

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081718\  
 Data File : BM016441.D  
 Acq On : 17 Aug 2018 16:28  
 Operator : SJ/JU  
 Sample : SSTDICV040  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 ICBM081718

Manual Integrations  
 APPROVED

Sohil  
 8/20/2018 4:17:08 PM

Quant Time: Aug 17 17:04:48 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM081718.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Aug 17 16:27:00 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.06  | 152  | 18571    | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 10.87 | 136  | 83164    | 20.00 | ng    | 0.00     |
| 38) Acenaphthene-d10      | 14.70 | 164  | 53961    | 20.00 | ng    | 0.00     |
| 63) Phenanthrene-d10      | 17.44 | 188  | 134401   | 20.00 | ng    | 0.00     |
| 75) Chrysene-d12          | 21.59 | 240  | 179624   | 20.00 | ng    | 0.00     |
| 86) Perylene-d12          | 23.90 | 264  | 147716   | 20.00 | ng    | 0.00     |

## System Monitoring Compounds

|                          |       |     |        |       |    |      |
|--------------------------|-------|-----|--------|-------|----|------|
| 5) 2-Fluorophenol        | 5.59  | 112 | 82725  | 74.86 | ng | 0.00 |
| 7) Phenol-d6             | 7.23  | 99  | 112179 | 75.53 | ng | 0.00 |
| 23) Nitrobenzene-d5      | 9.25  | 82  | 157224 | 78.73 | ng | 0.00 |
| 41) 2,4,6-Tribromophenol | 16.19 | 330 | 62963  | 77.25 | ng | 0.00 |
| 44) 2-Fluorobiphenyl     | 13.32 | 172 | 388024 | 74.77 | ng | 0.00 |
| 78) Terphenyl-d14        | 20.03 | 244 | 836008 | 70.34 | ng | 0.00 |

## Target Compounds

|                                |       |     |        |        |    | Qvalue |
|--------------------------------|-------|-----|--------|--------|----|--------|
| 2) 1,4-Dioxane                 | 3.40  | 88  | 21582  | 39.000 | ng | # 100  |
| 3) Pyridine                    | 3.83  | 79  | 46315  | 41.571 | ng | # 68   |
| 4) n-Nitrosodimethylamine      | 3.76  | 42  | 52394  | 40.815 | ng | # 22   |
| 6) Aniline                     | 7.39  | 93  | 59459  | 38.686 | ng | # 81   |
| 8) 2-Chlorophenol              | 7.62  | 128 | 50578  | 37.125 | ng | # 95   |
| 9) Benzaldehyde                | 7.21  | 77  | 38920  | 39.852 | ng | # 81   |
| 10) Phenol                     | 7.26  | 94  | 54187  | 37.662 | ng | # 83   |
| 11) bis(2-Chloroethyl)ether    | 7.49  | 93  | 41300  | 37.811 | ng | # 85   |
| 12) 1,3-Dichlorobenzene        | 7.94  | 146 | 57813  | 37.026 | ng | # 82   |
| 13) 1,4-Dichlorobenzene        | 8.09  | 146 | 57884  | 36.671 | ng | # 93   |
| 14) 1,2-Dichlorobenzene        | 8.41  | 146 | 59155  | 38.149 | ng | # 93   |
| 15) Benzyl Alcohol             | 8.32  | 79  | 49810  | 36.946 | ng | # 77   |
| 16) 2,2'-oxybis(1-Chloropropan | 8.59  | 45  | 52503  | 38.185 | ng | # 88   |
| 17) 2-Methylphenol             | 8.52  | 107 | 38366  | 36.903 | ng | # 75   |
| 18) Hexachloroethane           | 9.12  | 117 | 32181  | 39.590 | ng | # 100  |
| 19) n-Nitroso-di-n-propylamine | 8.88  | 70  | 44408  | 39.380 | ng | # 65   |
| 20) 3+4-Methylphenols          | 8.85  | 107 | 55873  | 39.107 | ng | # 80   |
| 22) Acetophenone               | 8.90  | 105 | 83799  | 37.571 | ng | # 73   |
| 24) Nitrobenzene               | 9.29  | 77  | 74512  | 38.709 | ng | # 94   |
| 25) Isophorone                 | 9.80  | 82  | 103586 | 38.892 | ng | # 91   |
| 26) 2-Nitrophenol              | 10.00 | 139 | 30762  | 37.989 | ng | # 34   |
| 27) 2,4-Dimethylphenol         | 10.05 | 122 | 41721  | 38.541 | ng | # 69   |
| 28) bis(2-Chloroethoxy)methane | 10.28 | 93  | 59000  | 38.295 | ng | # 97   |
| 29) 2,4-Dichlorophenol         | 10.53 | 162 | 56150  | 38.369 | ng | # 97   |
| 30) 1,2,4-Trichlorobenzene     | 10.73 | 180 | 65818  | 38.033 | ng | # 94   |
| 31) Naphthalene                | 10.92 | 128 | 160813 | 38.127 | ng | # 99   |
| 32) Benzoic acid               | 10.25 | 122 | 36223m | 41.197 | ng | #      |
| 33) 4-Chloroaniline            | 11.05 | 127 | 69304  | 38.338 | ng | # 86   |
| 34) Hexachlorobutadiene        | 11.17 | 225 | 57792  | 38.742 | ng | # 95   |
| 35) Caprolactam                | 11.87 | 113 | 14684  | 39.139 | ng | # 83   |
| 36) 4-Chloro-3-methylphenol    | 12.17 | 107 | 62746  | 37.600 | ng | # 85   |
| 37) 2-Methylnaphthalene        | 12.52 | 142 | 122128 | 39.030 | ng | # 82   |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.89 | 216 | 83206  | 37.666 | ng | # 100  |
| 40) Hexachlorocyclopentadiene  | 12.85 | 237 | 27181  | 38.343 | ng | # 94   |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081718\  
 Data File : BM016441.D  
 Acq On : 17 Aug 2018 16:28  
 Operator : SJ/JU  
 Sample : SSTDICV040  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 ICVBM081718

Manual Integrations  
 APPROVED

Sohil  
 8/20/2018 4:17:08 PM

Quant Time: Aug 17 17:04:48 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM081718.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Aug 17 16:27:00 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 42) 2,4,6-Trichlorophenol      | 13.15 | 196  | 51708    | 37.451 | nq    | 99       |
| 43) 2,4,5-Trichlorophenol      | 13.23 | 196  | 53040    | 39.376 | nq #  | 79       |
| 45) 1,1'-Biphenyl              | 13.53 | 154  | 186928   | 37.427 | nq    | 99       |
| 46) 2-Chloronaphthalene        | 13.58 | 162  | 148911   | 38.217 | nq #  | 86       |
| 47) 2-Nitroaniline             | 13.81 | 65   | 53601    | 39.669 | nq #  | 73       |
| 48) Acenaphthylene             | 14.43 | 152  | 225381   | 38.468 | nq    | 98       |
| 49) Dimethylphthalate          | 14.16 | 163  | 206612   | 39.044 | nq    | 99       |
| 50) 2,6-Dinitrotoluene         | 14.30 | 165  | 40145    | 39.970 | nq #  | 31       |
| 51) Acenaphthene               | 14.77 | 154  | 132674   | 37.524 | nq    | 98       |
| 52) 3-Nitroaniline             | 14.64 | 138  | 40251    | 40.675 | nq #  | 76       |
| 53) 2,4-Dinitrophenol          | 14.87 | 184  | 19016    | 40.125 | nq #  | 54       |
| 54) Dibenzofuran               | 15.10 | 168  | 241806   | 38.813 | nq #  | 76       |
| 55) 4-Nitrophenol              | 14.96 | 139  | 26824    | 41.499 | nq #  | 84       |
| 56) 2,4-Dinitrotoluene         | 15.10 | 165  | 65526    | 39.029 | nq #  | 53       |
| 57) Fluorene                   | 15.75 | 166  | 185107   | 38.890 | nq    | 96       |
| 58) 2,3,4,6-Tetrachlorophenol  | 15.33 | 232  | 51236    | 40.598 | nq #  | 100      |
| 59) Diethylphthalate           | 15.52 | 149  | 235106   | 39.154 | nq    | 95       |
| 60) 4-Chlorophenyl-phenylether | 15.73 | 204  | 119798   | 39.784 | nq #  | 84       |
| 61) 4-Nitroaniline             | 15.81 | 138  | 39477    | 41.260 | nq #  | 1        |
| 62) Azobenzene                 | 16.03 | 77   | 275258   | 39.750 | nq #  | 68       |
| 64) 4,6-Dinitro-2-methylphenol | 15.87 | 198  | 36888    | 39.358 | nq    | 84       |
| 65) n-Nitrosodiphenylamine     | 15.96 | 169  | 172386   | 37.904 | nq    | 97       |
| 66) 4-Bromophenyl-phenylether  | 16.63 | 248  | 67691    | 37.446 | nq #  | 73       |
| 67) Hexachlorobenzene          | 16.74 | 284  | 68922    | 38.852 | nq #  | 55       |
| 68) Atrazine                   | 16.90 | 200  | 66985    | 40.178 | nq    | 95       |
| 69) Pentachlorophenol          | 17.10 | 266  | 38694    | 38.095 | nq    | 98       |
| 70) Phenanthrene               | 17.49 | 178  | 289831   | 38.062 | nq    | 99       |
| 71) Anthracene                 | 17.57 | 178  | 285788   | 37.806 | nq    | 97       |
| 72) Carbazole                  | 17.85 | 167  | 287137   | 39.245 | nq #  | 95       |
| 73) Di-n-butylphthalate        | 18.38 | 149  | 415983   | 39.617 | nq #  | 95       |
| 74) Fluoranthene               | 19.49 | 202  | 409252   | 44.490 | nq    | 96       |
| 76) Benzidine                  | 19.67 | 184  | 155660   | 39.374 | nq    | 98       |
| 77) Pyrene                     | 19.84 | 202  | 427936   | 36.102 | nq    | 98       |
| 79) Butylbenzylphthalate       | 20.71 | 149  | 209413   | 37.695 | nq #  | 73       |
| 80) Benzo(a)anthracene         | 21.57 | 228  | 433954   | 37.340 | nq    | 98       |
| 81) 3,3'-Dichlorobenzidine     | 21.50 | 252  | 160986   | 38.911 | nq    | 99       |
| 82) Chrysene                   | 21.62 | 228  | 407275   | 37.290 | nq    | 98       |
| 83) Bis(2-ethylhexyl)phthalate | 21.46 | 149  | 314737   | 39.715 | nq    | 92       |
| 84) Di-n-octyl phthalate       | 22.36 | 149  | 499708   | 42.039 | nq #  | 86       |
| 85) Indeno(1,2,3-cd)pyrene     | 26.25 | 276  | 355681   | 39.814 | nq    | 99       |
| 87) Benzo(b)fluoranthene       | 23.20 | 252  | 359401   | 36.731 | nq #  | 92       |
| 88) Benzo(k)fluoranthene       | 23.24 | 252  | 377065   | 38.561 | nq #  | 96       |
| 89) Benzo(a)pyrene             | 23.80 | 252  | 341154   | 38.518 | nq #  | 97       |
| 90) Dibenzo(a,h)anthracene     | 26.24 | 278  | 302442   | 38.079 | nq #  | 91       |
| 91) Benzo(g,h,i)perylene       | 26.97 | 276  | 255155   | 35.744 | nq #  | 88       |

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

