

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092523\  
 Data File : BP017349.D  
 Acq On : 26 Sep 2023 07:45  
 Operator : MA/JU  
 Sample : 04472-09  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EZYE9

Quant Time: Sep 26 08:16:05 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 26 04:52:29 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.928  | 152  | 200659   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.746 | 136  | 852247   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.587 | 164  | 519186   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.357 | 188  | 1166084  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.463 | 240  | 1071309  | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.980 | 264  | 1195624  | 20.000 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 15961    | 3.267  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.746  | 84   | 26687    | 1.954  | ng/u1  | 0.02     |
| 7) Phenol-d5                  | 7.081  | 99   | 316891   | 19.229 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.269  | 67   | 205778   | 20.690 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.446  | 132  | 259864   | 19.881 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.634  | 113  | 159178   | 12.424 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.122  | 128  | 129013   | 20.939 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.834  | 143  | 138462   | 20.697 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.363 | 165  | 247720   | 19.599 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.916 | 131  | 230387   | 12.583 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.998 | 166  | 839559   | 22.573 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.287 | 160  | 899389   | 21.025 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.816 | 143  | 80250    | 12.385 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 751530   | 23.471 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.716 | 200  | 50280    | 7.650  | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.457 | 188  | 1135707  | 21.960 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.698 | 212  | 1494242  | 25.792 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.810 | 264  | 1389742  | 24.142 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        | Qvalue |          |
| 8) Phenol                     | 7.110  | 94   | 21770    | 1.313  | ng/u1# | 93       |
| 16) Acetophenone              | 8.775  | 105  | 131807   | 6.657  | ng/u1  | 98       |
| 72) Phenanthrene              | 17.398 | 178  | 582355   | 9.535  | ng/u1  | 99       |
| 74) Anthracene                | 17.492 | 178  | 97215    | 1.571  | ng/u1  | 99       |
| 80) Fluoranthene              | 19.369 | 202  | 900314   | 13.325 | ng/u1  | 99       |
| 82) Pyrene                    | 19.727 | 202  | 794468   | 11.061 | ng/u1  | 100      |
| 85) Benzo(a)anthracene        | 21.445 | 228  | 436427   | 6.036  | ng/u1  | 98       |
| 87) Chrysene                  | 21.498 | 228  | 413151   | 6.072  | ng/u1  | 97       |
| 90) Benzo(b)fluoranthene      | 23.192 | 252  | 581230   | 8.245  | ng/u1# | 95       |
| 91) Benzo(k)fluoranthene      | 23.239 | 252  | 210463   | 2.959  | ng/u1# | 95       |
| 93) Benzo(a)pyrene            | 23.863 | 252  | 461941   | 6.849  | ng/u1  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.627 | 276  | 338172   | 4.000  | ng/u1  | 97       |
| 95) Dibenzo(a,h)anthracene    | 26.639 | 278  | 83736    | 1.189  | ng/u1# | 80       |
| 96) Benzo(g,h,i)perylene      | 27.462 | 276  | 316785   | 4.647  | ng/u1  | 96       |
| -----                         |        |      |          |        |        |          |

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

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