

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG061022\  
 Data File : BG053879.D  
 Acq On : 10 Jun 2022 17:38  
 Operator : CG/JU  
 Sample : N3286-07  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-7D

Quant Time: Jun 11 02:04:20 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG060222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jun 02 23:22:27 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.088  | 152  | 17972    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.897 | 136  | 71466    | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.710 | 164  | 52951    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.454 | 188  | 149332   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.725 | 240  | 182577   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 24.998 | 264  | 182756   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.650  | 112  | 52402    | 53.544  | ng    | 0.00     |
| 7) Phenol-d6                | 7.236  | 99   | 40609    | 30.632  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.246  | 82   | 116472   | 107.023 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.196 | 330  | 126311   | 196.516 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.335 | 172  | 324075   | 91.560  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.063 | 244  | 802494   | 87.442  | ng    | 0.00     |

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

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

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