

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101422\  
 Data File : BF130663.D  
 Acq On : 14 Oct 2022 12:35  
 Operator : CG\JU  
 Sample : PB148312BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB148312BL

Quant Time: Oct 17 02:19:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Oct 12 06:52:52 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.581  | 152  | 87535    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 7.863  | 136  | 353350   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.610  | 164  | 192253   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.092 | 188  | 349072   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.721 | 240  | 266203   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.098 | 264  | 209674   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.193  | 112  | 623967   | 128.489 | ng    | 0.00     |
| 7) Phenol-d6                | 6.234  | 99   | 765350   | 125.353 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.151  | 82   | 489721   | 75.659  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.404 | 330  | 250195   | 132.999 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 8.939  | 172  | 1000038  | 78.792  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.680 | 244  | 1175218  | 73.273  | ng    | 0.00     |

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

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

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