

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031620\  
 Data File : BG044822.D  
 Acq On : 16 Mar 2020 20:56  
 Operator : CG/JU  
 Sample : L1859-02  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 TP-7-SA

Manual Integrations  
 APPROVED

mohammad  
 3/18/2020 8:44:22 AM

Quant Time: Mar 17 05:12:07 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Mar 16 17:37:15 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min)  |
|-----------------------------|-------|------|----------|--------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4   | 8.05  | 152  | 116449   | 20.00  | ng    | 0.00      |
| 21) Naphthalene-d8          | 10.86 | 136  | 478423   | 20.00  | ng    | 0.00      |
| 39) Acenaphthene-d10        | 14.68 | 164  | 311936   | 20.00  | ng    | 0.00      |
| 64) Phenanthrene-d10        | 17.43 | 188  | 643370   | 20.00  | ng    | 0.00      |
| 76) Chrysene-d12            | 21.70 | 240  | 504264   | 20.00  | ng    | 0.00      |
| 87) Perylene-d12            | 24.91 | 264  | 582244   | 20.00  | ng    | 0.00      |
| System Monitoring Compounds |       |      |          |        |       |           |
| 5) 2-Fluorophenol           | 5.62  | 112  | 602615   | 80.56  | ng    | 0.00      |
| 7) Phenol-d6                | 7.20  | 99   | 877454   | 80.73  | ng    | 0.00      |
| 23) Nitrobenzene-d5         | 9.20  | 82   | 574199   | 52.67  | ng    | 0.00      |
| 42) 2,4,6-Tribromophenol    | 16.17 | 330  | 336034   | 82.27  | ng    | 0.00      |
| 45) 2-Fluorobiphenyl        | 13.31 | 172  | 1182046  | 54.27  | ng    | 0.00      |
| 79) Terphenyl-d14           | 20.04 | 244  | 1409141  | 52.27  | ng    | 0.00      |
| Target Compounds            |       |      |          |        |       |           |
| 50) Dimethylphthalate       | 14.13 | 163  | 93229    | 3.753  | ng    | Qvalue 98 |
| 71) Phenanthrene            | 17.47 | 178  | 304153   | 8.847  | ng    | 98        |
| 72) Anthracene              | 17.56 | 178  | 68999    | 2.035  | ng    | 95        |
| 75) Fluoranthene            | 19.48 | 202  | 505498   | 12.349 | ng    | 97        |
| 78) Pyrene                  | 19.84 | 202  | 482876   | 13.052 | ng    | 98        |
| 81) Benzo(a)anthracene      | 21.68 | 228  | 233657   | 6.435  | ng    | 96        |
| 83) Chrysene                | 21.74 | 228  | 209435   | 6.038  | ng    | 96        |
| 86) Indeno(1,2,3-cd)pyrene  | 28.54 | 276  | 149778   | 3.294  | ng    | # 94      |
| 88) Benzo(b)fluoranthene    | 23.89 | 252  | 276902   | 7.204  | ng    | 99        |
| 89) Benzo(k)fluoranthene    | 23.95 | 252  | 79114m   | 2.175  | ng    |           |
| 90) Benzo(a)pyrene          | 24.75 | 252  | 210608   | 5.856  | ng    | 98        |
| 92) Benzo(a,h,i)perylene    | 29.68 | 276  | 152837   | 4.263  | ng    | 97        |

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

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