

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG021720\  
 Data File : BG044465.D  
 Acq On : 18 Feb 2020 2:49  
 Operator : CG/JU  
 Sample : L1528-03  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 FK-GF-02-048-B

Manual Integrations  
 APPROVED

mohammad  
 2/18/2020 3:17:09 PM

Quant Time: Feb 18 06:48:34 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG013020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Feb 10 16:11:50 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min)       |
|-----------------------------|-------|------|----------|--------|-------|----------------|
| 1) 1,4-Dichlorobenzene-d4   | 7.89  | 152  | 84772    | 20.00  | ng    | -0.01          |
| 21) Naphthalene-d8          | 10.69 | 136  | 344382   | 20.00  | ng    | -0.01          |
| 39) Acenaphthene-d10        | 14.53 | 164  | 206602   | 20.00  | ng    | -0.01          |
| 64) Phenanthrene-d10        | 17.27 | 188  | 440486   | 20.00  | ng    | -0.01          |
| 76) Chrysene-d12            | 21.52 | 240  | 448706   | 20.00  | ng    | -0.02          |
| 87) Perylene-d12            | 24.60 | 264  | 473411   | 20.00  | ng    | -0.03          |
| System Monitoring Compounds |       |      |          |        |       |                |
| 5) 2-Fluorophenol           | 5.47  | 112  | 561171   | 116.83 | ng    | 0.00           |
| 7) Phenol-d6                | 7.06  | 99   | 715873   | 110.98 | ng    | -0.01          |
| 23) Nitrobenzene-d5         | 9.04  | 82   | 440630   | 72.23  | ng    | -0.02          |
| 42) 2,4,6-Tribromophenol    | 16.02 | 330  | 262070   | 108.05 | ng    | -0.01          |
| 45) 2-Fluorobiphenyl        | 13.15 | 172  | 951295   | 77.10  | ng    | -0.01          |
| 79) Terphenyl-d14           | 19.89 | 244  | 1396544  | 64.02  | ng    | -0.01          |
| Target Compounds            |       |      |          |        |       |                |
| 50) Dimethylphthalate       | 13.98 | 163  | 47775    | 3.22   | ng    | Qvalue<br># 97 |
| 71) Phenanthrene            | 17.32 | 178  | 87511    | 4.10   | ng    | 96             |
| 75) Fluoranthene            | 19.33 | 202  | 243369   | 9.86   | ng    | 98             |
| 78) Pyrene                  | 19.68 | 202  | 226630   | 7.86   | ng    | 94             |
| 81) Benzo(a)anthracene      | 21.50 | 228  | 151754   | 5.49   | ng    | 95             |
| 83) Chrysene                | 21.56 | 228  | 142735   | 5.34   | ng    | 96             |
| 86) Indeno(1,2,3-cd)pyrene  | 28.07 | 276  | 103119   | 2.96   | ng    | 96             |
| 88) Benzo(b)fluoranthene    | 23.62 | 252  | 223961   | 8.21   | ng    | 98             |
| 89) Benzo(k)fluoranthene    | 23.68 | 252  | 80201m   | 3.02   | ng    |                |
| 90) Benzo(a)pyrene          | 24.45 | 252  | 149499   | 5.80   | ng    | 98             |
| 92) Benzo(g,h,i)perylene    | 29.16 | 276  | 100655   | 3.94   | ng    | # 93           |

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

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