

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF110922\  
 Data File : BF131095.D  
 Acq On : 09 Nov 2022 19:32  
 Operator : CG\JU  
 Sample : N5437-01  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 596C-3

Quant Time: Nov 10 04:32:29 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF110922.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 10 04:15:17 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.893  | 152  | 95263    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.175  | 136  | 357578   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.934  | 164  | 195620   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.428 | 188  | 348926   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.069 | 240  | 197238   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 15.563 | 264  | 169155   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.504  | 112  | 328105   | 58.573  | ng    | 0.00     |
| 7) Phenol-d6                | 6.510  | 99   | 256781   | 35.063  | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.457  | 82   | 629282   | 79.402  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.728 | 330  | 267483   | 132.307 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.251  | 172  | 1058324  | 75.524  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.016 | 244  | 1066323  | 94.789  | ng    | 0.00     |

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

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

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