

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102521\  
 Data File : BP007705.D  
 Acq On : 27 Oct 2021 06:55  
 Operator : CG/JU  
 Sample : M4303-10  
 Misc :  
 ALS Vial : 63 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 P013-TW001-03

Manual Integrations  
 APPROVED

mohammad  
 10/27/2021 11:51:19 AM

Quant Time: Oct 27 07:28:04 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP102521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Oct 25 16:57:16 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.934  | 152  | 201067   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.734 | 136  | 826412   | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.569 | 164  | 551998   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.328 | 188  | 1214432  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.416 | 240  | 1296813  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.863 | 264  | 1360121  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.370  | 96   | 30182    | 6.819  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 193485   | 14.455 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.099  | 99   | 17413    | 1.140  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.258  | 67   | 260565   | 30.730 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.463  | 132  | 12267    | 1.033  | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.640  | 113  | 79988    | 6.689  | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.087  | 128  | 174208   | 30.933 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.816  | 143  | 1011m    | 0.164  | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.357 | 165  | 12984    | 1.037  | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.869 | 131  | 447498   | 27.031 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.981 | 166  | 1122234  | 30.014 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.263 | 160  | 1343471  | 29.315 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.793 | 143  | 320m     | 0.047  | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.563 | 176  | 977992   | 31.026 | ng/u1 | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.687 | 200  | 468m     | 0.072  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.428 | 188  | 1671888  | 33.499 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.657 | 212  | 2019265  | 33.445 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.704 | 264  | 2046448  | 30.628 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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