

# PERIODIC TABLE OF THE ELEMENTS

|                   |                   |                   |                   |                   |                  |                   |                   |                   |                   |                   |                   |                   |                   |                   |                   |                   |                   |
|-------------------|-------------------|-------------------|-------------------|-------------------|------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| 1<br>H<br>1.0     | 2                 | 3                 | 4                 | 5                 | 6                | 7                 | 8                 | 9                 | 10                | 11                | 12                | 13                | 14                | 15                | 16                | 17<br>H<br>1.0    | 18<br>He<br>4.0   |
| 3<br>Li<br>6.9    | 4<br>Be<br>9.0    |                   |                   |                   |                  |                   |                   |                   |                   |                   |                   | 5<br>B<br>10.8    | 6<br>C<br>12.0    | 7<br>N<br>14.0    | 8<br>O<br>16.0    | 9<br>F<br>19.0    | 10<br>Ne<br>20.2  |
| 11<br>Na<br>23.0  | 12<br>Mg<br>24.3  |                   |                   |                   |                  |                   |                   |                   |                   |                   |                   | 13<br>Al<br>27.0  | 14<br>Si<br>28.1  | 15<br>P<br>31.0   | 16<br>S<br>32.1   | 17<br>Cl<br>35.5  | 18<br>Ar<br>39.9  |
| 19<br>K<br>39.1   | 20<br>Ca<br>40.1  | 21<br>Sc<br>45.0  | 22<br>Ti<br>47.9  | 23<br>V<br>50.9   | 24<br>Cr<br>52.0 | 25<br>Mn<br>54.9  | 26<br>Fe<br>55.8  | 27<br>Co<br>58.9  | 28<br>Ni<br>58.7  | 29<br>Cu<br>63.5  | 30<br>Zn<br>65.4  | 31<br>Ga<br>69.7  | 32<br>Ge<br>72.6  | 33<br>As<br>74.9  | 34<br>Se<br>79.0  | 35<br>Br<br>79.9  | 36<br>Kr<br>83.8  |
| 37<br>Rb<br>85.5  | 38<br>Sr<br>87.6  | 39<br>Y<br>88.9   | 40<br>Zr<br>91.2  | 41<br>Nb<br>92.9  | 42<br>Mo<br>95.9 | 43<br>Tc<br>98.0  | 44<br>Ru<br>101.1 | 45<br>Rh<br>102.9 | 46<br>Pd<br>106.4 | 47<br>Ag<br>107.9 | 48<br>Cd<br>112.4 | 49<br>In<br>114.8 | 50<br>Sn<br>118.7 | 51<br>Sb<br>121.8 | 52<br>Te<br>127.6 | 53<br>I<br>126.9  | 54<br>Xe<br>131.3 |
| 55<br>Cs<br>132.9 | 56<br>Ba<br>137.3 | 57<br>La<br>138.9 | 72<br>Hf<br>178.5 | 73<br>Ta<br>181.0 | 74<br>W<br>183.9 | 75<br>Re<br>186.2 | 76<br>Os<br>190.2 | 77<br>Ir<br>192.2 | 78<br>Pt<br>195.1 | 79<br>Au<br>197.0 | 80<br>Hg<br>200.6 | 81<br>Tl<br>204.4 | 82<br>Pb<br>207.2 | 83<br>Bi<br>209.0 | 84<br>Po<br>209.0 | 85<br>At<br>210.0 | 86<br>Rn<br>222.0 |

## Characteristic Infrared Absorptions of Some Functional Groups

| Functional Group | Bond | Frequency Range (cm <sup>-1</sup> ) | Functional Group | Bond | Frequency Range (cm <sup>-1</sup> ) |
|------------------|------|-------------------------------------|------------------|------|-------------------------------------|
| Alcohol          | O-H  | 3400 - 3650 (s, broad)              | Nitrile          | C≡N  | 2210 - 2260 (m)                     |
|                  | C-O  | 1050 - 1150 (s)                     | Carboxylic acid  | O-H  | 2500 - 3100 (s, broad)              |
| Ether            | C-O  | 1000 - 1260                         |                  | C=O  | 1700 - 1720 (s)                     |
| Amine            | N-H  | 3300 - 3500 (m)                     | Ester            | C=O  | 1710 - 1750 (s)                     |
| Alkane           | C-H  | 2850 - 2960 (m - s)                 | Acyl halide      | C=O  | 1770 - 1820 (s)                     |
| Alkene           | =C-H | 3020 - 3100 (m)                     | Acid anhydride   | C=O  | 1740 - 1790 (s)                     |
|                  | C=C  | 1640 - 1680 (m)                     |                  |      | 1800 - 1850 (s)                     |
| Alkyne           | ≡C-H | 3270 - 3330 (s)                     | Amide            | C=O  | 1630 - 1700 (s)                     |
|                  | C≡C  | 2100 - 2260 (m)                     | Aldehyde, ketone | C=O  | 1680 - 1730 (s)                     |

## Characteristic Hydrogen (<sup>1</sup>H NMR) Chemical Shifts

| Type of Hydrogen                | Structure                          | Chemical Shift, δ (ppm) | Type of Hydrogen | Structure                | Chemical Shift, δ (ppm) |
|---------------------------------|------------------------------------|-------------------------|------------------|--------------------------|-------------------------|
| Reference                       | (CH <sub>3</sub> ) <sub>4</sub> Si | 0.00                    | Amines           | —N—CH <sub>3</sub>       | 2.3 - 3.0               |
| Alkane, primary                 | —CH <sub>3</sub>                   | 0.7 - 1.3               | Alcohol, ether   | —O—CH <sub>3</sub>       | 3.3 - 4.0               |
| Alkane, secondary               | —CH <sub>2</sub> —                 | 1.2 - 1.4               | Ester            | —C(=O)—O—CH <sub>3</sub> | 3.7 - 4.2               |
| Alkane, tertiary                | —CH—                               | 1.4 - 1.7               | Vinyl            | C=CH <sub>2</sub>        | 5.0 - 6.5               |
| Allylic, primary                | C=C—CH <sub>3</sub>                | 1.6 - 1.9               | Aromatic         | Ar—H                     | 6.5 - 8.0               |
| Methyl ketone                   | —C(=O)—CH <sub>3</sub>             | 2.1 - 2.4               | Aldehyde         | —C(=O)—H                 | 9.7 - 10.0              |
| Aromatic methyl                 | Ar—CH <sub>3</sub>                 | 2.5 - 2.7               | Amine            | —NH <sub>2</sub>         | 1 - 5 variable          |
| Alkyne                          | —C≡CH                              | 2.5 - 2.7               | Alcohol          | —OH                      | 1 - 5 variable          |
| Alkyl halide<br>(X = Br, Cl, I) | —CH <sub>2</sub> —X                | 2.5 - 4.0               | Carboxylic acid  | —COOH                    | 11.0 - 12.0             |