

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF022321\  
 Data File : BF123281.D  
 Acq On : 23 Feb 2021 15:11  
 Operator : JU/CG  
 Sample : PB134772TB  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB134772TB

Quant Time: Feb 23 18:19:46 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF021921.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Feb 19 16:08:53 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.857  | 152  | 114612   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.139  | 136  | 465753   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.898  | 164  | 232233   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.386 | 188  | 440664   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.039 | 240  | 372873   | 20.00  | ng    | 0.00     |
| 86) Perylene-d12            | 15.510 | 264  | 348451   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.487  | 112  | 977621   | 122.78 | ng    | 0.01     |
| 7) Phenol-d6                | 6.493  | 99   | 1230137  | 113.72 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.416  | 82   | 814351   | 80.04  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.692 | 330  | 270258   | 114.29 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.222  | 172  | 1272836  | 78.73  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.986 | 244  | 1396199  | 72.00  | ng    | 0.00     |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

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

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