

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101715\  
 Data File : BF082337.D  
 Acq On : 16 Oct 2015 20:41  
 Operator : UM/IZ  
 Sample : G4046-02 10X  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 GOODLUCK58

Manual Integrations  
 APPROVED

MMdadoda  
 10/19/2015 5:32:26 PM

Quant Time: Oct 19 02:11:07 2015  
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Oct 12 01:46:21 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 6.93  | 152  | 52994m   | 20.00 | ng    | 0.09     |
| 21) Naphthalene-d8        | 8.21  | 136  | 243928m  | 20.00 | ng    | 0.08     |
| 38) Acenaphthene-d10      | 9.95  | 164  | 130707m  | 20.00 | ng    | 0.07     |
| 63) Phenanthrene-d10      | 11.45 | 188  | 216214m  | 20.00 | ng    | 0.08     |
| 75) Chrysene-d12          | 14.18 | 240  | 201193m  | 20.00 | ng    | 0.17     |
| 86) Perylene-d12          | 15.77 | 264  | 207898m  | 20.00 | ng    | 0.33     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|-------|------|----------|------|-------|----------|
| 5) 2-Fluorophenol           | 5.53  | 112  | 19749    | 7.52 | ng    | 0.10     |
| 7) Phenol-d6                | 6.56  | 99   | 26821    | 7.85 | ng    | 0.08     |
| 23) Nitrobenzene-d5         | 7.49  | 82   | 17075m   | 4.70 | ng    | 0.07     |
| 41) 2,4,6-Tribromophenol    | 10.75 | 330  | 9803m    | 8.00 | ng    | 0.08     |
| 44) 2-Fluorobiphenyl        | 9.27  | 172  | 48739m   | 6.18 | ng    | 0.06     |
| 78) Terphenyl-d14           | 13.04 | 244  | 53751m   | 7.08 | ng    | 0.08     |

| Target Compounds           | R.T.  | QIon | Response | Conc   | Units | Qvalue |
|----------------------------|-------|------|----------|--------|-------|--------|
| 27) 2,4-Dimethylphenol     | 8.00  | 122  | 14915m   | 4.72   | ng    |        |
| 31) Naphthalene            | 8.23  | 128  | 79793m   | 7.55   | ng    |        |
| 37) 2-Methylnaphthalene    | 8.91  | 142  | 142122m  | 19.39  | ng    |        |
| 45) 1,1'-Biphenyl          | 9.37  | 154  | 111541m  | 11.19  | ng    |        |
| 48) Acenaphthylene         | 9.82  | 152  | 93142m   | 7.62   | ng    |        |
| 51) Acenaphthene           | 9.99  | 154  | 482090m  | 65.70  | ng    |        |
| 54) Dibenzofuran           | 10.17 | 168  | 1089074m | 108.11 | ng    |        |
| 57) Fluorene               | 10.51 | 166  | 1336857m | 161.57 | ng    |        |
| 70) Phenanthrene           | 11.51 | 178  | 6305255m | 594.94 | ng    |        |
| 71) Anthracene             | 11.54 | 178  | 2023838m | 185.32 | ng    |        |
| 72) Carbazole              | 11.69 | 167  | 1247157m | 125.19 | ng    |        |
| 74) Fluoranthene           | 12.71 | 202  | 5953325m | 523.02 | ng    |        |
| 77) Pvrene                 | 12.93 | 202  | 4010573m | 335.90 | ng    |        |
| 80) Benzo(a)anthracene     | 14.17 | 228  | 1625626m | 155.29 | ng    |        |
| 82) Chrysene               | 14.22 | 228  | 1483214m | 134.47 | ng    |        |
| 85) Indeno(1,2,3-cd)pyrene | 17.26 | 276  | 147091m  | 16.78  | ng    |        |
| 87) Benzo(b)fluoranthene   | 15.34 | 252  | 1026529m | 81.16  | ng    |        |
| 88) Benzo(k)fluoranthene   | 15.35 | 252  | 417733m  | 39.29  | ng    |        |
| 89) Benzo(a)pvrene         | 15.70 | 252  | 342768m  | 31.53  | ng    |        |
| 90) Dibenzo(a,h)anthracene | 17.27 | 278  | 46104m   | 5.18   | ng    |        |
| 91) Benzo(g,h,i)perylene   | 17.70 | 276  | 100077m  | 11.56  | ng    |        |

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

