

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG020123\  
 Data File : BG056505.D  
 Acq On : 1 Feb 2023 18:24  
 Operator : CG/JU  
 Sample : PB150216BL  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB150216BL

Quant Time: Feb 02 05:04:47 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG012723.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Feb 01 01:15:52 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.276  | 152  | 33724    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.108 | 136  | 130042   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.892 | 164  | 106440   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.630 | 188  | 293353   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.919 | 240  | 291478   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.333 | 264  | 313759   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.797  | 112  | 275858   | 150.475 | ng    | 0.00     |
| 7) Phenol-d6                | 7.418  | 99   | 379714   | 139.793 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.451  | 82   | 203022   | 88.653  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.378 | 330  | 185508   | 131.096 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.523 | 172  | 700998   | 91.308  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.209 | 244  | 1512802  | 91.157  | ng    | 0.00     |

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

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

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