## Advanced (Subsidiary) GCE

CHEMISTRY A
Data Sheet

## Specimen



The information in this Sheet is for the use of candidates following Chemistry A H034 and H434.

A copy of this sheet will be included as an insert with each unit paper.
Copies of this sheet may be used for teaching.

## General information

- 1 mol of gas molecules occupies $24.0 \mathrm{dm}^{3}$ at room temperature and pressure, RTP.
- Avogadro constant, $N_{\mathrm{A}}=6.02 \times 10^{23} \mathrm{~mol}^{-1}$.
- Ionic product of water, $K_{\mathrm{w}}=1.00 \times 10^{-14} \mathrm{~mol}^{2} \mathrm{dm}^{-6}$.


## ${ }^{1} \mathrm{H}$ NMR chemical shifts relative to TMS

Chemical shifts are typical values and can vary slightly depending on the solvent, concentration and substituents.

| type of proton |  |  | chemical shift, $\delta / \mathrm{ppm}$ |
| :---: | :---: | :---: | :---: |
| $\mathrm{R}-\mathrm{CH}_{3}$ |  |  | 0.7-1.6 |
| $\mathrm{N}-\mathrm{H}$ |  | $\mathrm{R}-\mathrm{OH}$ | 1.0-5.5* |
| $\mathrm{R}-\mathrm{CH}_{2}-\mathrm{R}$ |  |  | 1.2-1.4 |
| $\mathrm{R}_{3} \mathrm{CH}$ |  |  | 1.6-2.0 |
|  |  |  | 2.0-2.9 |
|    <br> 2.3-2.7 |  |  |  |
| $\mathrm{N}-\mathrm{CH}_{3} \quad \mathrm{~N}-\mathrm{CH}_{2} \mathrm{R}$ |  |  | 2.3-2.9 |
| $\mathrm{O}-\mathrm{CH}_{3} \quad \mathrm{O}-\mathrm{CH}_{2} \mathrm{R}$ |  | $\mathrm{O}-\mathrm{CHR}_{2}$ | 3.3-4.3 |
| Br or $\mathrm{Cl}-\mathrm{CH}_{3}$ | Br or $\mathrm{Cl}-\mathrm{CH}_{2} \mathrm{R}$ | Br or $\mathrm{Cl}-\mathrm{CHR}_{2}$ | 3.0-4.2 |
|  |  |  | 4.5-10.0* |
| $-\mathrm{CH}=\mathrm{CH}-$ |  |  | 4.5-6.0 |
|  |  |  | 5.0-12.0* |
|  |  |  | 6.5-8.0 |
|  |  |  | 9.0-10 |
|  |  |  | 11.0-12.0* |

* OH and NH chemical shifts are very variable (sometimes outside these limits) and are often broad. Signals are not usually seen as split peaks.
${ }^{13} \mathrm{C}$ NMR chemical shifts relative to TMS
Chemical shifts are typical values and can vary slightly depending on the solvent, concentration and substituents.

| type of carbon | chemical shift, $\delta$ I ppm |
| :---: | :---: |
| C-C (alkanes) | 10-35 |
|  | 20-30 |
| $\mathrm{C}-\mathrm{Cl}$ or $\mathrm{C}-\mathrm{Br}$ | 30-70 |
| $\mathrm{C}-\mathrm{N}$ (amines) | 35-60 |
| $\mathrm{C}-\mathrm{OH}$ | 50-65 |
| $\mathbf{C = C}$ (alkenes) | 115-140 |
| aromatic | 125-150 |
| carbonyl (ester, carboxylic acid, amide) | 160-185 |
| carbonyl (aldehyde, ketone) | 190-220 |

## Characteristic infrared absorptions in organic molecules

| bond | location | wavenumber/cm $^{\mathbf{- 1}}$ |
| :--- | :--- | :---: |
| C-O | alcohols, esters, carboxylic acids | $1000-1300$ |
| C=O | aldehydes, ketones, carboxylic acids, esters, <br> amides | $1640-1750$ |
| C-H | organic compound with a C-H bond | $2850-3100$ |
| O-H | carboxylic acids | $2500-3300$ (very broad) |
| N-H | amines, amides | $3200-3500$ |
| O-H | alcohols, phenols | $3200-3550$ (broad) |

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## The Periodic Table of the Elements



| $\begin{gathered} 140.1 \\ \text { Ce } \\ \text { cerium } \\ 58 \end{gathered}$ | 140.9 Pr prasedymum 59 | $\begin{gathered} 144.2 \\ \mathrm{Nd} \\ \text { neodymium } \\ 60 \end{gathered}$ | $\begin{gathered} 144.9 \\ \text { Pm } \\ \text { promethium } \\ 61 \end{gathered}$ | $\begin{gathered} 150.4 \\ \mathbf{S m} \\ \text { samarium } \\ 62 \end{gathered}$ | $\begin{gathered} 152.0 \\ \text { Eu } \\ \text { europium } \\ 63 \end{gathered}$ | $\begin{gathered} 157.2 \\ \text { Gd } \\ \text { gadolinium } \\ 64 \end{gathered}$ | $\begin{gathered} 158.9 \\ \mathrm{~Tb} \\ \text { terbium } \\ 65 \end{gathered}$ | $\begin{gathered} 162.5 \\ \text { Dy } \\ \text { dysprosium } \\ 66 \end{gathered}$ | $\begin{gathered} 164.9 \\ \text { Ho } \\ \text { nolmium } \\ 67 \end{gathered}$ | $\begin{gathered} 167.3 \\ \text { Er } \\ \text { erbium } \\ 68 \end{gathered}$ | $\begin{gathered} 168.9 \\ \mathrm{Tm} \\ \text { tuxuium } \\ 69 \end{gathered}$ | $\begin{gathered} 173.0 \\ \text { Yb } \\ \text { ytterbium } \\ 70 \end{gathered}$ | $\begin{gathered} 175.0 \\ \text { Lu } \\ \text { luetium } \\ 71 \end{gathered}$ |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 232.0 | [231] | 238.1 | [237] | [242] | [243] | [247] | [245] | [251] | [254] | [253] | [256] | [254] | [257] |
| $\begin{gathered} \text { Th } \\ \text { thorium } \\ 90 \end{gathered}$ | $\begin{aligned} & \mathrm{Pa} \\ & \text { protacinium } \\ & 91 \end{aligned}$ | $\begin{gathered} \mathbf{U} \\ \text { uranium } \\ 92 \end{gathered}$ | $\begin{gathered} \text { Np } \\ \text { neptunium } \\ 93 \end{gathered}$ | $\begin{gathered} \mathrm{Pu} \\ \text { plutonium } \\ 94 \end{gathered}$ | $\begin{gathered} \text { americium } \\ 95 \end{gathered}$ | $\begin{gathered} \text { Curium } \\ 96 \end{gathered}$ | $\begin{gathered} \text { Bk } \\ \text { berkelium } \\ 97 \end{gathered}$ | $\begin{gathered} \text { Cf } \\ \text { californium } \\ 98 \end{gathered}$ | Es einsteinium 99 | $\begin{aligned} & \text { Fm } \\ & \text { fermium } \\ & 100 \end{aligned}$ | $\begin{gathered} \mathbf{M d} \\ \text { mendelevium } \\ 101 \end{gathered}$ | $\begin{aligned} & \text { No } \\ & \text { nobelium } \\ & 102 \end{aligned}$ | $\begin{gathered} \mathbf{L r} \\ \text { lawencium } \\ 103 \end{gathered}$ |

