

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF110522\  
 Data File : BF131019.D  
 Acq On : 05 Nov 2022 14:50  
 Operator : CG\JU  
 Sample : N5340-05RE  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MW-9-102722RE

Quant Time: Nov 07 00:48:35 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF102722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Nov 07 00:46:02 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.904  | 152  | 70223    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.186  | 136  | 276200   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.945  | 164  | 124404   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.439 | 188  | 225347   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.086 | 240  | 140474   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.586 | 264  | 122637   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.516  | 112  | 22959    | 5.668   | ng    | 0.00     |
| 7) Phenol-d6                | 6.522  | 99   | 137378   | 26.098  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.469  | 82   | 503955   | 110.021 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.739 | 330  | 8710     | 8.317   | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.269  | 172  | 866798   | 105.342 | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.033 | 244  | 785967   | 102.528 | ng    | 0.00     |

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

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

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