

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG012623\  
 Data File : BG056441.D  
 Acq On : 26 Jan 2023 13:43  
 Operator : CG/JU  
 Sample : 01283-05  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MH-10

Quant Time: Jan 27 00:58:22 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG122722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jan 25 04:54:56 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.293  | 152  | 33460    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.125 | 136  | 122478   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.909 | 164  | 97950    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.647 | 188  | 280027   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.936 | 240  | 286793   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 25.368 | 264  | 304435   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.820  | 112  | 253230   | 135.129 | ng    | 0.00     |
| 7) Phenol-d6                | 7.436  | 99   | 314191   | 109.173 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.469  | 82   | 189986   | 79.346  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.390 | 330  | 203992   | 164.441 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.534 | 172  | 641380   | 90.422  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.221 | 244  | 1517539  | 94.772  | ng    | 0.00     |

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

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

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