

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042121\  
 Data File : BG048820.D  
 Acq On : 21 Apr 2021 15:11  
 Operator : CG/JU  
 Sample : M2090-23  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SEDGWICK-COMP-5

Quant Time: Apr 21 16:39:05 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG041421.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 20 17:33:30 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.062  | 152  | 20665    | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.888 | 136  | 75951    | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.719 | 164  | 51381    | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.480 | 188  | 132578   | 20.00  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.775 | 240  | 139414   | 20.00  | ng    | # 0.00   |
| 86) Perylene-d12            | 25.106 | 264  | 130467   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.612  | 112  | 155037   | 125.48 | ng    | 0.00     |
| 7) Phenol-d6                | 7.227  | 99   | 183797   | 105.17 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.231  | 82   | 169416   | 99.86  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.217 | 330  | 105038   | 152.47 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.338 | 172  | 364813   | 102.54 | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.095 | 244  | 746994   | 92.22  | ng    | 0.00     |

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

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

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