

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121922\  
 Data File : BG056018.D  
 Acq On : 19 Dec 2022 18:08  
 Operator : CG/JU  
 Sample : N6089-01  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MH-24

Quant Time: Dec 20 01:07:17 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG121322.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Dec 20 01:01:12 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min)  |
|-----------------------------|--------|------|----------|---------|-------|-----------|
| -----                       |        |      |          |         |       |           |
| Internal Standards          |        |      |          |         |       |           |
| 1) 1,4-Dichlorobenzene-d4   | 8.364  | 152  | 10229    | 20.000  | ng    | 0.00      |
| 21) Naphthalene-d8          | 11.202 | 136  | 35697    | 20.000  | ng    | # 0.00    |
| 39) Acenaphthene-d10        | 14.974 | 164  | 29884    | 20.000  | ng    | 0.00      |
| 64) Phenanthrene-d10        | 17.712 | 188  | 80835    | 20.000  | ng    | 0.00      |
| 76) Chrysene-d12            | 22.019 | 240  | 82998    | 20.000  | ng    | -0.01     |
| 86) Perylene-d12            | 25.509 | 264  | 98375    | 20.000  | ng    | -0.02     |
| System Monitoring Compounds |        |      |          |         |       |           |
| 5) 2-Fluorophenol           | 5.891  | 112  | 79358    | 181.942 | ng    | 0.00      |
| 7) Phenol-d6                | 7.501  | 99   | 93270    | 166.263 | ng    | 0.00      |
| 23) Nitrobenzene-d5         | 9.539  | 82   | 63118    | 123.563 | ng    | 0.00      |
| 42) 2,4,6-Tribromophenol    | 16.455 | 330  | 119585   | 165.954 | ng    | 0.00      |
| 45) 2-Fluorobiphenyl        | 13.605 | 172  | 227214   | 106.634 | ng    | 0.00      |
| 79) Terphenyl-d14           | 20.286 | 244  | 510046   | 108.611 | ng    | 0.00      |
| Target Compounds            |        |      |          |         |       |           |
| 92) Benzo(g,h,i)perylene    | 30.773 | 276  | 17569    | 2.744   | ng    | Qvalue 95 |
| -----                       |        |      |          |         |       |           |

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

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