

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121922\  
 Data File : BP013188.D  
 Acq On : 21 Dec 2022 03:52  
 Operator : CG/JU  
 Sample : N5984-01  
 Misc :  
 ALS Vial : 70 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0LC3

Quant Time: Dec 21 05:36:45 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  | 558992   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.763 | 136  | 2381465  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.593 | 164  | 1530983  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.351 | 188  | 3114713  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.433 | 240  | 2291792  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.916 | 264  | 2664263  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.334  | 96   | 73218    | 5.600  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.764  | 84   | 1076293  | 28.981 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.111  | 99   | 1419073  | 31.167 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 891416   | 32.455 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.481  | 132  | 1138185  | 31.755 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.664  | 113  | 1139277  | 30.538 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.122  | 128  | 573027   | 32.750 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.846  | 143  | 656577   | 33.331 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165  | 1192008  | 31.154 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.905 | 131  | 1537939  | 28.472 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.999 | 166  | 3568267  | 30.326 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.287 | 160  | 4084141  | 31.589 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.793 | 143  | 460560   | 21.760 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.587 | 176  | 3084371  | 30.985 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.704 | 200  | 395350   | 19.724 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.451 | 188  | 4492446  | 31.970 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.681 | 212  | 4561819  | 34.678 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.757 | 264  | 4197330  | 31.607 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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