

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110322\  
 Data File : BG055444.D  
 Acq On : 3 Nov 2022 19:26  
 Operator : CG/JU  
 Sample : N5306-03  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 20221168-COMP-03-MODIFIED-ELUTRIATE

Quant Time: Nov 04 05:30:43 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG110122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 04 05:28:43 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.235  | 152  | 28487    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.067 | 136  | 115587   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.851 | 164  | 83906    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.583 | 188  | 200241   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.854 | 240  | 200421   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.227 | 264  | 216638   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.761  | 112  | 103161   | 65.034  | ng    | 0.00     |
| 7) Phenol-d6                | 7.383  | 99   | 107787   | 48.999  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.410  | 82   | 172550   | 79.559  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.331 | 330  | 146739   | 128.213 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.476 | 172  | 444979   | 78.630  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.162 | 244  | 800304   | 77.648  | ng    | 0.00     |

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

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

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