

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111822\  
 Data File : BG055689.D  
 Acq On : 18 Nov 2022 13:42  
 Operator : CG/JU  
 Sample : N5557-01  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-14

Manual Integrations  
 APPROVED

Reviewed By :Christian Giraldo 11/21/2022  
 Supervised By :Jagrut Upadhyay 11/21/2022

Quant Time: Nov 18 23:55:36 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG111622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 17 02:02:46 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.255  | 152  | 44397    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 11.082 | 136  | 172255   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.877 | 164  | 200389m  | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.621 | 188  | 239568   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 21.916 | 240  | 238362   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12            | 25.330 | 264  | 252823   | 20.000 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.788  | 112  | 182720   | 74.390 | ng    | 0.00     |
| 7) Phenol-d6                | 7.398  | 99   | 192435   | 58.855 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.425  | 82   | 253108   | 77.664 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.364 | 330  | 170482   | 60.417 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.502 | 172  | 572988   | 42.837 | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.206 | 244  | 640098   | 53.711 | ng    | 0.00     |
|                             |        |      |          |        |       |          |
| Target Compounds            |        |      |          |        |       | Qvalue   |
| 20) 3+4-Methylphenols       | 9.013  | 107  | 8397     | 2.313  | ng    | 95       |
| 31) Naphthalene             | 11.134 | 128  | 140618   | 15.101 | ng    | # 97     |
| 37) 2-Methylnaphthalene     | 12.715 | 142  | 29382    | 4.214  | ng    | 97       |
| 38) 1-Methylnaphthalene     | 12.938 | 142  | 418557   | 61.839 | ng    | 99       |
| 51) 2,6-Dinitrotoluene      | 14.548 | 165  | 9115     | 2.791  | ng    | # 20     |
| 58) Fluorene                | 15.917 | 166  | 33317    | 2.214  | ng    | 96       |
| 71) Phenanthrene            | 17.662 | 178  | 35526    | 2.748  | ng    | # 91     |
| 73) Carbazole               | 18.021 | 167  | 210826   | 18.670 | ng    | 98       |
| -----                       |        |      |          |        |       |          |

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

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