

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG083023\  
 Data File : BG058904.D  
 Acq On : 30 Aug 2023 20:47  
 Operator : MA/JU  
 Sample : 04196-23  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OR-345-COMP-04

Quant Time: Aug 31 01:41:15 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG081823.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Aug 31 01:27:35 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.755  | 152  | 13224    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.557 | 136  | 57391    | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.412 | 164  | 47320    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.161 | 188  | 118162   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.397 | 240  | 98260    | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 24.429 | 264  | 114642   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.334  | 112  | 128296   | 156.126 | ng    | 0.00     |
| 7) Phenol-d6                | 6.932  | 99   | 166755   | 144.637 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.924  | 82   | 124048   | 103.918 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.910 | 330  | 118654   | 159.918 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.031 | 172  | 322583   | 99.151  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.788 | 244  | 657147   | 120.086 | ng    | 0.00     |

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

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

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