

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP113022\  
 Data File : BP012907.D  
 Acq On : 01 Dec 2022 08:02  
 Operator : CG/JU  
 Sample : N5770-05  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH1G2

Quant Time: Dec 01 08:31:23 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 29 22:53:34 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.010  | 152  | 611450   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.822 | 136  | 2834267  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.645 | 164  | 2002302  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.398 | 188  | 4373630  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.486 | 240  | 3833627  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.992 | 264  | 3263752  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.393  | 96   | 63571    | 4.455  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.822  | 84   | 173624   | 4.066  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.158  | 99   | 403407   | 8.029  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.328  | 67   | 1005017  | 32.461 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.534  | 132  | 1018374  | 26.599 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.710  | 113  | 735617   | 18.171 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.175  | 128  | 657873   | 37.924 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.899  | 143  | 673369   | 36.557 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.434 | 165  | 1241239  | 28.881 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 562907   | 9.565  | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.051 | 166  | 4623615  | 33.548 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.339 | 160  | 5214972  | 32.429 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.822 | 143  | 105399   | 4.289  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.634 | 176  | 4192750  | 34.737 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.745 | 200  | 458649   | 21.941 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.498 | 188  | 6942415  | 36.060 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.727 | 212  | 7794914  | 34.993 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.833 | 264  | 5865826  | 36.590 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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