

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF110422\  
 Data File : BF130992.D  
 Acq On : 04 Nov 2022 15:08  
 Operator : CG\JU  
 Sample : N5340-03  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-7-102722

Quant Time: Nov 05 00:37:28 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 04 06:20:24 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.910  | 152  | 93999    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.198  | 136  | 406327   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.963  | 164  | 186554   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.457 | 188  | 281933   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.104 | 240  | 187115   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.615 | 264  | 149863   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.534  | 112  | 326402   | 60.198  | ng    | 0.00     |
| 7) Phenol-d6                | 6.534  | 99   | 252408   | 35.823  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.481  | 82   | 748773   | 111.118 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.757 | 330  | 248747   | 158.396 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.280  | 172  | 1299019  | 105.276 | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.051 | 244  | 1054329  | 103.252 | ng    | 0.00     |

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

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

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