

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111120\  
 Data File : BG047297.D  
 Acq On : 11 Nov 2020 23:29  
 Operator : CG/JU  
 Sample : L4680-06  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PAV-B7-110920-10-20

Quant Time: Nov 12 02:10:17 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG110420.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 04 16:19:40 2020  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.008  | 152  | 57671    | 20.00  | ng    | # 0.00   |
| 21) Naphthalene-d8          | 10.817 | 136  | 232247   | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.642 | 164  | 160724   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.391 | 188  | 394232   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.651 | 240  | 405469   | 20.00  | ng    | # 0.00   |
| 86) Perylene-d12            | 24.847 | 264  | 394793   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.576  | 112  | 376822   | 105.91 | ng    | 0.01     |
| 7) Phenol-d6                | 7.174  | 99   | 451624   | 87.69  | ng    | 0.01     |
| 23) Nitrobenzene-d5         | 9.172  | 82   | 388577   | 71.80  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.134 | 330  | 311339   | 115.68 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.267 | 172  | 889124   | 70.95  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.000 | 244  | 1552793  | 67.34  | ng    | 0.00     |

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

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

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