

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122222\  
 Data File : BG056067.D  
 Acq On : 22 Dec 2022 18:18  
 Operator : CG/JU  
 Sample : PB149663BL  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB149663BL

Quant Time: Dec 23 05:19:51 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG121322.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Dec 23 05:09:54 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.356  | 152  | 7652     | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.188 | 136  | 30404    | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.960 | 164  | 29965    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.698 | 188  | 93225    | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.999 | 240  | 112813   | 20.000  | ng    | #-0.01   |
| 86) Perylene-d12            | 25.483 | 264  | 121209   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.871  | 112  | 50223    | 153.923 | ng    | 0.00     |
| 7) Phenol-d6                | 7.487  | 99   | 63708    | 151.812 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.525  | 82   | 41980    | 96.489  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.441 | 330  | 105264   | 145.685 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.591 | 172  | 167355   | 78.329  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.278 | 244  | 529025   | 82.880  | ng    | 0.00     |

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

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

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