

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG041522\  
 Data File : BG053174.D  
 Acq On : 15 Apr 2022 21:15  
 Operator : CG/JU  
 Sample : N2394-08  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 CS-4-SOIL-PILE

Quant Time: Apr 18 03:18:28 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   | 8.116  | 152  | 27630    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.930 | 136  | 104359   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.737 | 164  | 77497    | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 17.480 | 188  | 203584   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 21.762 | 240  | 241011   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 25.070 | 264  | 239619   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.684  | 112  | 213337   | 132.356 | ng    | 0.00     |
| 7) Phenol-d6                | 7.276  | 99   | 288385   | 116.021 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.279  | 82   | 232008   | 90.268  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.223 | 330  | 177872   | 144.266 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 13.362 | 172  | 511712   | 91.049  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.088 | 244  | 1246001  | 87.314  | ng    | 0.00     |

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

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

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