

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120822\  
 Data File : BG055927.D  
 Acq On : 8 Dec 2022 14:25  
 Operator : CG/JU  
 Sample : PB149440TB  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB149440TB

Quant Time: Dec 09 08:09:42 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG111622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Dec 09 07:57:14 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.141  | 152  | 24303    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.973 | 136  | 107051   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.774 | 164  | 84862    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.518 | 188  | 221153   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.796 | 240  | 219133   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.127 | 264  | 228207   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.673  | 112  | 201954   | 150.202 | ng    | 0.00     |
| 7) Phenol-d6                | 7.301  | 99   | 264404   | 147.727 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.316  | 82   | 170040   | 83.955  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.267 | 330  | 150924   | 126.300 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.400 | 172  | 448888   | 79.245  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.109 | 244  | 966413   | 88.209  | ng    | 0.00     |

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

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

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