

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091423\  
 Data File : BF135247.D  
 Acq On : 14 Sep 2023 11:28  
 Operator : CG\JU  
 Sample : PB155361BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB155361BL

Quant Time: Sep 15 01:05:28 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Sep 12 00:19:07 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.763  | 152  | 115872   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.040  | 136  | 445119   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.798  | 164  | 236615   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.292 | 188  | 478649   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.945 | 240  | 264407   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.416 | 264  | 242573   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.404  | 112  | 955181   | 134.075 | ng    | 0.02     |
| 7) Phenol-d6                | 6.410  | 99   | 1192096  | 131.594 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.328  | 82   | 765505   | 93.434  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.598 | 330  | 339100   | 144.213 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.122  | 172  | 1374844  | 87.664  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.892 | 244  | 1560049  | 91.464  | ng    | 0.00     |

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

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

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