

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\  
 Data File : BG041111.D  
 Acq On : 7 Jun 2019 23:41  
 Operator : HP/JU  
 Sample : PB120405BS  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB120405BS

Manual Integrations  
 APPROVED

mohammad  
 6/11/2019 8:54:28 AM

Quant Time: Jun 08 18:04:40 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Jun 08 17:33:39 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.11  | 152  | 35978    | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 10.93 | 136  | 151343   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 14.74 | 164  | 120332   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 17.48 | 188  | 316283   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 21.76 | 240  | 316323   | 20.00 | ng    | 0.00     |
| 87) Perylene-d12          | 25.08 | 264  | 337667   | 20.00 | ng    | 0.00     |

## System Monitoring Compounds

|                          |       |     |         |        |    |      |
|--------------------------|-------|-----|---------|--------|----|------|
| 5) 2-Fluorophenol        | 5.64  | 112 | 284703  | 142.69 | ng | 0.00 |
| 7) Phenol-d6             | 7.25  | 99  | 418595  | 135.85 | ng | 0.00 |
| 23) Nitrobenzene-d5      | 9.28  | 82  | 278409  | 81.67  | ng | 0.00 |
| 42) 2,4,6-Tribromophenol | 16.23 | 330 | 239279  | 133.94 | ng | 0.00 |
| 45) 2-Fluorobiphenyl     | 13.36 | 172 | 685345  | 82.02  | ng | 0.00 |
| 79) Terphenyl-d14        | 20.07 | 244 | 1284364 | 86.17  | ng | 0.00 |

## Target Compounds

|                                |       |     |        |        |    | Qvalue |
|--------------------------------|-------|-----|--------|--------|----|--------|
| 2) 1,4-Dioxane                 | 3.49  | 88  | 36560  | 40.381 | ng | 98     |
| 3) Pyridine                    | 3.91  | 79  | 100467 | 37.501 | ng | 99     |
| 4) n-Nitrosodimethylamine      | 3.83  | 42  | 56763  | 45.653 | ng | 94     |
| 6) Aniline                     | 7.43  | 93  | 132859 | 32.302 | ng | 96     |
| 8) 2-Chlorophenol              | 7.67  | 128 | 118337 | 49.749 | ng | 99     |
| 9) Benzaldehyde                | 7.24  | 77  | 36567  | 19.893 | ng | 93     |
| 10) Phenol                     | 7.28  | 94  | 163688 | 49.544 | ng | 97     |
| 11) bis(2-Chloroethyl)ether    | 7.52  | 93  | 106287 | 44.345 | ng | 94     |
| 12) 1,3-Dichlorobenzene        | 7.99  | 146 | 119541 | 42.077 | ng | 95     |
| 13) 1,4-Dichlorobenzene        | 8.14  | 146 | 122518 | 42.604 | ng | 99     |
| 14) 1,2-Dichlorobenzene        | 8.46  | 146 | 121188 | 43.403 | ng | 99     |
| 15) Benzyl Alcohol             | 8.35  | 79  | 135067 | 47.385 | ng | 99     |
| 16) 2,2'-oxybis(1-Chloropropan | 8.62  | 45  | 162012 | 41.366 | ng | 95     |
| 17) 2-Methylphenol             | 8.54  | 107 | 114776 | 47.980 | ng | 98     |
| 18) Hexachloroethane           | 9.19  | 117 | 47163  | 42.552 | ng | 99     |
| 19) n-Nitroso-di-n-propylamine | 8.91  | 70  | 100763 | 42.492 | ng | 98     |
| 20) 3+4-Methylphenols          | 8.87  | 107 | 158951 | 47.969 | ng | 95     |
| 22) Acetophenone               | 8.93  | 105 | 193250 | 45.520 | ng | 98     |
| 24) Nitrobenzene               | 9.33  | 77  | 159895 | 46.024 | ng | 97     |
| 25) Isophorone                 | 9.84  | 82  | 297491 | 45.854 | ng | 98     |
| 26) 2-Nitrophenol              | 10.03 | 139 | 71188  | 49.704 | ng | 95     |
| 27) 2,4-Dimethylphenol         | 10.08 | 122 | 129605 | 54.791 | ng | 95     |
| 28) bis(2-Chloroethoxy)methane | 10.32 | 93  | 153734 | 43.904 | ng | 99     |
| 29) 2,4-Dichlorophenol         | 10.57 | 162 | 147521 | 49.112 | ng | 97     |
| 30) 1,2,4-Trichlorobenzene     | 10.79 | 180 | 149032 | 43.614 | ng | 98     |
| 31) Naphthalene                | 10.98 | 128 | 376344 | 46.315 | ng | 99     |
| 32) Benzoic acid               | 10.23 | 122 | 80627  | 44.541 | ng | 95     |
| 33) 4-Chloroaniline            | 11.09 | 127 | 90695  | 24.055 | ng | 98     |
| 34) Hexachlorobutadiene        | 11.24 | 225 | 108370 | 41.448 | ng | 97     |
| 35) Caprolactam                | 11.88 | 113 | 48615m | 49.011 | ng |        |
| 36) 4-Chloro-3-methylphenol    | 12.20 | 107 | 170391 | 49.669 | ng | 96     |
| 37) 2-Methylnaphthalene        | 12.57 | 142 | 295839 | 47.271 | ng | 100    |
| 38) 1-Methylnaphthalene        | 12.80 | 142 | 280324 | 47.533 | ng | 99     |
| 40) 1,2,4,5-Tetrachlorobenzene | 12.94 | 216 | 215104 | 44.351 | ng | 99     |

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\  
 Data File : BG041111.D  
 Acq On : 7 Jun 2019 23:41  
 Operator : HP/JU  
 Sample : PB120405BS  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
 PB120405BS

Manual Integrations  
 APPROVED

mohammad  
 6/11/2019 8:54:28 AM

Quant Time: Jun 08 18:04:40 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Jun 08 17:33:39 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc    | Units | Dev(Min) |
|--------------------------------|-------|------|----------|---------|-------|----------|
| 41) Hexachlorocyclopentadiene  | 12.90 | 237  | 278697   | 104.130 | nq    | 99       |
| 43) 2,4,6-Trichlorophenol      | 13.17 | 196  | 145233   | 46.289  | nq    | 96       |
| 44) 2,4,5-Trichlorophenol      | 13.25 | 196  | 160109   | 48.386  | nq    | 95       |
| 46) 1,1'-Biphenyl              | 13.58 | 154  | 410688   | 46.107  | nq    | 97       |
| 47) 2-Chloronaphthalene        | 13.62 | 162  | 340787   | 44.567  | nq    | 99       |
| 48) 2-Nitroaniline             | 13.83 | 65   | 125680   | 45.814  | nq    | 95       |
| 49) Acenaphthylene             | 14.46 | 152  | 531644   | 46.195  | nq    | 99       |
| 50) Dimethylphthalate          | 14.19 | 163  | 458738   | 45.035  | nq    | 99       |
| 51) 2,6-Dinitrotoluene         | 14.32 | 165  | 103131   | 46.958  | nq    | 97       |
| 52) Acenaphthene               | 14.80 | 154  | 318530   | 45.688  | nq    | 97       |
| 53) 3-Nitroaniline             | 14.65 | 138  | 69900    | 31.829  | nq    | # 97     |
| 54) 2,4-Dinitrophenol          | 14.86 | 184  | 137968   | 105.977 | nq    | 98       |
| 55) Dibenzofuran               | 15.13 | 168  | 554858   | 47.297  | nq    | 100      |
| 56) 4-Nitrophenol              | 14.95 | 139  | 159903   | 106.152 | nq    | 98       |
| 57) 2,4-Dinitrotoluene         | 15.10 | 165  | 149841   | 48.299  | nq    | 96       |
| 58) Fluorene                   | 15.78 | 166  | 439626   | 47.056  | nq    | 99       |
| 59) 2,3,4,6-Tetrachlorophenol  | 15.35 | 232  | 164579   | 50.610  | nq    | 97       |
| 60) Diethylphthalate           | 15.54 | 149  | 478236   | 46.005  | nq    | 98       |
| 61) 4-Chlorophenyl-phenylether | 15.77 | 204  | 278488   | 45.860  | nq    | 96       |
| 62) 4-Nitroaniline             | 15.82 | 138  | 103489   | 44.511  | nq    | 97       |
| 63) Azobenzene                 | 16.06 | 77   | 440933   | 46.669  | nq    | 99       |
| 65) 4,6-Dinitro-2-methylphenol | 15.86 | 198  | 91683    | 53.005  | nq    | 95       |
| 66) n-Nitrosodiphenylamine     | 15.99 | 169  | 402168   | 45.233  | nq    | 97       |
| 67) 4-Bromophenyl-phenylether  | 16.66 | 248  | 180722   | 42.340  | nq    | 98       |
| 68) Hexachlorobenzene          | 16.77 | 284  | 191026   | 42.029  | nq    | 98       |
| 69) Atrazine                   | 16.93 | 200  | 170837   | 49.797  | nq    | 97       |
| 70) Pentachlorophenol          | 17.13 | 266  | 230736   | 93.979  | nq    | 99       |
| 71) Phenanthrene               | 17.52 | 178  | 761365   | 46.214  | nq    | 99       |
| 72) Anthracene                 | 17.62 | 178  | 774006   | 47.636  | nq    | 99       |
| 73) Carbazole                  | 17.89 | 167  | 668891   | 47.977  | nq    | 100      |
| 74) Di-n-butylphthalate        | 18.42 | 149  | 781427   | 45.096  | nq    | 100      |
| 75) Fluoranthene               | 19.53 | 202  | 984636   | 48.387  | nq    | 98       |
| 77) Benzidine                  | 19.71 | 184  | 408439   | 54.127  | nq    | 99       |
| 78) Pyrene                     | 19.89 | 202  | 974231   | 47.378  | nq    | 98       |
| 80) Butylbenzylphthalate       | 20.76 | 149  | 369158   | 46.132  | nq    | 100      |
| 81) Benzo(a)anthracene         | 21.74 | 228  | 945366   | 44.897  | nq    | 98       |
| 82) 3,3'-Dichlorobenzidine     | 21.66 | 252  | 220397   | 29.759  | nq    | 100      |
| 83) Chrysene                   | 21.81 | 228  | 925605   | 45.935  | nq    | 100      |
| 84) Bis(2-ethylhexyl)phthalate | 21.61 | 149  | 505120   | 44.840  | nq    | 99       |
| 85) Di-n-octyl phthalate       | 22.85 | 149  | 858569   | 45.763  | nq    | 100      |
| 86) Indeno(1,2,3-cd)pyrene     | 28.91 | 276  | 1128099  | 45.703  | nq    | 100      |
| 88) Benzo(b)fluoranthene       | 24.02 | 252  | 946075   | 46.194  | nq    | 98       |
| 89) Benzo(k)fluoranthene       | 24.09 | 252  | 938403   | 46.302  | nq    | 99       |
| 90) Benzo(a)pyrene             | 24.93 | 252  | 939565   | 48.084  | nq    | 100      |
| 91) Dibenzo(a,h)anthracene     | 28.98 | 278  | 921311   | 47.187  | nq    | 99       |
| 92) Benzo(g,h,i)perylene       | 30.12 | 276  | 930973   | 49.079  | nq    | 99       |

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

