

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG102221\  
 Data File : BG050687.D  
 Acq On : 22 Oct 2021 13:12  
 Operator : CG/JU  
 Sample : PB140104BL  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB140104BL

Quant Time: Oct 22 14:05:54 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Oct 14 10:56:22 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.252  | 152  | 37873    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.073 | 136  | 162296   | 20.000  | ng    | #-0.01   |
| 39) Acenaphthene-d10        | 14.868 | 164  | 105358   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.612 | 188  | 232251   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.913 | 240  | 218832   | 20.000  | ng    | #-0.01   |
| 86) Perylene-d12            | 25.327 | 264  | 204709   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.791  | 112  | 353074   | 139.422 | ng    | 0.00     |
| 7) Phenol-d6                | 7.395  | 99   | 460913   | 125.710 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.422  | 82   | 327601   | 83.559  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.355 | 330  | 121379   | 120.784 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.493 | 172  | 594978   | 84.992  | ng    | -0.01    |
| 79) Terphenyl-d14           | 20.203 | 244  | 970500   | 86.429  | ng    | 0.00     |

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

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

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