

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111723\  
 Data File : BG059730.D  
 Acq On : 17 Nov 2023 13:47  
 Operator : MA/JU  
 Sample : 05388-04  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 TP-3

Quant Time: Nov 17 14:23:27 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG111623.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 17 01:51:38 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.342  | 152  | 31322    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.186 | 136  | 132412   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.970 | 164  | 86556    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.720 | 188  | 167287   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 22.044 | 240  | 157440   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.622 | 264  | 231234   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.839  | 112  | 233897   | 125.688 | ng    | 0.01     |
| 7) Phenol-d6                | 7.461  | 99   | 287280   | 104.173 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.541  | 82   | 340112   | 98.100  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.451 | 330  | 135403   | 137.873 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.589 | 172  | 541065   | 84.142  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.293 | 244  | 940701   | 88.205  | ng    | 0.00     |

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

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

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