

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042023\  
 Data File : BG057150.D  
 Acq On : 20 Apr 2023 17:41  
 Operator : CG/JU  
 Sample : PB152259BL  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB152259BL

Quant Time: Apr 21 06:06:40 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG042023.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Apr 21 05:53:40 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.393  | 152  | 8199     | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.230 | 136  | 33693    | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 15.001 | 164  | 40958    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.739 | 188  | 101662   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 22.051 | 240  | 94758    | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.570 | 264  | 99869    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.902  | 112  | 63516    | 138.746 | ng    | 0.00     |
| 7) Phenol-d6                | 7.523  | 99   | 93184    | 129.125 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.568  | 82   | 66655    | 90.385  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.488 | 330  | 87032    | 122.316 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.633 | 172  | 236150   | 76.442  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.312 | 244  | 547322   | 89.401  | ng    | 0.00     |

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

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

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