

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062922\  
 Data File : BG054145.D  
 Acq On : 29 Jun 2022 14:56  
 Operator : CG/JU  
 Sample : N3517-03  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-3-20220627

Quant Time: Jun 29 17:19:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061322.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 27 00:27:54 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.071  | 152  | 24154    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.874 | 136  | 95047    | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.693 | 164  | 71788    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.437 | 188  | 190931   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.714 | 240  | 196661   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 24.969 | 264  | 200046   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.639  | 112  | 72825    | 57.645  | ng    | 0.00     |
| 7) Phenol-d6                | 7.231  | 99   | 58582    | 34.182  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.229  | 82   | 174062   | 102.914 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.179 | 330  | 187042   | 162.166 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.318 | 172  | 494226   | 100.110 | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.046 | 244  | 955927   | 97.900  | ng    | 0.00     |

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

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

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