

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051122\  
 Data File : BP010305.D  
 Acq On : 12 May 2022 02:43  
 Operator : CG/JU  
 Sample : N2707-02  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EXL85

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022  
 Supervised By :Yogesh Patel 05/17/2022

Quant Time: May 12 05:33:06 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu May 12 01:10:15 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.905  | 152  | 137681   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.705 | 136  | 579508   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.534 | 164  | 412710   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.287 | 188  | 932771   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.369 | 240  | 936821   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.798 | 264  | 906061   | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.346  | 96   | 16235m   | 5.009  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.775  | 84   | 45975    | 5.098  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.069  | 99   | 64236    | 5.023  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.234  | 67   | 231871   | 29.311 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.440  | 132  | 206993   | 21.979 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.616  | 113  | 131042   | 12.465 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.063  | 128  | 141800   | 30.517 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.787  | 143  | 145061   | 27.381 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.328 | 165  | 247038   | 25.619 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.846 | 131  | 210287   | 14.772 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.946 | 166  | 1021289  | 31.062 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.228 | 160  | 1118716  | 28.538 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.740 | 143  | 22223m   | 3.524  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.528 | 176  | 890793   | 32.096 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.651 | 200  | 171548   | 26.857 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.387 | 188  | 1595264  | 34.924 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.616 | 212  | 1920961  | 37.106 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.645 | 264  | 1696796  | 35.116 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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