

Quantities, Units, Symbols and Nomenclature used in NCEA Chemistry Level 3 and Scholarship Examination Papers

NCEA Chemistry examinations will use the following information, which has been based on International Union of Pure and Applied Chemistry (IUPAC) recommendations. Candidates should be encouraged to use this IUPAC terminology, but those who use other terminology will not be penalised if their answers indicate a clear understanding of the chemistry involved.

General Chemistry

Symbols for the physical quantities, *M*, *V*, *H*, *s*, *K*, are written in italics (sloping letters). Any following subscripts will be in upright type.

Symbols / Expressions	Units in common use
<i>M</i> , molar mass, is the mass of one mole of a defined substance and will be used for elements and compounds. <i>M_r</i> , relative molecular mass, and <i>A_r</i> , relative atomic mass, will not be used.	g mol ⁻¹
<i>V</i> , volume. A looped <i>ℓ</i> is not used in these abbreviations.	L and mL
<i>n</i> , amount of substance, expressed in moles. It is incorrect to use the term 'number of moles'. (See details under 'Amount of Substance' below.)	mol
<i>c</i> , amount concentration, is expressed as moles per litre, also denoted by the format []. Concentrations may also be written as <i>mass concentration</i> , expressed as grams per litre.	mol L ⁻¹
<i>Composition of a mixture</i> , commonly expressed as % w/ <i>V</i> , % w/w and % <i>V/V</i> , will be used only after giving a clear definition of their meaning (eg grams per 100 mL, grams per 100 g, mL per 100 mL respectively).	g L ⁻¹
<i>s</i> (<i>italic s</i>), solubility, units as for concentration.	mol L ⁻¹

Amount of Substance

This is a physical quantity, symbol *n* (*italic n*), measured in a unit called the mole, which has the abbreviation mol.

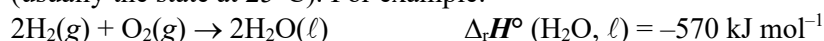
The term 'number of moles' is to be avoided in favour of the 'amount of substance in moles'. In the same manner, the size of an object can be described in terms of its 'length in metres', rather than its 'number of metres'.

Graph Axes and Table Headings

Labelled as quantity / unit, eg *c* / mol L⁻¹. Only values will then be written on the axes or in a table.

Enthalpy changes, ΔH Units commonly used kJ mol^{-1}

$\Delta_r H^\circ$, standard enthalpy of reaction when reactants and products are in their standard state (usually the state at 25°C). For example:

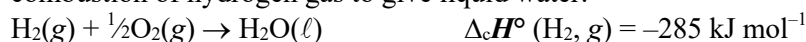


The term mol^{-1} means per mole of reaction, which is determined by the chemical equation; ie 2 mol of H_2 reacting with 1 mol of O_2 to give 2 mol of H_2O .

$\Delta_f H^\circ$, standard enthalpy of formation, per mole of product. For example, the standard enthalpy of formation of liquid water:



$\Delta_c H^\circ$, standard enthalpy of combustion, per mole of substance burnt. For example, the standard enthalpy of combustion of hydrogen gas to give liquid water:



- Note (i) The superscript $^\circ$ denotes a defined standard state.
(ii) The alternative superscript $^\theta$ (plimsoll) is acceptable.
(iii) A space is always left between any value and its unit, as well as between units for composite units.

$\Delta_{\text{fus}} H$, enthalpy of fusion (melting)

$\Delta_{\text{vap}} H$, enthalpy of vaporisation

$\Delta_{\text{sub}} H$, enthalpy of sublimation

Standard Electrode Potential

Electrode potentials are defined as standard electrode potentials, E° .

Units are volts, symbol V.

eg	Redox couple	E° / V
	$\text{Zn}^{2+} / \text{Zn}$	-0.76
	$\text{Fe}^{3+} / \text{Fe}^{2+}$	+0.77

A half cell is an electrode and the couple it is in contact with. When the oxidant and reductant are in different phases, a vertical line in the cell diagram is used to represent the phase boundary.

For example

$\text{Zn}(\text{s}) \mid \text{Zn}^{2+}(\text{aq})$ Oxidant and reductant are in different phases. Metal electrode is part of redox couple.

OR

$\text{Fe}^{3+}(\text{aq}), \text{Fe}^{2+}(\text{aq}) \mid \text{Pt}$ Oxidant and reductant are in the same phase. An inert electrode is used.

The vertical line represents a phase boundary.

Equilibrium Constant, K

Constants will be dimensionless, ie have no units, in keeping with current IUPAC conventions. They will include:

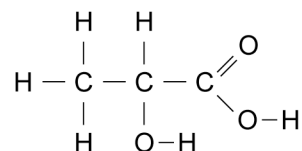
- K_c General equilibrium constant in which the equilibrium composition is expressed in terms of concentration of species
- K_a Acid association constant or acidity constant
- K_w Dissociation constant of water
- K_s Solubility product or solubility constant

– p notation will be restricted to: $\text{p}K_a$ for $-\log_{10} K_a$
and pH for $-\log_{10} [\text{H}_3\text{O}^+]$

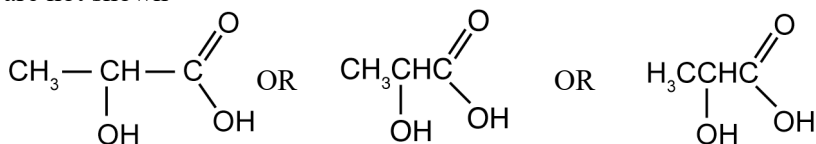
Organic Chemical Formulae

	Information conveyed	Example: lactic acid
empirical formula	Stoichiometric proportions of atoms only. Simplest ratio formula.	CH_2O
molecular formula	Formula of the actual molecule.	$\text{C}_3\text{H}_6\text{O}_3$
structural formula	Shows how atoms are connected. It may be drawn in different ways.	

(a) All atoms and bonds are shown.

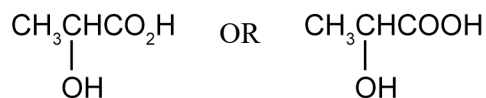


(b) Bonds to hydrogen are not shown

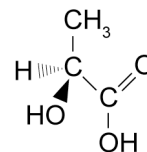


OR

Only bonds to substituents are shown.



(c) Stereochemistry (3-D arrangement of atoms) is shown.



The structural formulae in (b) are referred to as condensed structural formulae.

Organic Chemical Nomenclature

IUPAC conventions will be followed. There is ongoing discussion on some of the following naming. Candidates will be given full credit for alternative naming if an unambiguous structure is implied. Some examples are:

Structure	IUPAC name
$\begin{array}{c} \text{CH}_3 - \text{CH} - \text{CH}_2 - \text{CH}_2 - \text{CH}_3 \\ \\ \text{CH}_3 \end{array}$	2-methylpentane
$\begin{array}{c} \text{OH} \\ \\ \text{CH}_3 - \text{CH} - \text{CH} - \text{CH}_3 \\ \\ \text{CH}_3 \end{array}$	3-methylbutan-2-ol
$\begin{array}{c} \text{CH}_3 - \text{CH} - \text{CH}_2 - \text{C} - \text{OH} \\ \qquad \qquad \qquad \\ \text{CH}_2 \qquad \qquad \qquad \text{O} \\ \\ \text{CH}_3 \end{array}$	3-methylpentanoic acid
$\begin{array}{c} \text{O} \\ \\ \text{Br} - \text{CH}_2 - \text{CH} - \text{CH}_2 - \text{C} - \text{CH}_3 \\ \\ \text{Cl} \end{array}$	5-bromo-4-chloropentan-2-one
$\begin{array}{c} \text{CH}_3 - \text{CH}_2 - \text{C} - \text{O} - \text{CH}_2 - \text{CH}_3 \\ \\ \text{O} \end{array}$	ethyl propanoate
$\text{CH}_3 - \text{CH}_2 - \text{CH}_2 - \text{NH}_2$	propan-1-amine
$\begin{array}{c} \text{CH}_3 - \text{C} - \text{NH}_2 \\ \\ \text{O} \end{array}$	ethanamide

References

P Akins and L Jones, *Chemistry – Molecules, Matter and Change* (3rd edition), WH Freeman, 1997.
 Aylward & Findlay, *SI Chemical Data* (6th Edition), John Wiley & Sons, Australia, 2008.