

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF032922\  
 Data File : BF127558.D  
 Acq On : 29 Mar 2022 11:37  
 Operator : CG\JU  
 Sample : PB143673BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB143673BL

Quant Time: Mar 29 15:53:41 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF032122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 29 07:13:31 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.834  | 152  | 112787   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.116  | 136  | 418014   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.869  | 164  | 245483   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.363 | 188  | 450660   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.010 | 240  | 323719   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.492 | 264  | 232933   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.469  | 112  | 865950   | 122.554 | ng    | 0.01     |
| 7) Phenol-d6                | 6.475  | 99   | 1188167  | 123.018 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.404  | 82   | 815774   | 83.420  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.669 | 330  | 319047   | 122.006 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.192  | 172  | 1394350  | 82.942  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.951 | 244  | 1601808  | 81.002  | ng    | 0.00     |

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

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

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