

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\  
 Data File : BF094123.D  
 Acq On : 3 Apr 2017 16:21  
 Operator : SJ/MA  
 Sample : PB97988BS  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB97988BS

Manual Integrations  
 APPROVED

mohammad  
 4/4/2017 5:33:57 PM

Quant Time: Apr 04 01:29:29 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  | 171204   | 20.00 | ng    | -0.02    |
| 21) Naphthalene-d8        | 7.39  | 136  | 682321   | 20.00 | ng    | -0.02    |
| 38) Acenaphthene-d10      | 9.53  | 164  | 321746   | 20.00 | ng    | -0.02    |
| 63) Phenanthrene-d10      | 11.36 | 188  | 556570   | 20.00 | ng    | -0.02    |
| 75) Chrysene-d12          | 14.62 | 240  | 446430   | 20.00 | ng    | -0.01    |
| 86) Perylene-d12          | 16.24 | 264  | 334874   | 20.00 | ng    | -0.01    |

## System Monitoring Compounds

|                          |       |     |         |        |    |       |
|--------------------------|-------|-----|---------|--------|----|-------|
| 5) 2-Fluorophenol        | 4.43  | 112 | 1271563 | 125.45 | ng | 0.00  |
| 7) Phenol-d6             | 5.50  | 99  | 1592066 | 126.96 | ng | 0.00  |
| 23) Nitrobenzene-d5      | 6.54  | 82  | 1011816 | 84.63  | ng | -0.02 |
| 41) 2,4,6-Tribromophenol | 10.52 | 330 | 360872  | 141.94 | ng | 0.00  |
| 44) 2-Fluorobiphenyl     | 8.72  | 172 | 1864472 | 88.39  | ng | -0.02 |
| 78) Terphenyl-d14        | 13.34 | 244 | 1586041 | 79.63  | ng | -0.01 |

## Target Compounds

|                                |      |     |         |       |    | Qvalue |
|--------------------------------|------|-----|---------|-------|----|--------|
| 2) 1,4-Dioxane                 | 2.29 | 88  | 163470  | 33.57 | ng | 99     |
| 3) Pyridine                    | 2.77 | 79  | 499811  | 37.66 | ng | 95     |
| 4) n-Nitrosodimethylamine      | 2.73 | 42  | 222205  | 40.69 | ng | 88     |
| 6) Aniline                     | 5.51 | 93  | 455294  | 26.10 | ng | # 43   |
| 8) 2-Chlorophenol              | 5.66 | 128 | 502073  | 45.06 | ng | 95     |
| 9) Benzaldehyde                | 5.38 | 77  | 381837  | 45.10 | ng | 99     |
| 10) Phenol                     | 5.52 | 94  | 580364  | 40.67 | ng | 90     |
| 11) bis(2-Chloroethyl)ether    | 5.60 | 93  | 464360  | 41.64 | ng | 96     |
| 12) 1,3-Dichlorobenzene        | 5.82 | 146 | 534588  | 42.78 | ng | 98     |
| 13) 1,4-Dichlorobenzene        | 5.91 | 146 | 541450  | 42.78 | ng | 97     |
| 14) 1,2-Dichlorobenzene        | 6.09 | 146 | 502439  | 43.10 | ng | 94     |
| 15) Benzyl Alcohol             | 6.06 | 79  | 407114  | 44.36 | ng | 98     |
| 16) 2,2'-oxybis(1-Chloropropan | 6.22 | 45  | 656338  | 38.20 | ng | 94     |
| 17) 2-Methylphenol             | 6.20 | 107 | 416992  | 43.70 | ng | 95     |
| 18) Hexachloroethane           | 6.47 | 117 | 190461  | 42.81 | ng | # 85   |
| 19) n-Nitroso-di-n-propylamine | 6.37 | 70  | 347045  | 43.06 | ng | 96     |
| 20) 3+4-Methylphenols          | 6.38 | 107 | 537044  | 46.47 | ng | # 63   |
| 22) Acetophenone               | 6.36 | 105 | 702343  | 46.18 | ng | # 86   |
| 24) Nitrobenzene               | 6.56 | 77  | 513287  | 43.53 | ng | 93     |
| 25) Isophorone                 | 6.85 | 82  | 928809  | 44.21 | ng | 95     |
| 26) 2-Nitrophenol              | 6.93 | 139 | 282689  | 50.27 | ng | 93     |
| 27) 2,4-Dimethylphenol         | 7.00 | 122 | 461668  | 47.65 | ng | 98     |
| 28) bis(2-Chloroethoxy)methane | 7.11 | 93  | 606434  | 43.77 | ng | 99     |
| 29) 2,4-Dichlorophenol         | 7.23 | 162 | 439491  | 48.51 | ng | 98     |
| 30) 1,2,4-Trichlorobenzene     | 7.33 | 180 | 423301  | 45.36 | ng | 97     |
| 31) Naphthalene                | 7.42 | 128 | 1418693 | 43.24 | ng | 100    |
| 32) Benzoic acid               | 7.16 | 122 | 167114  | 25.91 | ng | 91     |
| 33) 4-Chloroaniline            | 7.48 | 127 | 255572  | 19.22 | ng | 94     |
| 34) Hexachlorobutadiene        | 7.57 | 225 | 214924  | 44.91 | ng | 100    |
| 35) Caprolactam                | 7.93 | 113 | 140319m | 48.57 | ng |        |
| 36) 4-Chloro-3-methylphenol    | 8.11 | 107 | 461872  | 48.79 | ng | 92     |
| 37) 2-Methylnaphthalene        | 8.26 | 142 | 1003695 | 45.89 | ng | 98     |
| 39) 1,2,4,5-Tetrachlorobenzene | 8.47 | 216 | 369635  | 42.00 | ng | 99     |
| 40) Hexachlorocyclopentadiene  | 8.46 | 237 | 367969  | 89.34 | ng | 99     |

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\  
 Data File : BF094123.D  
 Acq On : 3 Apr 2017 16:21  
 Operator : SJ/MA  
 Sample : PB97988BS  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleID :  
 PB97988BS

Manual Integrations  
 APPROVED

mohammad  
 4/4/2017 5:33:57 PM

Quant Time: Apr 04 01:29:29 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) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 42) 2,4,6-Trichlorophenol      | 8.62  | 196  | 292220   | 46.93 | nq    | 99       |
| 43) 2,4,5-Trichlorophenol      | 8.67  | 196  | 296567   | 44.01 | nq    | 93       |
| 45) 1,1'-Biphenyl              | 8.84  | 154  | 1118687  | 43.11 | nq    | 98       |
| 46) 2-Chloronaphthalene        | 8.86  | 162  | 890224   | 43.82 | nq    | 95       |
| 47) 2-Nitroaniline             | 8.99  | 65   | 269869   | 44.78 | nq    | # 79     |
| 48) Acenaphthylene             | 9.36  | 152  | 1414022  | 41.88 | nq    | 99       |
| 49) Dimethylphthalate          | 9.23  | 163  | 1067655  | 46.68 | nq    | 100      |
| 50) 2,6-Dinitrotoluene         | 9.29  | 165  | 240971   | 45.96 | nq    | # 90     |
| 51) Acenaphthene               | 9.57  | 154  | 843804   | 42.83 | nq    | 100      |
| 52) 3-Nitroaniline             | 9.49  | 138  | 144792   | 23.52 | nq    | 88       |
| 53) 2,4-Dinitrophenol          | 9.63  | 184  | 194974   | 70.11 | nq    | # 69     |
| 54) Dibenzofuran               | 9.79  | 168  | 1222670  | 45.59 | nq    | 100      |
| 55) 4-Nitrophenol              | 9.74  | 139  | 374146   | 77.60 | nq    | # 61     |
| 56) 2,4-Dinitrotoluene         | 9.79  | 165  | 294494   | 44.72 | nq    | # 69     |
| 57) Fluorene                   | 10.21 | 166  | 924820   | 41.58 | nq    | 100      |
| 58) 2,3,4,6-Tetrachlorophenol  | 9.95  | 232  | 214777   | 46.45 | nq    | 99       |
| 59) Diethylphthalate           | 10.10 | 149  | 1032744  | 45.69 | nq    | 99       |
| 60) 4-Chlorophenyl-phenylether | 10.22 | 204  | 403167   | 42.55 | nq    | 97       |
| 61) 4-Nitroaniline             | 10.25 | 138  | 268042   | 45.37 | nq    | 93       |
| 62) Azobenzene                 | 10.42 | 77   | 999418   | 46.74 | nq    | 95       |
| 64) 4,6-Dinitro-2-methylphenol | 10.29 | 198  | 148736   | 43.64 | nq    | # 71     |
| 65) n-Nitrosodiphenylamine     | 10.37 | 169  | 881887   | 45.82 | nq    | 99       |
| 66) 4-Bromophenyl-phenylether  | 10.82 | 248  | 254509   | 46.49 | nq    | 96       |
| 67) Hexachlorobenzene          | 10.89 | 284  | 261470   | 47.19 | nq    | # 87     |
| 68) Atrazine                   | 11.04 | 200  | 263086   | 45.63 | nq    | 94       |
| 69) Pentachlorophenol          | 11.14 | 266  | 220363   | 63.89 | nq    | 96       |
| 70) Phenanthrene               | 11.39 | 178  | 1384947  | 44.73 | nq    | 99       |
| 71) Anthracene                 | 11.46 | 178  | 1419831  | 44.54 | nq    | 99       |
| 72) Carbazole                  | 11.66 | 167  | 1335039  | 40.84 | nq    | 99       |
| 73) Di-n-butylphthalate        | 12.11 | 149  | 1613539  | 47.46 | nq    | 99       |
| 74) Fluoranthene               | 12.86 | 202  | 1421989  | 44.85 | nq    | 98       |
| 76) Benzidine                  | 13.02 | 184  | 825874   | 46.74 | nq    | # 93     |
| 77) Pyrene                     | 13.13 | 202  | 1429370  | 44.12 | nq    | 99       |
| 79) Butylbenzylphthalate       | 13.94 | 149  | 689047   | 49.91 | nq    | # 85     |
| 80) Benzo(a)anthracene         | 14.60 | 228  | 1183005  | 45.44 | nq    | 99       |
| 81) 3,3'-Dichlorobenzidine     | 14.57 | 252  | 197490   | 21.22 | nq    | # 96     |
| 82) Chrysene                   | 14.64 | 228  | 1042657  | 43.16 | nq    | 99       |
| 83) Bis(2-ethylhexyl)phthalate | 14.66 | 149  | 941639   | 50.00 | nq    | 99       |
| 84) Di-n-octyl phthalate       | 15.42 | 149  | 1510076  | 47.99 | nq    | 98       |
| 85) Indeno(1,2,3-cd)pyrene     | 17.59 | 276  | 834900   | 39.90 | nq    | 100      |
| 87) Benzo(b)fluoranthene       | 15.84 | 252  | 992061   | 48.33 | nq    | 99       |
| 88) Benzo(k)fluoranthene       | 15.87 | 252  | 891794   | 44.40 | nq    | # 96     |
| 89) Benzo(a)pyrene             | 16.19 | 252  | 871153   | 46.09 | nq    | 99       |
| 90) Dibenzo(a,h)anthracene     | 17.61 | 278  | 688822   | 43.03 | nq    | 98       |
| 91) Benzo(g,h,i)perylene       | 17.99 | 276  | 693010   | 42.96 | nq    | 98       |

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\  
Data File : BF094123.D  
Acq On : 3 Apr 2017 16:21  
Operator : SJ/MA  
Sample : PB97988BS  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
PB97988BS

Manual Integrations  
APPROVED

mohammad  
4/4/2017 5:33:57 PM

Quant Time: Apr 04 01:29:29 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

