

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF080321\  
 Data File : BF124927.D  
 Acq On : 3 Aug 2021 21:33  
 Operator : JU/CG  
 Sample : M3241-03  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SPLP-AMENDED-COM-6

Quant Time: Aug 04 05:09:00 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF072221.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 02 11:15:51 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.845  | 152  | 6635     | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 8.128  | 136  | 26144    | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.881  | 164  | 15131    | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.369 | 188  | 26036    | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 14.004 | 240  | 18681    | 20.000  | ng    | -0.02    |
| 86) Perylene-d12            | 15.468 | 264  | 14279    | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.457  | 112  | 23216    | 59.244  | ng    | -0.02    |
| 7) Phenol-d6                | 6.469  | 99   | 18117    | 36.878  | ng    | -0.02    |
| 23) Nitrobenzene-d5         | 7.410  | 82   | 44421    | 101.126 | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol    | 10.675 | 330  | 17924    | 137.968 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.204  | 172  | 90665    | 88.007  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.957 | 244  | 109714   | 100.672 | ng    | 0.00     |

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

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

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