

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF041123\  
 Data File : BF132854.D  
 Acq On : 11 Apr 2023 17:28  
 Operator : CG\JU  
 Sample : PB151349BL  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB151349BL

Quant Time: Apr 12 04:05:30 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF041123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 12 03:34:19 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.769  | 152  | 248529   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.051  | 136  | 984664   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.804  | 164  | 498007   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.280 | 188  | 902948   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.916 | 240  | 661241   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.345 | 264  | 599357   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.387  | 112  | 1968422  | 123.044 | ng    | 0.02     |
| 7) Phenol-d6                | 6.398  | 99   | 2484779  | 120.429 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.334  | 82   | 1530120  | 80.411  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.592 | 330  | 583722   | 131.887 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.133  | 172  | 2491095  | 79.131  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.874 | 244  | 2945455  | 77.180  | ng    | 0.00     |

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

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

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