

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111522\  
 Data File : BG055633.D  
 Acq On : 15 Nov 2022 22:32  
 Operator : CG/JU  
 Sample : N5431-06  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DISSOLVED-20221176-COMP-11-MODIFIED-ELUTRIATE

Quant Time: Nov 16 01:46:37 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG110922.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 16 01:36:26 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.256  | 152  | 40507    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.082 | 136  | 169333   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.878 | 164  | 123332   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.616 | 188  | 310284   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.917 | 240  | 300396   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.330 | 264  | 309962   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.783  | 112  | 107512   | 50.554  | ng    | 0.00     |
| 7) Phenol-d6                | 7.387  | 99   | 116675   | 39.718  | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 9.425  | 82   | 173043   | 69.277  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.358 | 330  | 155404   | 129.142 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.503 | 172  | 419297   | 51.282  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.207 | 244  | 793702   | 53.070  | ng    | 0.00     |

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

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

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