

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF042021\  
 Data File : BF123647.D  
 Acq On : 20 Apr 2021 17:26  
 Operator : JU/CG  
 Sample : M2070-02  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 JGD-3150-D3550-A

Quant Time: Apr 21 01:20:05 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF041521.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 20 15:49:40 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.851  | 152  | 244953   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.134  | 136  | 959442   | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.892  | 164  | 476127   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.380 | 188  | 922874   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.027 | 240  | 611103   | 20.00  | ng    | 0.00     |
| 86) Perylene-d12            | 15.504 | 264  | 546783   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.487  | 112  | 1831151  | 120.82 | ng    | 0.02     |
| 7) Phenol-d6                | 6.487  | 99   | 2176085  | 103.62 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.416  | 82   | 1917039  | 91.85  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.686 | 330  | 791988   | 146.89 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.216  | 172  | 2854521  | 89.37  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.974 | 244  | 3129955  | 99.54  | ng    | 0.00     |

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

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

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