

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG080223\  
 Data File : BG058582.D  
 Acq On : 2 Aug 2023 17:28  
 Operator : MA/JU  
 Sample : 03813-04  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PLAUDERVILLE-2

Quant Time: Aug 03 02:44:09 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG072823.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jul 28 22:29:07 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.812  | 152  | 58734    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.615 | 136  | 253143   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.457 | 164  | 172916   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.207 | 188  | 377747   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 21.443 | 240  | 330884   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 24.516 | 264  | 417875   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.385  | 112  | 255720   | 66.547 | ng    | 0.00     |
| 7) Phenol-d6                | 6.984  | 99   | 384093   | 71.833 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.975  | 82   | 216314   | 41.978 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.950 | 330  | 169667   | 59.864 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.076 | 172  | 486675   | 42.178 | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.821 | 244  | 762922   | 43.359 | ng    | 0.00     |

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

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

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