

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG060822\  
 Data File : BG053850.D  
 Acq On : 8 Jun 2022 13:10  
 Operator : CG/JU  
 Sample : PB145354TB  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB145354TB

Quant Time: Jun 09 02:29:30 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG060222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jun 02 23:22:27 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.095  | 152  | 25434    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.897 | 136  | 103512   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.711 | 164  | 74369    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.454 | 188  | 194188   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.726 | 240  | 228851   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 24.993 | 264  | 220407   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.657  | 112  | 193644   | 139.812 | ng    | 0.00     |
| 7) Phenol-d6                | 7.243  | 99   | 247404   | 131.867 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.246  | 82   | 163842   | 103.942 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.197 | 330  | 145591   | 161.278 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.342 | 172  | 445323   | 89.581  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.063 | 244  | 1017724  | 88.471  | ng    | 0.00     |

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

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

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