

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG112221\  
 Data File : BG051166.D  
 Acq On : 22 Nov 2021 10:47  
 Operator : CG/JU  
 Sample : PB140911TB  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB140911TB

Quant Time: Nov 22 17:16:21 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG111921.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 19 16:26:22 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.207  | 152  | 24742    | 20.000  | ng    | # 0.00   |
| 21) Naphthalene-d8          | 11.034 | 136  | 108719   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.835 | 164  | 78323    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.585 | 188  | 186678   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.886 | 240  | 181043   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 25.293 | 264  | 176622   | 20.000  | ng    | 0.01     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.746  | 112  | 224544   | 140.292 | ng    | 0.00     |
| 7) Phenol-d6                | 7.361  | 99   | 300331   | 133.226 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.383  | 82   | 202174   | 84.888  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.327 | 330  | 111788   | 134.284 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.460 | 172  | 440513   | 84.869  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.176 | 244  | 838716   | 84.690  | ng    | 0.00     |

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

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

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