

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG041322\  
 Data File : BG053141.D  
 Acq On : 13 Apr 2022 16:45  
 Operator : CG/JU  
 Sample : N2385-02  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 N48963

Quant Time: Apr 14 02:57:22 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG040822.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Apr 08 15:31:24 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.122  | 152  | 34848    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.930 | 136  | 144802   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.743 | 164  | 104795   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.486 | 188  | 226813   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.768 | 240  | 223302   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 25.064 | 264  | 231907   | 20.000  | ng    | #-0.01   |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.684  | 112  | 132126   | 64.993  | ng    | 0.00     |
| 7) Phenol-d6                | 7.276  | 99   | 125902   | 40.160  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.279  | 82   | 319888   | 89.699  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.229 | 330  | 232054   | 139.184 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.368 | 172  | 713480   | 93.881  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.088 | 244  | 1009169  | 76.327  | ng    | 0.00     |

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

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

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