

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110922\  
 Data File : BG055528.D  
 Acq On : 9 Nov 2022 18:07  
 Operator : CG/JU  
 Sample : N5360-12  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 20221172-COMP-07-MODIFIED-ELUTRIATE

Quant Time: Nov 10 04:30:37 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG110922.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 10 03:39:46 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.273  | 152  | 55934    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.099 | 136  | 231518   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.883 | 164  | 157250   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.627 | 188  | 349109   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.922 | 240  | 326443   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.342 | 264  | 345511   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.794  | 112  | 176241   | 60.016  | ng    | 0.00     |
| 7) Phenol-d6                | 7.398  | 99   | 171794   | 42.351  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.443  | 82   | 305219   | 89.373  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.364 | 330  | 215967   | 140.760 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.514 | 172  | 730178   | 70.041  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.212 | 244  | 603249   | 37.117  | ng    | 0.00     |

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

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

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