

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG081222\  
 Data File : BG054618.D  
 Acq On : 12 Aug 2022 11:59  
 Operator : CG/JU  
 Sample : N4046-05  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BEACONPOINT-BACKFILL-03

Quant Time: Aug 13 00:36:06 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Aug 13 00:32:40 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.965  | 152  | 13651    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.779 | 136  | 49873    | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.610 | 164  | 41141    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.360 | 188  | 119276   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.625 | 240  | 154576   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 24.827 | 264  | 160654   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.556  | 112  | 103422   | 158.181 | ng    | 0.00     |
| 7) Phenol-d6                | 7.172  | 99   | 135842   | 151.627 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.134  | 82   | 90889    | 99.243  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.108 | 330  | 123443   | 142.906 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.229 | 172  | 270204   | 92.651  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.968 | 244  | 771727   | 99.042  | ng    | 0.00     |

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

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

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