

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP011421\  
 Data File : BP004560.D  
 Acq On : 14 Jan 2021 16:55  
 Operator : CG/JU  
 Sample : M1004-07  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BG-2

Quant Time: Jan 14 21:51:02 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270E-BP011221.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jan 12 15:19:32 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.810  | 152  | 196072   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.610 | 136  | 719780   | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.463 | 164  | 433418   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.228 | 188  | 927340   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.322 | 240  | 1094444  | 20.00  | ng    | 0.00     |
| 86) Perylene-d12            | 23.680 | 264  | 1223378  | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.405  | 112  | 1214628  | 107.97 | ng    | 0.00     |
| 7) Phenol-d6                | 6.987  | 99   | 1450177  | 94.12  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.969  | 82   | 862859   | 67.77  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.963 | 330  | 653053   | 86.66  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.081 | 172  | 1842439  | 55.26  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.792 | 244  | 2290342  | 38.23  | ng    | 0.00     |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 50) Dimethylphthalate       | 13.916 | 163  | 78828    | 2.19   | ng    | 99       |
| 71) Phenanthrene            | 17.269 | 178  | 315212   | 5.93   | ng    | 99       |
| 75) Fluoranthene            | 19.239 | 202  | 637171   | 9.85   | ng    | 99       |
| 78) Pyrene                  | 19.598 | 202  | 534249   | 7.17   | ng    | 99       |
| 81) Benzo(a)anthracene      | 21.304 | 228  | 319600   | 4.29   | ng    | 98       |
| 83) Chrysene                | 21.357 | 228  | 345324   | 4.87   | ng    | 98       |
| 87) Indeno(1,2,3-cd)pyrene  | 26.104 | 276  | 219802   | 2.33   | ng    | # 94     |
| 88) Benzo(b)fluoranthene    | 22.963 | 252  | 486950   | 6.15   | ng    | 97       |
| 90) Benzo(a)pyrene          | 23.574 | 252  | 312197   | 4.23   | ng    | 97       |
| 92) Benzo(g,h,i)perylene    | 26.845 | 276  | 217820   | 2.84   | ng    | 97       |
| -----                       |        |      |          |        |       |          |

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

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