

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101922\  
 Data File : BF130716.D  
 Acq On : 19 Oct 2022 17:57  
 Operator : CG\JU  
 Sample : PB148423BL  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB148423BL

Quant Time: Oct 20 02:45:10 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Oct 18 03:15:54 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.569  | 152  | 74473    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 7.851  | 136  | 301454   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.598  | 164  | 163600   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.080 | 188  | 302522   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.710 | 240  | 195935   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 15.086 | 264  | 156189   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.187  | 112  | 851519   | 206.102 | ng    | 0.02     |
| 7) Phenol-d6                | 6.228  | 99   | 860797   | 165.713 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.139  | 82   | 538400   | 97.499  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.392 | 330  | 285591   | 178.403 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 8.928  | 172  | 1089020  | 100.831 | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.668 | 244  | 1298952  | 110.032 | ng    | 0.00     |

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

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

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