

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF020223\  
 Data File : BF132302.D  
 Acq On : 02 Feb 2023 20:43  
 Operator : CG\JU  
 Sample : 01386-12  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 JPP-14-013123

Quant Time: Feb 03 00:25:58 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF013123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 02 01:30:34 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.066  | 152  | 191393   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.348  | 136  | 729613   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 10.113 | 164  | 382344   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.613 | 188  | 693281   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 14.277 | 240  | 316175   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 15.860 | 264  | 347040   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.701  | 112  | 1312941  | 110.266 | ng    | 0.02     |
| 7) Phenol-d6                | 6.684  | 99   | 1508190  | 102.740 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.631  | 82   | 1079479  | 82.331  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.913 | 330  | 585484   | 148.907 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.431  | 172  | 1911631  | 77.209  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.213 | 244  | 2010140  | 122.984 | ng    | 0.00     |

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

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

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