

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091521\  
 Data File : BF125526.D  
 Acq On : 16 Sep 2021 2:59  
 Operator : JU/CG  
 Sample : M3687-27 5X  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 A011145

Quant Time: Sep 16 04:16:03 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF091521.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 15 15:08:32 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min)     |
|-----------------------------|--------|------|----------|--------|-------|--------------|
| -----                       |        |      |          |        |       |              |
| Internal Standards          |        |      |          |        |       |              |
| 1) 1,4-Dichlorobenzene-d4   | 6.869  | 152  | 157250   | 20.000 | ng    | 0.00         |
| 21) Naphthalene-d8          | 8.151  | 136  | 566203   | 20.000 | ng    | # 0.00       |
| 39) Acenaphthene-d10        | 9.910  | 164  | 256047   | 20.000 | ng    | 0.00         |
| 64) Phenanthrene-d10        | 11.398 | 188  | 392348   | 20.000 | ng    | 0.00         |
| 76) Chrysene-d12            | 14.051 | 240  | 274545   | 20.000 | ng    | # 0.00       |
| 86) Perylene-d12            | 15.551 | 264  | 220874   | 20.000 | ng    | 0.02         |
| System Monitoring Compounds |        |      |          |        |       |              |
| 5) 2-Fluorophenol           | 5.487  | 112  | 162003   | 16.369 | ng    | 0.00         |
| 7) Phenol-d6                | 6.498  | 99   | 194329   | 14.774 | ng    | -0.01        |
| 23) Nitrobenzene-d5         | 7.428  | 82   | 101160   | 8.821  | ng    | -0.01        |
| 42) 2,4,6-Tribromophenol    | 10.698 | 330  | 32773    | 10.509 | ng    | 0.00         |
| 45) 2-Fluorobiphenyl        | 9.222  | 172  | 227199   | 13.021 | ng    | -0.01        |
| 79) Terphenyl-d14           | 12.980 | 244  | 180388   | 10.654 | ng    | 0.00         |
| Target Compounds            |        |      |          |        |       |              |
| 32) Benzoic acid            | 7.910  | 122  | 31902    | 5.954  | ng    | Qvalue<br>89 |
| -----                       |        |      |          |        |       |              |

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

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