

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061221\  
 Data File : BF124295.D  
 Acq On : 12 Jun 2021 19:55  
 Operator : JU/CG  
 Sample : M2677-03 10X  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 T2-CON

Quant Time: Jun 14 04:28:00 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF060321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jun 10 01:53:56 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.840  | 152  | 31890    | 20.000 | ng    | -0.01    |
| 21) Naphthalene-d8          | 8.122  | 136  | 106376   | 20.000 | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.881  | 164  | 55508    | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.369 | 188  | 80032    | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 14.016 | 240  | 67195    | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 15.504 | 264  | 42035    | 20.000 | ng    | 0.01     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.475  | 112  | 9630     | 4.546  | ng    | 0.00     |
| 7) Phenol-d6                | 6.498  | 99   | 12776    | 4.487  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.404  | 82   | 11243    | 4.192  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.675 | 330  | 3206     | 4.077  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.198  | 172  | 20338    | 4.611  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.951 | 244  | 20145    | 4.116  | ng    | -0.01    |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 78) Pyrene                  | 12.816 | 202  | 21888    | 3.392  | ng    | 95       |
| 90) Benzo(a)pyrene          | 15.439 | 252  | 8024     | 2.829  | ng    | # 58     |
| 92) Benzo(g,h,i)perylene    | 17.451 | 276  | 10975    | 7.236  | ng    | # 82     |
| -----                       |        |      |          |        |       |          |

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

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