

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052422\  
 Data File : BG053679.D  
 Acq On : 24 May 2022 10:47  
 Operator : CG/JU  
 Sample : PB145019BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB145019BL

Quant Time: May 24 15:00:00 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue May 24 05:05:32 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.055  | 152  | 25135    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.869 | 136  | 102078   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.687 | 164  | 75081    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.430 | 188  | 186986   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.713 | 240  | 198909   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 24.973 | 264  | 187965   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.629  | 112  | 200904   | 135.652 | ng    | 0.00     |
| 7) Phenol-d6                | 7.232  | 99   | 285067   | 126.827 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.218  | 82   | 201522   | 84.038  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.179 | 330  | 136587   | 123.345 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.307 | 172  | 451264   | 87.828  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.039 | 244  | 972110   | 91.023  | ng    | 0.00     |

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

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

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