

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG041823\  
 Data File : BG057135.D  
 Acq On : 18 Apr 2023 22:31  
 Operator : CG/JU  
 Sample : PB152179BL  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB152179BL

Quant Time: Apr 19 05:00:46 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG041823.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 19 04:21:14 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.396  | 152  | 13379    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.240 | 136  | 37096    | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 15.005 | 164  | 39765    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.748 | 188  | 98584    | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 22.060 | 240  | 87748    | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.591 | 264  | 93999    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.906  | 112  | 78581    | 100.032 | ng    | 0.00     |
| 7) Phenol-d6                | 7.533  | 99   | 157893   | 135.969 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.571  | 82   | 105216   | 124.909 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.491 | 330  | 90550    | 134.223 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.636 | 172  | 276268   | 91.532  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.316 | 244  | 571944   | 98.609  | ng    | 0.00     |

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

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

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