

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG012224\  
 Data File : BG060382.D  
 Acq On : 22 Jan 2024 12:55  
 Operator : MA/JU  
 Sample : PB158504TB  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB158504TB

Quant Time: Jan 22 13:27:08 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG010824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jan 09 04:07:12 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.926  | 152  | 183747   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.734 | 136  | 730329   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.565 | 164  | 578574   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.315 | 188  | 1630778  | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.575 | 240  | 1505646  | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 24.742 | 264  | 1643762  | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.505  | 112  | 1505200  | 134.736 | ng    | 0.00     |
| 7) Phenol-d6                | 7.092  | 99   | 2294448  | 138.701 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.089  | 82   | 1376097  | 88.973  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.058 | 330  | 1212743  | 142.466 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.185 | 172  | 3462253  | 83.198  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.935 | 244  | 5581004  | 99.837  | ng    | 0.00     |

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

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

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