

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011323\  
 Data File : BG056284.D  
 Acq On : 13 Jan 2023 19:19  
 Operator : CG/JU  
 Sample : 01140-01  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MH-5

Quant Time: Jan 16 00:24:12 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG122722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jan 12 03:38:40 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.317  | 152  | 40620    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.149 | 136  | 155354   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.927 | 164  | 130093   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.665 | 188  | 328820   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.960 | 240  | 314632   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.403 | 264  | 335640   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.837  | 112  | 271122   | 119.175 | ng    | 0.01     |
| 7) Phenol-d6                | 7.453  | 99   | 390816   | 111.861 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.492  | 82   | 207821   | 68.427  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.407 | 330  | 171661   | 104.188 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.558 | 172  | 655757   | 69.607  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.238 | 244  | 1220205  | 69.460  | ng    | 0.00     |

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

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

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