1912397  | 19.100 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.704 | 200  | 195590   | 9.688  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.445 | 188  | 3225139  | 22.787 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.680 | 212  | 3108811  | 25.801 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.751 | 264  | 2699739  | 22.583 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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