

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF071423\  
 Data File : BF134284.D  
 Acq On : 14 Jul 2023 11:22  
 Operator : CG\JU  
 Sample : PB154028BL  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB154028BL

Quant Time: Jul 15 00:22:37 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF062623.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jul 14 22:51:27 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.840  | 152  | 84351    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.122  | 136  | 320267   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.881  | 164  | 166040   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.375 | 188  | 342713   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.039 | 240  | 278472   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.545 | 264  | 282511   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.481  | 112  | 655275   | 129.948 | ng    | 0.01     |
| 7) Phenol-d6                | 6.487  | 99   | 770300   | 124.214 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.410  | 82   | 539936   | 90.879  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.681 | 330  | 291756   | 140.145 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.204  | 172  | 1041637  | 91.209  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.980 | 244  | 1357597  | 78.462  | ng    | 0.00     |

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

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

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