

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF111021\  
 Data File : BF126099.D  
 Acq On : 10 Nov 2021 10:15  
 Operator : JU/CG  
 Sample : PB140520BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB140520BL

Quant Time: Nov 10 13:33:56 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF110321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 03 17:21:25 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.851  | 152  | 225789   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.134  | 136  | 840729   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.886  | 164  | 508383   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.375 | 188  | 944668   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.010 | 240  | 724873   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.474 | 264  | 554517   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.481  | 112  | 1623038  | 123.443 | ng    | 0.02     |
| 7) Phenol-d6                | 6.487  | 99   | 2045999  | 110.389 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.416  | 82   | 1525728  | 85.585  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.680 | 330  | 762420   | 130.863 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.216  | 172  | 2891945  | 76.872  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.963 | 244  | 3557611  | 79.142  | ng    | 0.00     |

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

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

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