

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\  
 Data File : BF094166.D  
 Acq On : 4 Apr 2017 13:41  
 Operator : SJ/MA  
 Sample : I2510-09DL 10X  
 Misc :  
 ALS Vial : 48 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 E2-M10-ESWDL

Manual Integrations  
 APPROVED

mohammad  
 4/4/2017 5:35:48 PM

Quant Time: Apr 04 14:05:19 2017  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Mar 31 14:38:29 2017  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 5.89  | 152  | 159550   | 20.00 | ng    | -0.02    |
| 21) Naphthalene-d8          | 7.39  | 136  | 585959   | 20.00 | ng    | -0.02    |
| 38) Acenaphthene-d10        | 9.53  | 164  | 227180   | 20.00 | ng    | -0.02    |
| 63) Phenanthrene-d10        | 11.36 | 188  | 356447   | 20.00 | ng    | -0.02    |
| 75) Chrysene-d12            | 14.62 | 240  | 306572   | 20.00 | ng    | -0.01    |
| 86) Perylene-d12            | 16.24 | 264  | 212713   | 20.00 | ng    | 0.00     |
| System Monitoring Compounds |       |      |          |       |       |          |
| 5) 2-Fluorophenol           | 4.42  | 112  | 82711    | 8.76  | ng    | -0.02    |
| 7) Phenol-d6                | 5.49  | 99   | 107611   | 9.21  | ng    | -0.02    |
| 23) Nitrobenzene-d5         | 6.53  | 82   | 49711    | 4.84  | ng    | -0.03    |
| 41) 2,4,6-Tribromophenol    | 10.50 | 330  | 12829    | 7.15  | ng    | -0.02    |
| 44) 2-Fluorobiphenyl        | 8.71  | 172  | 126321   | 8.48  | ng    | -0.02    |
| 78) Terphenyl-d14           | 13.33 | 244  | 80067    | 5.85  | ng    | -0.02    |
| Target Compounds            |       |      |          |       |       | Qvalue   |
| 48) Acenaphthylene          | 9.36  | 152  | 114360   | 4.80  | ng    | 99       |
| 70) Phenanthrene            | 11.39 | 178  | 108273   | 5.46  | ng    | 98       |
| 71) Anthracene              | 11.45 | 178  | 92247    | 4.52  | ng    | 99       |
| 74) Fluoranthene            | 12.85 | 202  | 486534   | 23.96 | ng    | 98       |
| 77) Pyrene                  | 13.13 | 202  | 433059   | 19.46 | ng    | 97       |
| 80) Benzo(a)anthracene      | 14.60 | 228  | 214897m  | 12.02 | ng    |          |
| 82) Chrysene                | 14.64 | 228  | 248579   | 14.98 | ng    | 98       |
| 85) Indeno(1,2,3-cd)pyrene  | 17.59 | 276  | 95525    | 6.65  | ng    | 94       |
| 87) Benzo(b)fluoranthene    | 15.84 | 252  | 292985m  | 22.47 | ng    |          |
| 88) Benzo(k)fluoranthene    | 15.86 | 252  | 94079m   | 7.37  | ng    |          |
| 89) Benzo(a)pyrene          | 16.19 | 252  | 127474   | 10.62 | ng    | 98       |
| 90) Dibenzo(a,h)anthracene  | 17.60 | 278  | 25371    | 2.50  | ng    | # 86     |
| 91) Benzo(a,h,i)perylene    | 17.99 | 276  | 79318    | 7.74  | ng    | 98       |

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\  
Data File : BF094166.D  
Acq On : 4 Apr 2017 13:41  
Operator : SJ/MA  
Sample : I2510-09DL 10X  
Misc :  
ALS Vial : 48 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
E2-M10-ESWDL

Manual Integrations  
APPROVED

mohammad  
4/4/2017 5:35:48 PM

Quant Time: Apr 04 14:05:19 2017  
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Fri Mar 31 14:38:29 2017  
Response via : Initial Calibration

