

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG053023\  
 Data File : BG057608.D  
 Acq On : 30 May 2023 16:08  
 Operator : CG/JU  
 Sample : 02955-16  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DUP07

Quant Time: May 31 03:28:27 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG052723.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 31 03:25:39 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.340  | 152  | 17200    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.177 | 136  | 70095    | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.954 | 164  | 51225    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.698 | 188  | 158785   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.998 | 240  | 161270   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.487 | 264  | 175876   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.861  | 112  | 57491    | 58.069  | ng    | 0.00     |
| 7) Phenol-d6                | 7.488  | 99   | 53334    | 34.422  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.521  | 82   | 124145   | 82.792  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.441 | 330  | 117230   | 142.941 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.580 | 172  | 308641   | 86.589  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.265 | 244  | 779669   | 83.745  | ng    | 0.00     |

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

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

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