

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF092022\  
 Data File : BF130337.D  
 Acq On : 20 Sep 2022 11:07  
 Operator : CG\JU  
 Sample : PB147679BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB147679BL

Quant Time: Sep 21 05:10:02 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Sep 19 15:06:14 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.657  | 152  | 106766   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 7.939  | 136  | 432741   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.692  | 164  | 235358   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.169 | 188  | 416219   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.804 | 240  | 306467   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.192 | 264  | 229684   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.275  | 112  | 908211   | 147.543 | ng    | 0.02     |
| 7) Phenol-d6                | 6.304  | 99   | 1111367  | 144.295 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.228  | 82   | 722190   | 85.260  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.480 | 330  | 375043   | 166.303 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.022  | 172  | 1391809  | 86.113  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.763 | 244  | 1700314  | 91.904  | ng    | 0.00     |

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

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

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