

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG040622\  
 Data File : BG053063.D  
 Acq On : 6 Apr 2022 20:23  
 Operator : CG/JU  
 Sample : N2254-05  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 WC-14A-1A-4-3

Quant Time: Apr 07 04:11:04 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG031522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 06 14:35:39 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.129  | 152  | 33507    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.942 | 136  | 127616   | 20.000 | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.755 | 164  | 86801    | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.498 | 188  | 172439   | 20.000 | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.792 | 240  | 175593   | 20.000 | ng    | # 0.01   |
| 86) Perylene-d12            | 25.123 | 264  | 184297   | 20.000 | ng    | # 0.04   |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.691  | 112  | 165389   | 85.960 | ng    | 0.00     |
| 7) Phenol-d6                | 7.289  | 99   | 234039   | 80.685 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.286  | 82   | 185782   | 73.066 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.241 | 330  | 101142   | 86.715 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.380 | 172  | 343620   | 58.385 | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.101 | 244  | 533018   | 53.997 | ng    | 0.00     |
| Target Compounds            |        |      |          |        |       |          |
|                             |        |      |          |        |       | Qvalue   |
| 52) Acenaphthene            | 14.820 | 154  | 14132    | 2.895  | ng    | # 72     |
| 58) Fluorene                | 15.801 | 166  | 38936    | 5.991  | ng    | # 88     |
| 71) Phenanthrene            | 17.545 | 178  | 103513   | 11.221 | ng    | # 91     |
| 78) Pyrene                  | 19.913 | 202  | 77542    | 6.366  | ng    | # 86     |
| 81) Benzo(a)anthracene      | 21.769 | 228  | 31242    | 2.657  | ng    | # 33     |
| 83) Chrysene                | 21.834 | 228  | 57057    | 5.162  | ng    | # 61     |
| -----                       |        |      |          |        |       |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG040622\  
Data File : BG053063.D  
Acq On : 6 Apr 2022 20:23  
Operator : CG/JU  
Sample : N2254-05  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
WC-14A-1A-4-3

Quant Time: Apr 07 04:11:04 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG031522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Wed Apr 06 14:35:39 2022  
Response via : Initial Calibration

