

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052522\  
 Data File : BG053708.D  
 Acq On : 26 May 2022 00:25  
 Operator : CG/JU  
 Sample : N2897-10  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 P042-SS001-2430-01

Quant Time: May 26 02:53:27 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 25 15:11:41 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.990  | 152  | 26521    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.798 | 136  | 106977   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.622 | 164  | 79892    | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.372 | 188  | 171487   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 21.636 | 240  | 173369   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 24.838 | 264  | 175800   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.552  | 112  | 144560   | 92.507 | ng    | 0.00     |
| 7) Phenol-d6                | 7.162  | 99   | 223468   | 94.226 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.153  | 82   | 159739   | 63.564 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.114 | 330  | 116638   | 98.987 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.248 | 172  | 351784   | 64.344 | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.980 | 244  | 541035   | 58.122 | ng    | 0.00     |

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

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

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