With 2019 NCEA Exam included

2020 Version

Chemistry AS 91164
C2.4 Bonding, Structure and Energy

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AS 91164
C2.4

Bonding and structure are limited to:

❑ Lewis structures, shape and polarity of simple molecules.

❑ Intermolecular forces (the distinction between the different types of intermolecular forces is not required)

❑ Ionic, covalent and metallic bonding

❑ Molecular, ionic, metallic and covalent network substances

❑ Properties are limited to hardness (including malleability and ductility), electrical conductivity, melting and boiling points and solubility.

Energy changes are limited to:

❑ exothermic and endothermic reactions including energy (enthalpy) changes associated with differing amounts of substances and changes of state and enthalpy changes associated with the making and breaking of chemical bonds
❑ calculations of energy changes using $\Delta_r H$ and reaction stoichiometry, and bond enthalpy

Chemistry is the study of matter and energy and the interaction between them. The elements are the building blocks of all types of matter in the universe. Each element is made up of only one type of atom, each with its specific number of protons known as its atomic number.
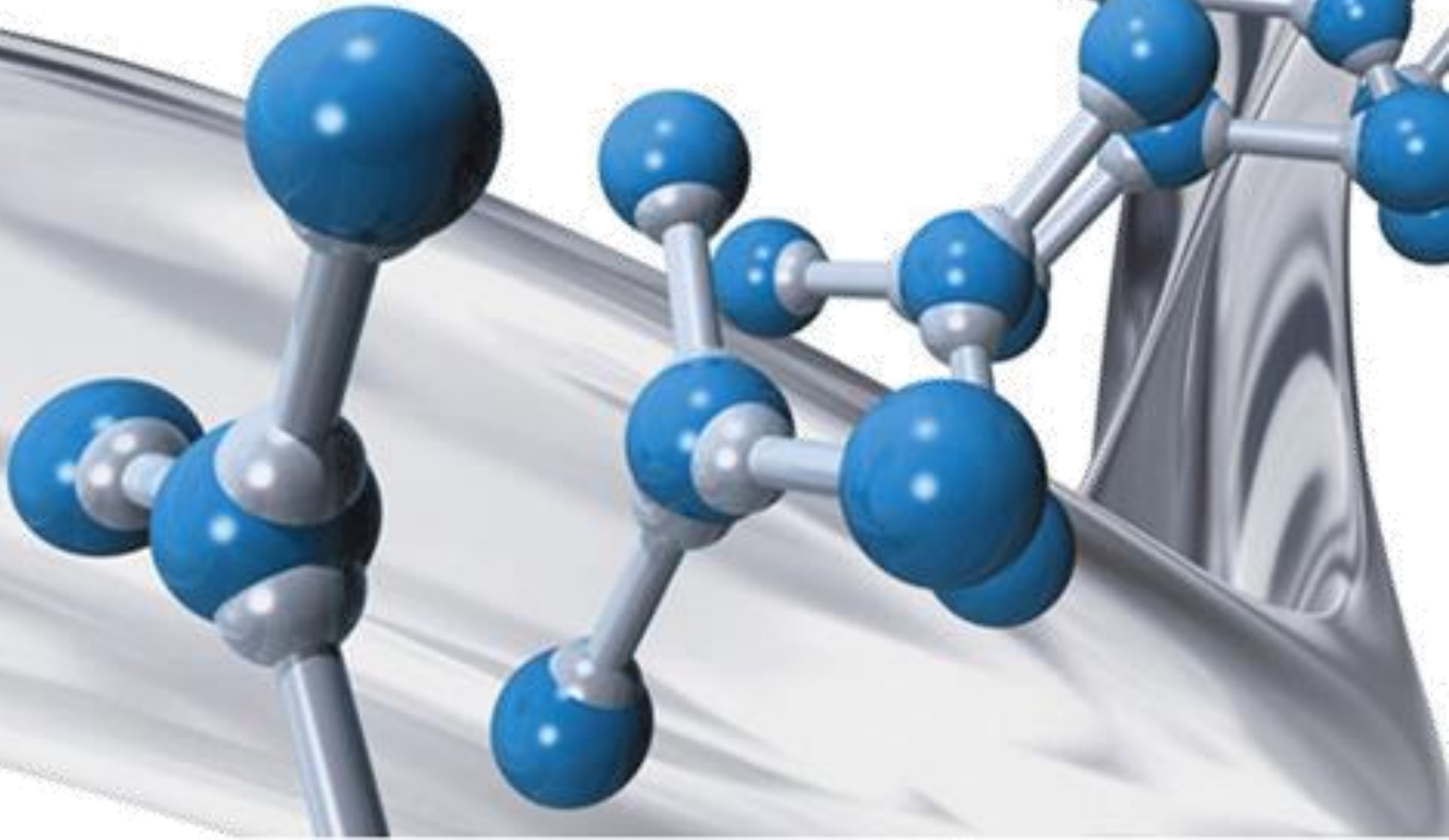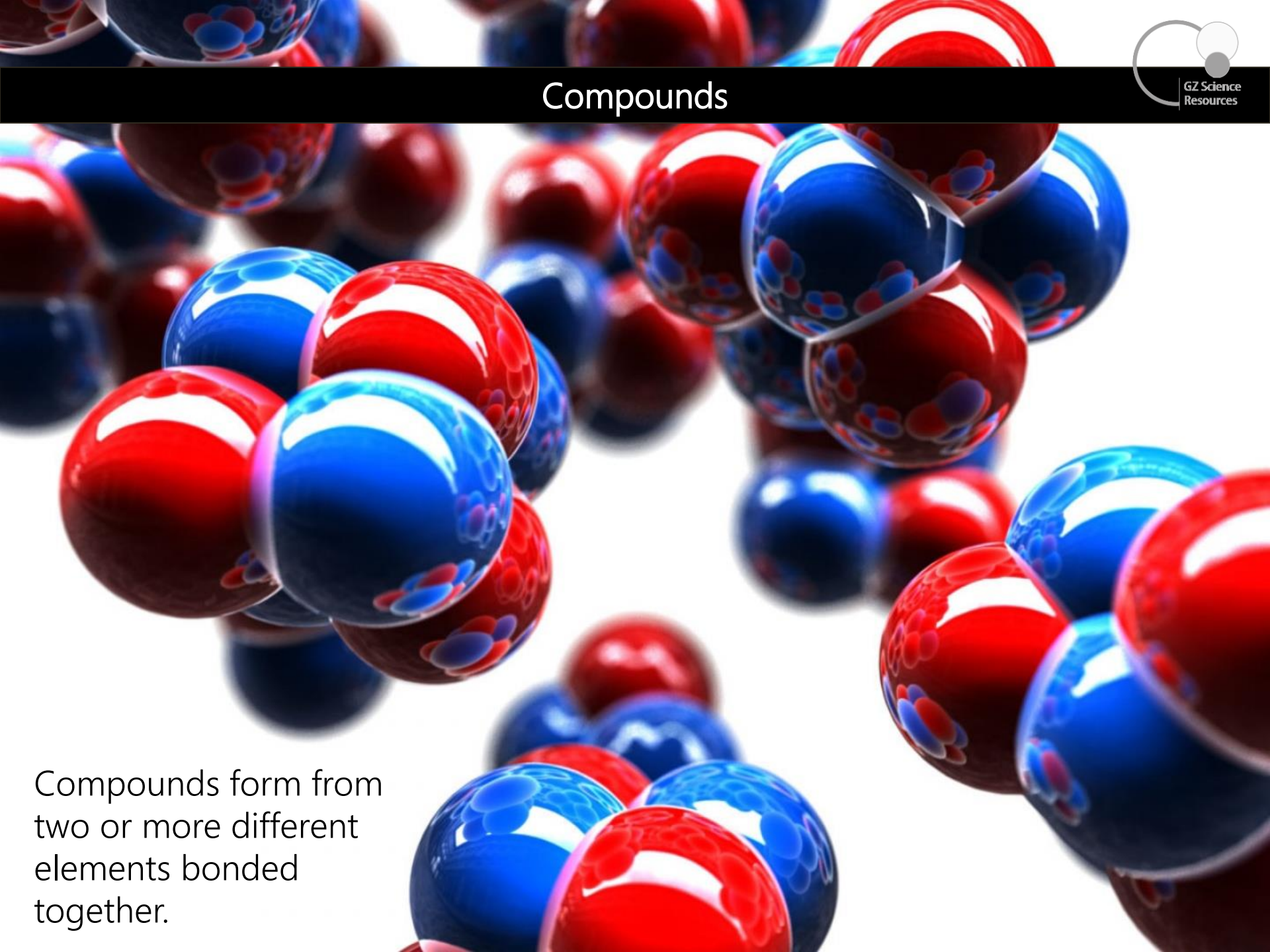
A large amount of energy is required to break an atom down into smaller particles. The elements occur in widely varying quantities on earth. The ten most abundant elements make up 98% of the mass of earth. Many elements occur only in traces, and a few elements are synthetic and highly unstable.

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Compounds form from two or more different elements bonded together.

Background Knowledge
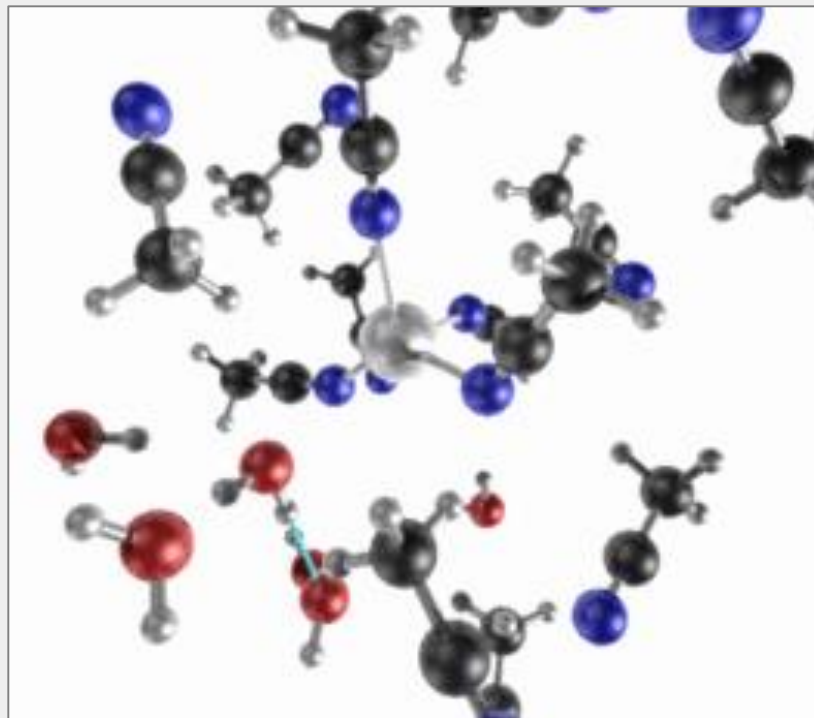
The compounds are often more stable than the elements they originated from and may release this extra energy in the form of heat and/or light when bonding together.

There are two main types of bonding holding atoms together in a compound; Ionic and Covalent. (we also discuss metallic bonding in detail later)

Ionic Bonding holds together oppositely charged ions due to electrostatic attraction. This type of bonding occurs when **metal** and **non-metal atoms** react and there is a complete **transfer of electrons** to form negative (anion) and positive (cation) ions.

The ions then combine in a set ratio to form a neutral compound (usually) with negative and positive charges balanced out.
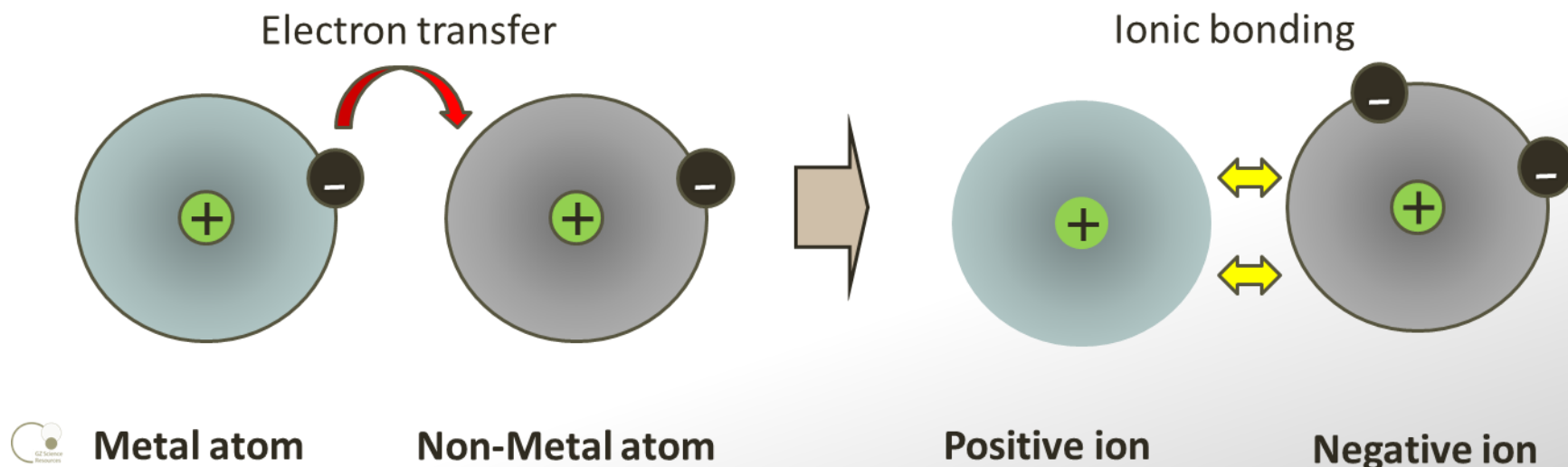
Electron transfer   Ionic bonding

Metal atom   **Non-Metal atom**   **Positive ion**   **Negative ion**

Some compounds are **ionic compounds**, since they are made up of cations and anions.
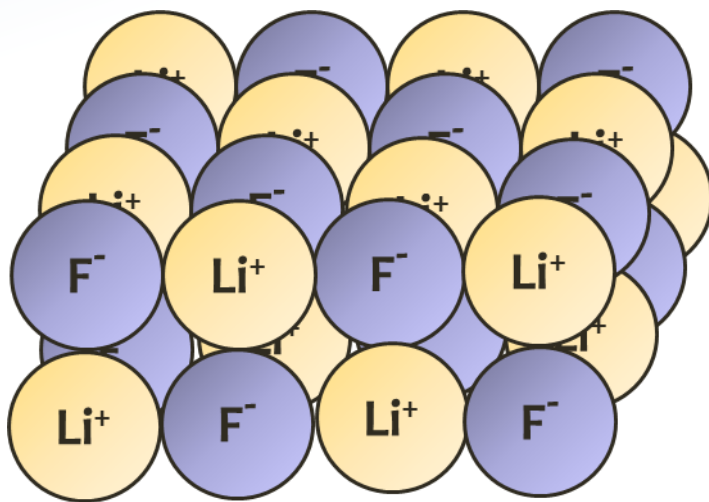
A model of an ionic compound in solid state

The Anion (F^-) takes an electron from the Cation (Li^+) so their outer energy levels have a stable 8 electrons each.

Anions and cations have a strong <u>electrostatic attraction</u> for each other so they bond together as a compound.

Compounds are neutral substances. For ionic compounds, the charges of the positive ions are balanced by the charges of the negative ions.

Metallic solids also involve strong electrostatic attraction, but between ions and free moving valence electrons (this will be explained in detail later)

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Covalent Bonding occurs when electrons are 'shared' between neighbouring atoms. This often occurs when two or more **non-metals** react. No ions are formed and there is no transfer of electrons. The compound formed is neutral with no charge.

Non-Metal atom   Non-Metal atom
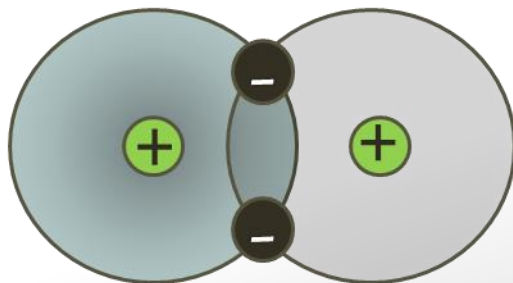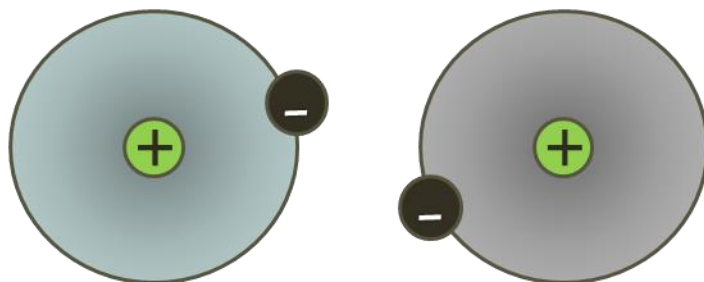
covalent bonding

The valance electrons (electrons in outside energy level) are involved in bonding. These electrons orbit in pairs. The negative charge of the electron pair will attract the positively positive nucleus of other atoms, and this holds the atoms together in a molecule.

Covalent bonding occurs in both **molecular solids**, just within the molecules, and in **covalent network solids**, between each and every individual atom in a solid.

All covalent bonds are strong. That is, they require a large amount of energy to 'break' the bond. However, some covalent bonds are stronger than others. The greater the overlap of valence orbitals (the area the valence electrons orbit the nucleus) the stronger the bond.


Bonding atoms

Overlap between valence orbitals

(a) H–F versus H–I   (b) F–F versus I–I
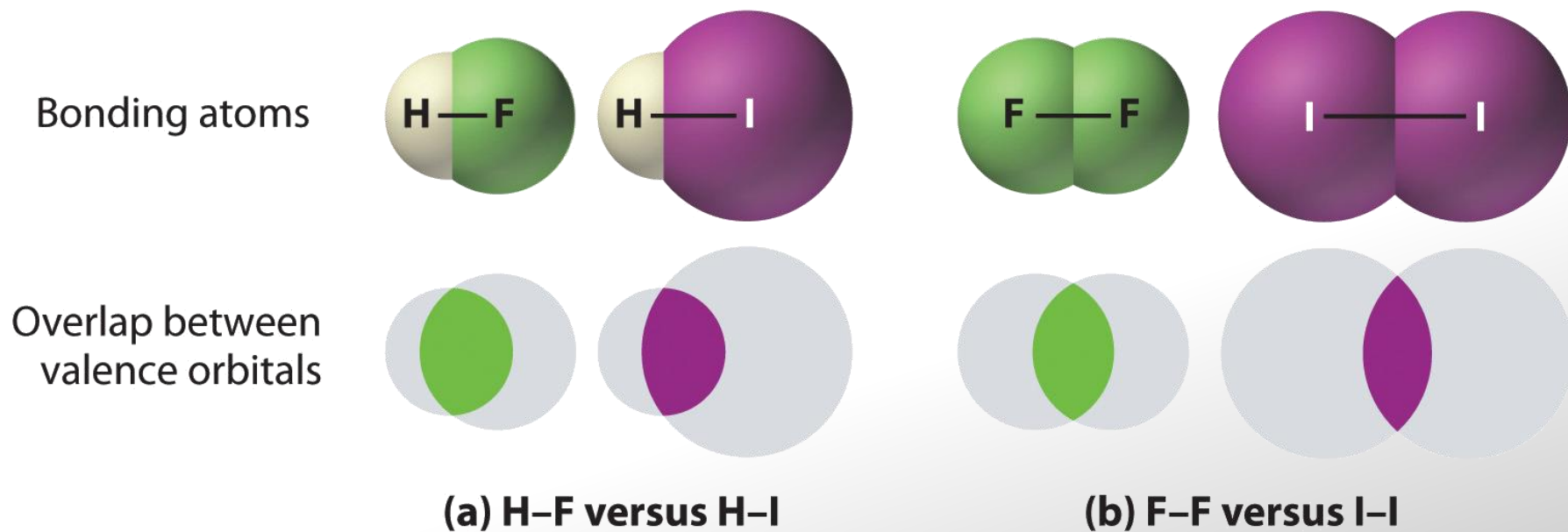
The electron-pair must lie between the nuclei for the attraction to outweigh the repulsion of the two nuclei. This 'sharing' of electrons between atoms creates a covalent bond – giving both atoms the stability of a full outer shell.
Covalent bonds are normally formed between pairs of non-metallic atoms.

Only one pair of electrons holding the nuclei together
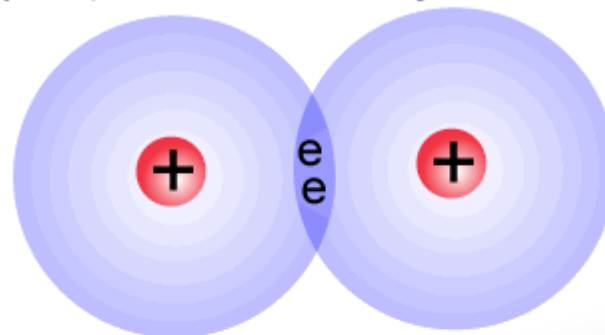
Some covalent bonds involve only one pair of electrons and are known as single bonds. Other covalent bonds can involve two pairs of electrons; double bonds and three pairs of electrons; triple bonds.

Two pair of electrons hold the nuclei tighter and closer

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Background
Knowledge
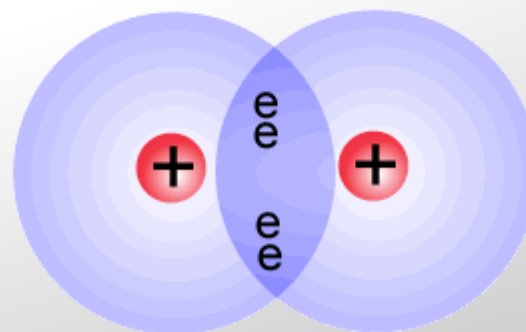
Lavoisier – French Chemist 1789

Lavoisier devised a system of naming compounds based on their chemical composition. If the compound is formed between a Metal **cation** (+ve) and a Non-Metal **anion** (-ve), then the compound name joins the two names together with the <u>metal name first</u>. Names of the ions need to be remembered.
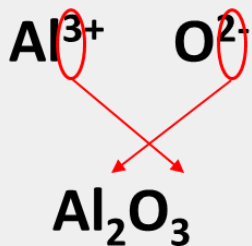
Sodium

+   ⟶   **Sodium hydroxide**

hydroxide

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1. Write down the ions (with charges) that react to form the compound. Cation comes before Anion.

Al^{3+} O^{2-}

Al_2O_3

2. Cross and drop the charge numbers.

3. Place brackets around a compound ion.

4. If the numbers are both the same remove.

5. If any of the numbers are a 1 they are removed

6. Remove any brackets if not followed by a number

H^+ SO_4^{-2}

$H_2(SO_4)_1$ $\longrightarrow$ H_2SO_4
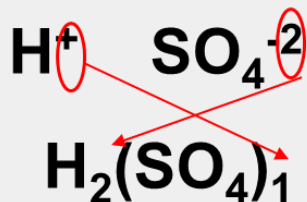
G Lewis – American Chemist 1916

G Lewis devised a system of drawing covalent molecules showing arrangement of atoms and valence electrons – both those involved in bonding and those that are not (called non-bonding pairs). Electrons in inner shells are <u>not</u> involved in bonding. These diagrams are called **Lewis structures**. (or diagrams) The Lewis structure is drawn so that each atom has eight electrons associated with it (except for hydrogen, which has two). This is the *octet rule*.
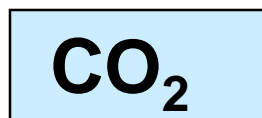
Lewis structure of H_2O (water)

Hydrogen electron

Bonded pair

H

O

H

Oxygen electron

Lone pair
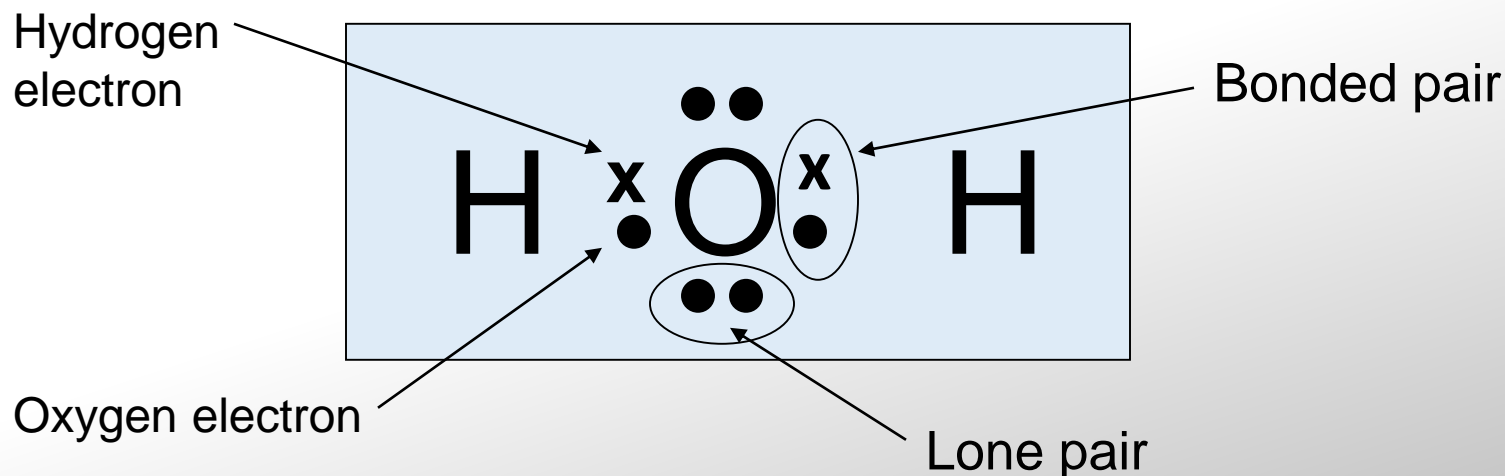
1. Calculate valence electrons of all atoms. If the molecule is an ion, then subtract the charge from the total electrons and place the charge outside of square brackets of the Lewis structure at the end. *Example carbon dioxide.*
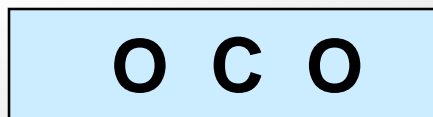
C = 4

O = 6

O = 6

16

CO_2

2. Write down number of pairs of electrons.

16 / 2 = 8 pairs

3. Place atom with least filled valence shell in the centre with the other atoms arranged around the outside (periphery)

O   C   O

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4. Bond all atoms together (either x● or x x   or ——   = one pair of electrons)

O —— C —— O
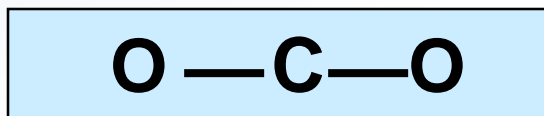
8 pairs – 2 pairs = 6 pairs remaining

5. Place remaining e- pairs around the periphery atoms so each has 4 pairs (including bond pair) around it.
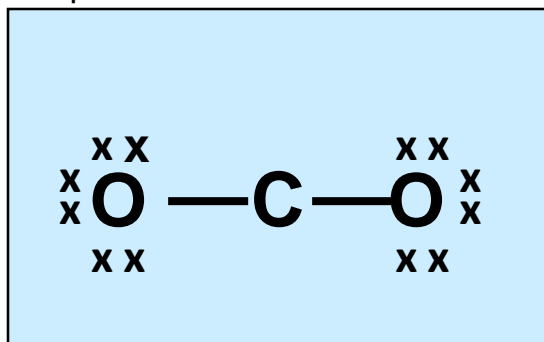
6 pairs – 6 pairs = 0 pairs remaining

6. If there any remaining pairs place them around the outside of the central atom.

7. Rearrange lone pairs (pairs not bonded) into bonded pairs if the central atom does not have 4 pairs around it.

❑ The number of covalent bonds an atom forms is called its *valence*.
❑ Some atoms have *fixed valence*. E.g.: H = 1, C = 4, F = 1. (most halogens = 1)
❑ Some atoms have *variable valence*. For example:
    O = 2 (sometimes 3), B, N = 3 (sometimes 4).
❑ An atom bonded to only one other atom is *peripheral* (monovalent atoms such as H and F are always peripheral).
❑ An atom bonded to two or more other atoms is *central*.
❑ Often, the formula is written to indicate connectivity. For example: HCN = H bonded to C, C bonded to N, H and N are not bonded.

extra info
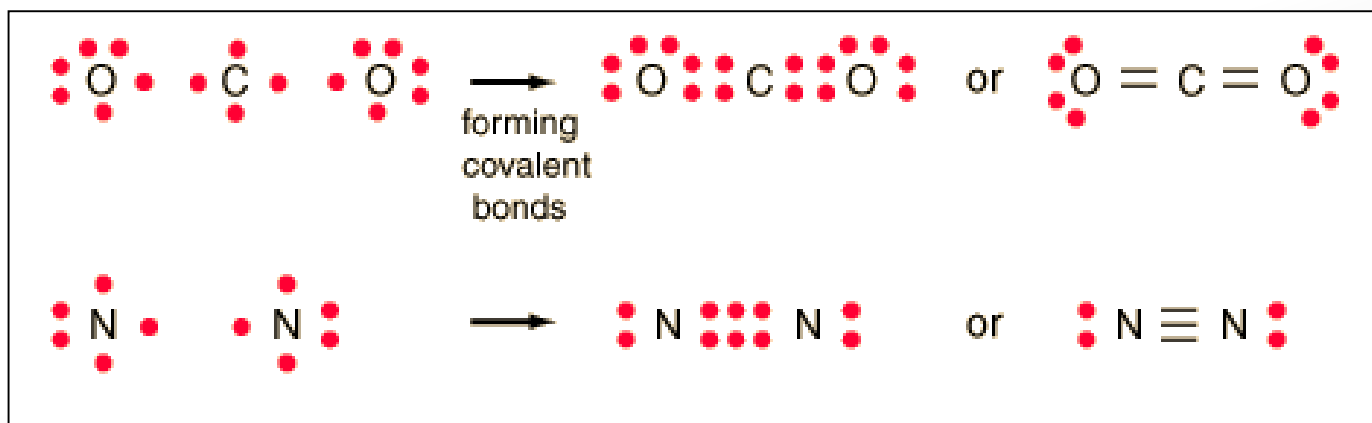
If there are extra non-bonding pairs of electrons left after all of the periphery atoms are filled in accordance with the *octet rule* then they are placed around the central atom(s) according to the **Rule of Orbitals.**

The Rule of Orbitals: the total number of non-bonding pairs and bond pairs (LP+BP) associated with an atom cannot exceed the number of Valence Shell Orbitals (VSO = n^2, where n is the row of the Periodic Table in which that atom resides).
n = 1 (H): <u>maximum</u> VSE pairs (LP+BP) = VSO = 1;
n = 2 (B, C, N, O, F): <u>maximum</u> VSE pairs (LP+BP) = VSO = 4 ("octet rule")
n = 3 ((Al, Si, P, S, Cl): <u>maximum</u> VSE pairs (LP+BP) = VSO = 9 etc.

Boron and Beryllium often are found with only **3** (or 2) non-bonded + bonded pairs of electrons around them

Sidgwick and Powell – 1940

Sidgewick and Powell devised a theory to predict the shapes molecule formed. It is based on the following ideas:

Electron pairs form regions of negative charge

Negative charges repel each other

Regions of negative charge will be spaced as far apart as possible around a central atom.
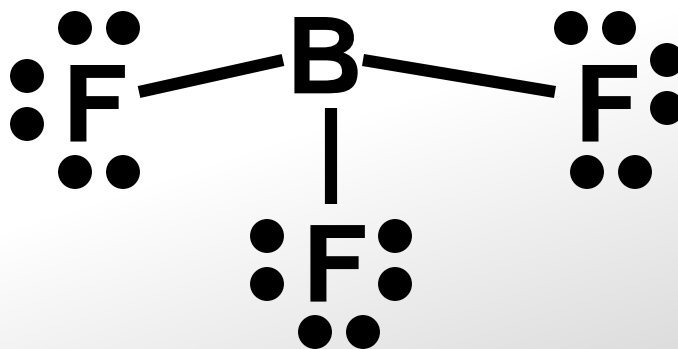
This theory is called **valence-shell electron-pair repulsion (VSEPR) theory**

The shapes of molecules are determined by the way the regions of negative charge are arranged around the central atom in the molecule. A region may consist of one non-bonded pair of electrons or one bonded pair or two bonded pairs or three bonded pairs. All of these electron arrangements occupy the same region of space.
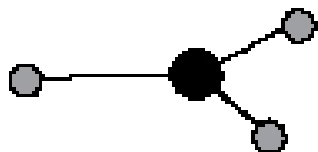
Three pairs
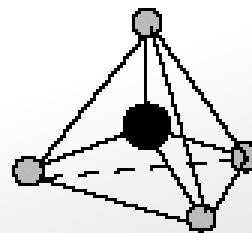
Four pairs

Since regions of electrons are negatively charged, they repel each other as far apart as possible (**VSEPR**) into a **linear geometry / arrangement**. Two negative regions arrange themselves on opposite sides of the central atom.
The bond angle will be 180°.

The shape name is **linear.**
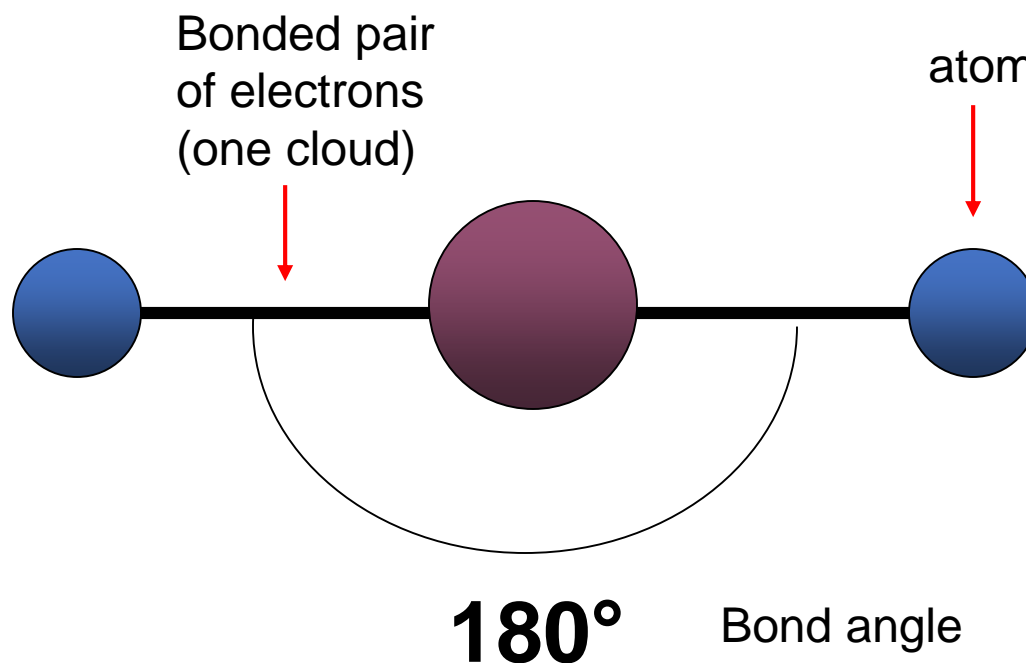
Bonded pair
of electrons
(one cloud)

atom

180°   Bond angle

Three regions of negative charge will cause a bond angle of 120° as they repel each other into a **trigonal planar geometry / arrangement.**
All the atoms still lie on a flat plane (like a sheet of paper).
As there are no non-bonded pairs of electrons, the final shape is also a **trigonal planar**. (or triangular planar)
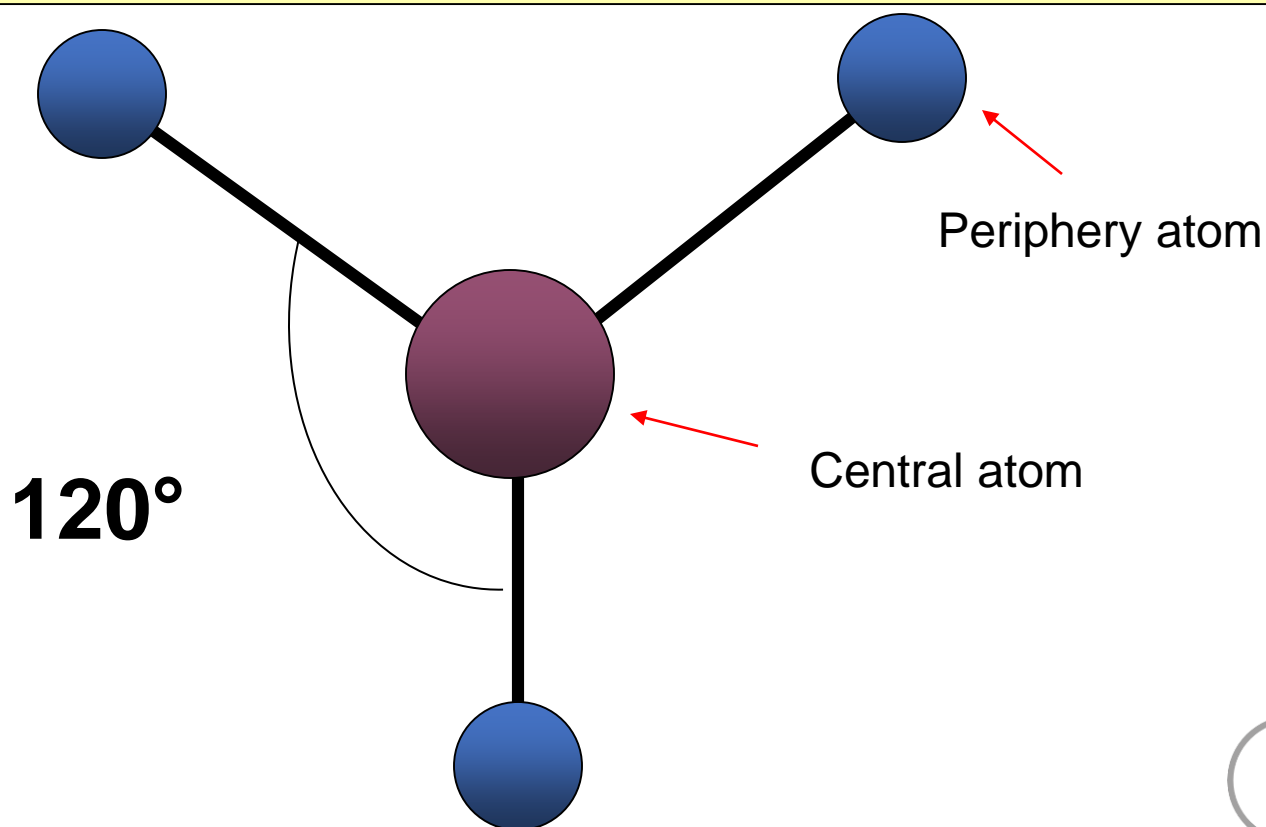
Periphery atom

Central atom

120°

The regions of negative charge repel to a **trigonal planar geometry.** The bond angle between the remaining pairs is approximately 120°. As there is one non-bonded pair of electrons, the final shape formed by the atoms is called **bent**.

Non-bonded pair

Bonded pair

~120°

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i extra info
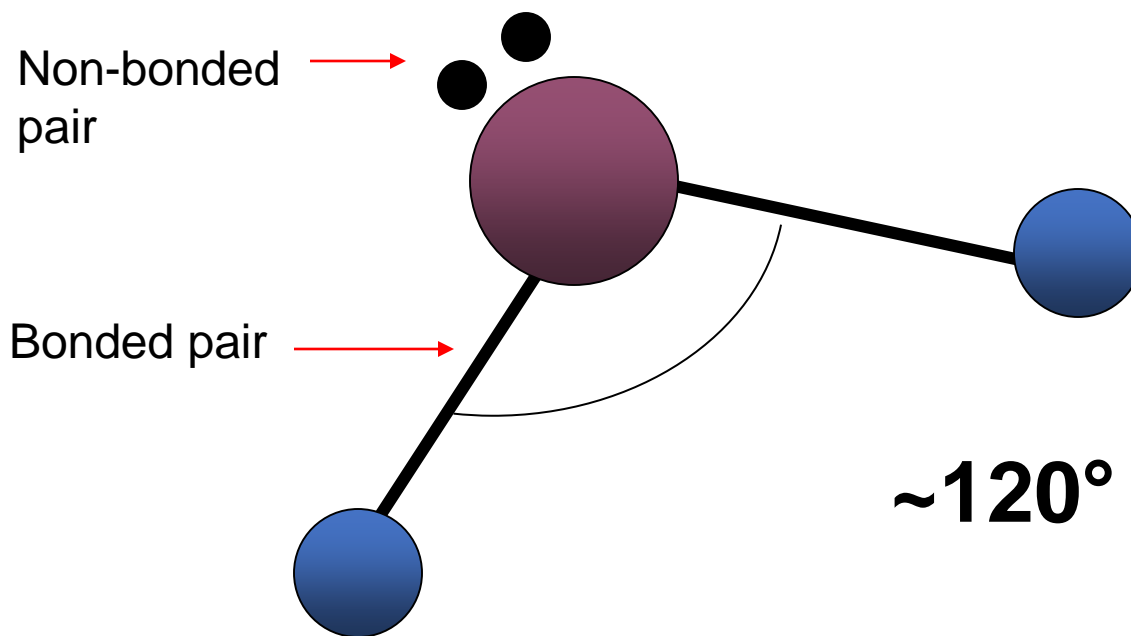
When one of the regions of electrons is a *non-bonding pair* it will have a slightly greater push to the bonded pairs. This is because the non-bonding pair is only orbiting around one positive nucleus and their negative charge is less 'neutralised' than if they had another nucleus to orbit around.

When four regions of negative charge are around a central atom, they repel each other into a 3-dimensional **tetrahedral geometry / arrangement.** The bond angle is now 109.5°. This is because it is a 3-dimensional sphere divided into 4 rather than a circle.

This final shape is **tetrahedral.**

~109.5°

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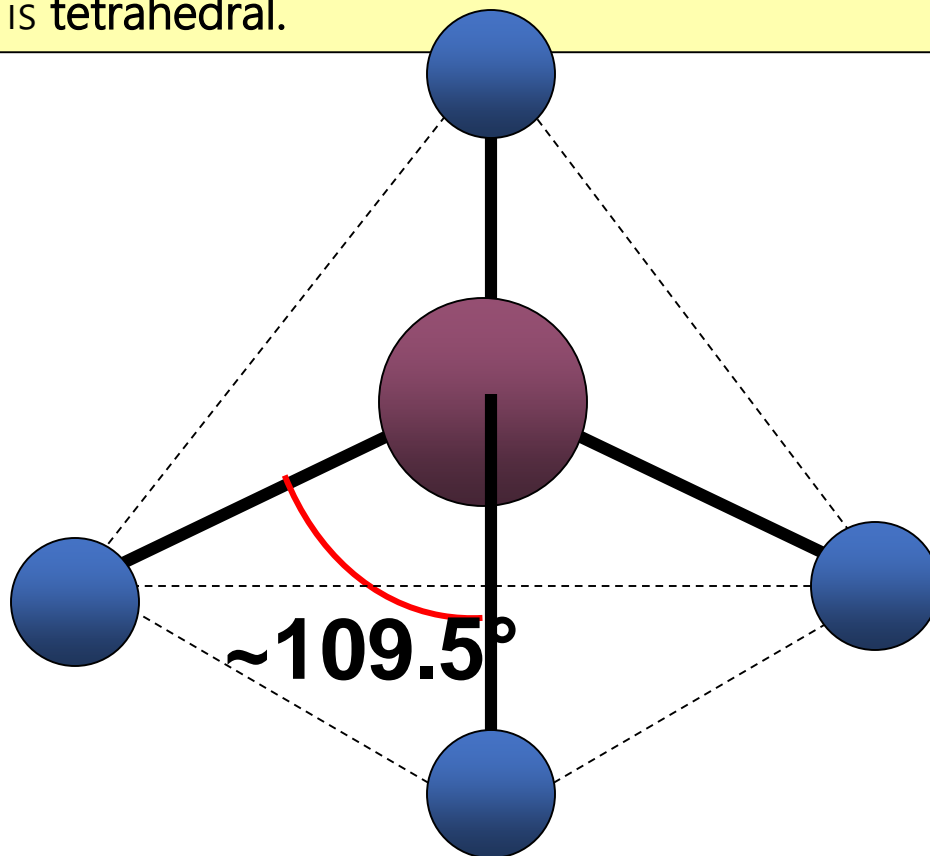
The four regions of negative charge still occupy a 3-dimensional **tetrahedral geometry / arrangement.** (The lone pair, however, exerts a stronger repulsion to the remaining bonded pairs). The bond angle is approximately 109.5°.
The final shape the bonded atoms form, due to one non-bonded pair, is a **trigonal pyramid** (or a triangular pyramid)

Non-bonding pair

~109.5°
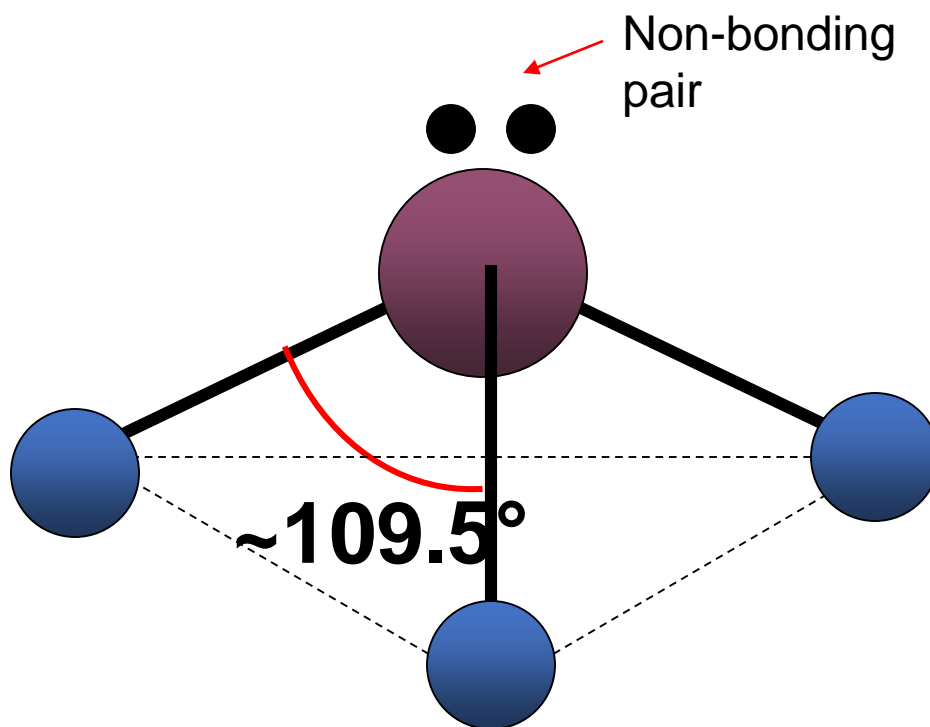
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The 4 regions of negative charge repel each other to a (warped) **tetrahedral geometry / arrangement.** However, the two non-bonding pairs create a much stronger repulsion than one non-bonding pair and the bond angle between the remaining bonded pairs is smaller again at approximately 109.5° (compared to 120° of the bent shape with only 3 regions of negative charge). The final shape the bonded atoms form is called **Bent**.

Non-bonding pairs

~109.5°

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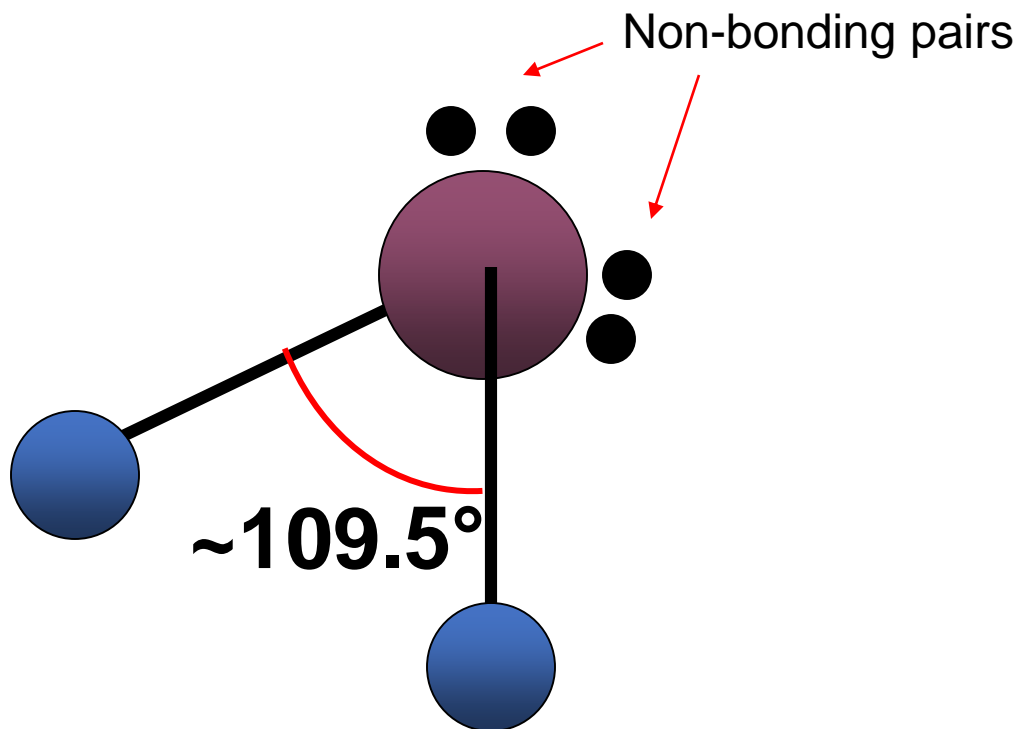
Determining Molecular Shapes

Electron regions of negative charge (bonding or non-bonding pairs)

2 regions

3 regions

4 regions

No non-bonding pair
Linear
180°
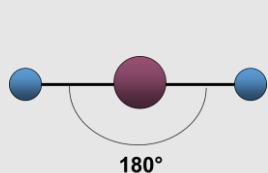
No non-bonding pair
Trigonal Planar
120°
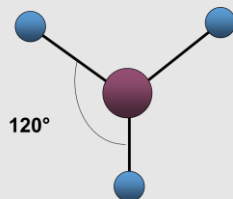
1 non-bonding pair
Bent
120°
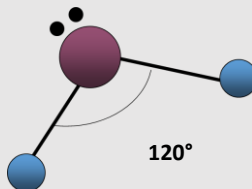
No non-bonding pair
Tetrahedral
109.5°
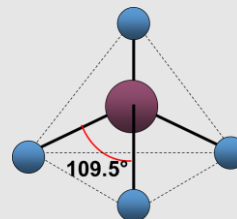
1 non-bonding pair
Trigonal Pyramid
Lone pair
109.5°
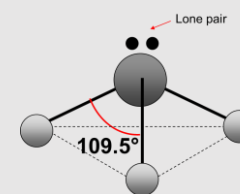
2 non-bonding pairs
Bent
109.5°

CO_2
180°

BF_3
120°

SO_2
120°

CH_4
~109.5°

NH_3
~109.5°

H_2O
~109.5°

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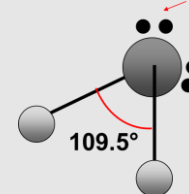
Note: make sure the question asks you to draw a shape and **not** a Lewis structure

1. Atom on **same plane** as central atom – straight solid line

2. Atom **receding** from central atom – lines starting large and getting smaller

3. Atom **approaching** from central atom – solid triangle starting small and getting larger

approaching

receding

Same plane
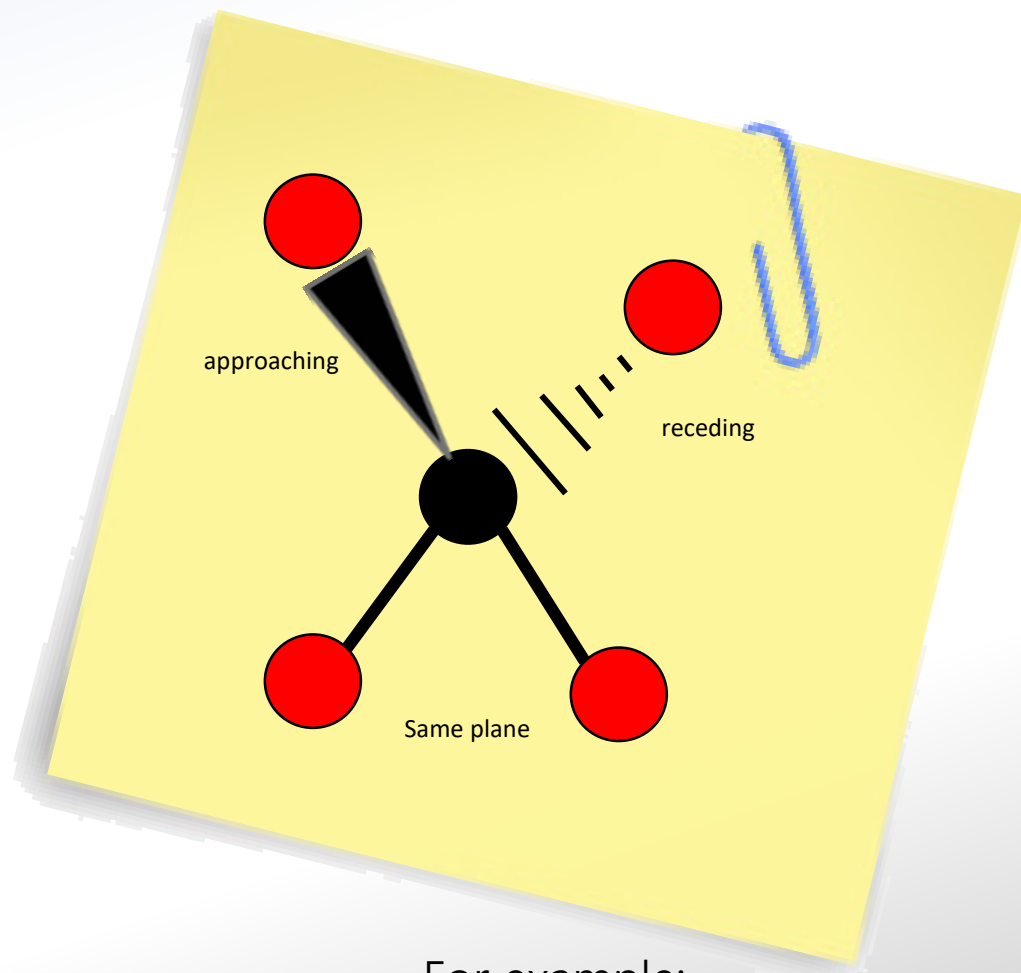
For example:
Tetrahedral shape

Explain why the shape of the CO_2 molecule is linear but the shape of H_2O is bent?

1. The C (central atom) of CO_2 has 2 **regions of negative charge** around it in the form of double bonds connected to an O atom. (draw Lewis structure)
2. Each of the regions of **negative charge repel each other the furthest away** from each other in 3 dimensional space into a linear **geometry / arrangement**
3. There are no non-bonding pairs so the CO_2 molecule therefore also **forms a *linear shape***

1. The O molecule (central atom) of H_2O has 4 **regions of negative charge** around it in the form of two single bonds connected to a H atom and two non-bonding pairs. (draw Lewis structure)
2. Each of the regions of **negative charge repel each other the furthest away** from each other in 3 dimensional space and form a tetrahedral **geometry / arrangement**.
3. However with only 2 of the regions bonded to atoms the shape the H_2O molecule **forms is a *bent shape*.**

Question 1b (i) : The Lewis structure for a molecule containing atoms of boron, oxygen, and hydrogen, is shown below.
The following table describes the **shapes around two of the atoms** in the molecule above.
Complete the table with the approximate bond angles *x* and *y*.
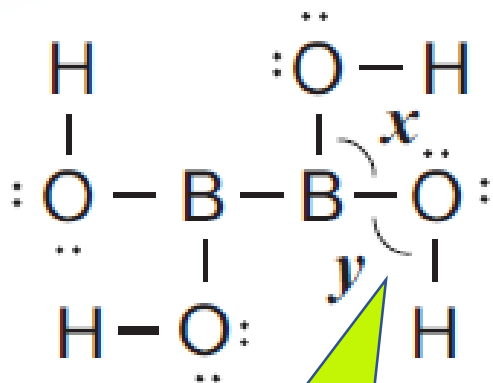
Central atom	Shape formed by bonds around the central atom	Approximate bond angle
B	trigonal planar	$x =$ 120°
O	bent	$y =$ 109.5°

Even though this is an unknown molecule focus on each atom and the number of regions of charge around each

Bond angles

2 regions = 180°

3 regions = 120°

4 regions = 109.5°

Question 1b (ii): The bond angles x and y in the molecule above are different. Elaborate on why the bond angles are different.
In your answer you should include:
• factors which determine the shape around the:
- B atom for bond angle x
- O atom for bond angle y
• reference to the arrangement of electrons around the B and O atoms.
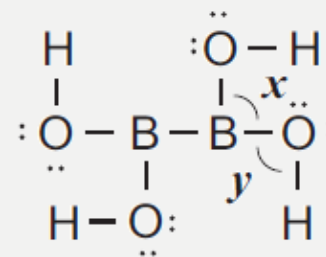
Answer 1b (ii): The B atom has **three regions of electron density** around it. These are all bonding regions. The regions of electron density are arranged to minimise repulsion / are arranged as **far apart as possible** from each other in a **trigonal planar** arrangement. This is why the bond angle is 120°

The O atom has **four regions of electron density** around it. The regions of electron density are arranged to minimise repulsion / are arranged **as far apart** as possible from each other in a **tetrahedral** arrangement / two of these are bonding (and two are non-bonding). This is why the bond angle is 109.5°.

Achieved Question

Q 1a: Draw the Lewis Structure for each of the following molecules

Draw the Lewis structure (electron dot diagram) for each of the following molecules.

Molecule	O_2	OCl_2	CH_2O
Lewis structure	:O=O:	:Cl—O—Cl:	O (double bond) H—C—H

Look further through the exam paper for hints on drawing

In the past the Lewis structure is a "stand alone" question – which case there is an Achieved point given for the majority correct.

In past years the following molecules have been used:

HCN CH_2Br_2 AsH_3 CH_4 H_2O N_2 PCl_3 CO_2 H_2S

These have one central atom (except N_2) and 2 or more outside attached atoms

Question 1b: Carbon atoms can bond with different atoms to form many different compounds. The following table shows the Lewis structure for two molecules containing carbon as the central atom, CCl_4 and $COCl_2$. These molecules have different bond angles and shapes.

Evaluate the Lewis structure of each molecule to determine why they have different bond angles and shapes.

In your answer you should include:
- The approximate bond angle in each molecule
- The shape of each molecule
- Factors that determine the shape and bond angle for each molecule.

Lewis structures are given in the shape and bond angle questions

Make sure you clearly state the correct shape and bond angle **beside each Lewis structure** - before starting the discussion. Use the formula as the "name" for each molecule – you do not have to know the name unless it is given.

Molecule	CCl_4	$COCl_2$
Lewis structure		

Excellence Question

Answer 1b: In each CCl_4 molecule, there are four negative electron clouds / regions around the central C atom. These repel each other as far away from each other as possible in a tetrahedral (base) arrangement, resulting in a 109.5° bond angle. All of these regions of electrons are bonding, without any non-bonding regions, so the shape of the molecule is tetrahedral.

In each $COCl_2$ molecule, there are three negative electron clouds / regions around the central C atom. These repel as far away from each other as possible in a triangular / trigonal planar (base) shape, resulting in a 120° bond angle. All of these regions of electrons are bonding, without any non-bonding regions, so the shape of the molecule is trigonal planar.

Bond angles

2 regions = 180°

3 regions = 120°

4 regions = 109.5°

Discuss each molecule separately using the same steps

❏ State number of electron regions around central atom
❏ State base shape they repel to and the angle
❏ State number of bonding/non-bonding regions
❏ State final shape (will be the same as base shape if no lone pairs)

Question 3a (i) : Draw the Lewis structure (electron dot diagram) for each of the following molecules, and name their shapes.

Molecule	H_2O	CS_2	PH_3
Lewis structure	H – Ö – H	S̈=C=S̈	H – P̈ – H | H
Name of shape	bent	linear	Trigonal pyramid
Approximate bond angle around the central atom	109.5°	180°	109.5°

Question 3a (ii): Compare and contrast the shapes and bond angles of H_2O, CS_2 and PH_3.

Answer 3a (ii): Bond angle is determined by the number of electron clouds / areas of negative charge around the central atom, which are arranged to minimise repulsion / are arranged as far apart from each other as possible (maximum separation).

Both H_2O and PH_3 have 4 electron clouds / areas of negative charge around the central atom, so the bond angle is that of a tetrahedral arrangement of 109.5°, whereas there are only 2 electron clouds / areas of negative charge around the central atom in CS_2, which means minimum repulsion is at 180°, resulting in CS_2's shape being linear.
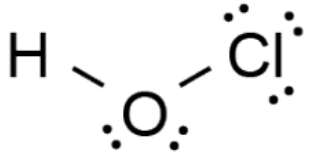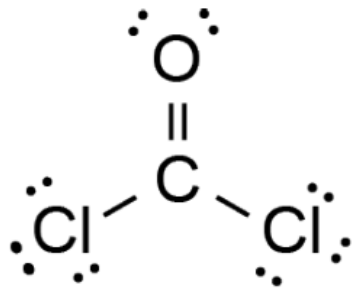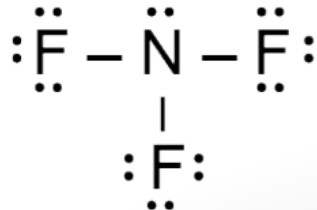
The shapes of H_2O and PH_3 differ despite having the same tetrahedral arrangement because water has two non-bonding pairs of electrons around the central atom, while phosphine only has one non-bonding pair. The resulting shapes are bent or v-shaped for H_2O, while PH_3 is trigonal pyramid
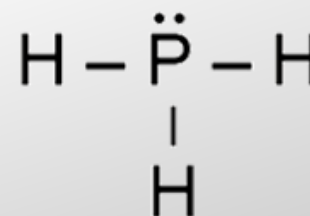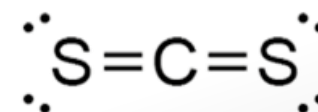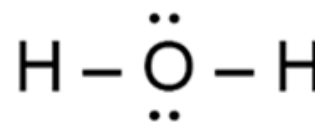
$$H - \ddot{O} - H$$

$$\ddot{S} = C = \ddot{S}$$

$$H - \ddot{P} - H$$
$$|$$
$$H$$

Question 2a (i) : Draw the Lewis structure (electron dot diagram) for each of the following molecules, and name their shapes.

Molecule	HOCl	COCl$_2$	NF$_3$
Lewis structure			
Name of shape	bent / v-shaped	trigonal planar	trigonal pyramid
Approximate bond angle around the central atom	109.5°	120°	109.5°

Question 2a (ii): Justify the shapes and bond angles of HOCl and $COCl_2$.

Answer 2a (ii): Bond angle is determined by the **number of electron density regions** around the central atom, which are arranged into a position to minimise repulsion / are **arranged as far apart from each other as possible** (maximum separation).

HOCl has 4 electron density regions / areas of negative charge around the central O atom.

This means the electron density regions around the central atom is arranged with maximum separation in a **tetrahedral shape** with a bond angle of 109.5°, to minimise (electron-electron) repulsion. Due to the presence of two non-bonding pairs of electrons / regions (or two bonding regions) on the central O atom, HOCl has an actual shape that is **bent** / v-shaped / angular.

$COCl_2$ **has only 3 electron** density regions / areas of negative charge around its central C atom so the electron density regions around the central atom is arranged with maximum separation in a **trigonal planar shape** with a bond angle of 120°, to minimise (electron-electron) repulsion. Since $COCl_2$ has only bonding electron pairs (no non-bonding pairs) on its central atom, the actual shape is **trigonal planar** (with bond angles of 120°).

Question 2a. Draw the Lewis structure (electron dot diagram) for each of the following molecules, and name their shapes.

Molecule	H_2S	NH_3	BF_3
Lewis Structure	H – S̈ – H	H | H – N – H	:F̈ – B – F̈: | :F̈:
Name of Shape	Bent	Trigonal Pyramid	Trigonal Planar
Approximate bond angle around central atom	109.5°	109.5°	120°

Excellence
Question

Question 2b. Compare and contrast the shapes and bond angles of NH_3 and BF_3.

NH_3 has **four electron clouds / regions of negative charge** around its central N atom. As the electron clouds maximise separation to minimise repulsion they take a **tetrahedral geometry** with a **bond angle of 109.5°**. Three of the regions are bonded and one is non-bonded, so the **overall shape is trigonal pyramid.**
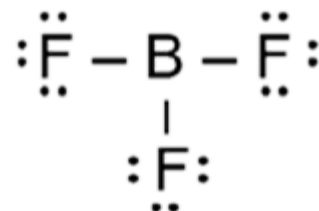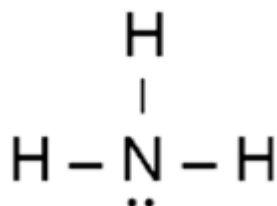
In contrast, BF_3 only has **three regions of negative charge** around its central B atom. As the electron clouds maximise separation to minimise repulsion they take a **trigonal planar geometry** with the bond angle of 120°. While BF_3 has three bonded regions like NH_3, because there is no non-bonding regions BF_3's **shape is trigonal planar.**

So although both molecules have three bonded areas to the central atom, ammonia has a fourth region of negative charge, which is not bonded. This affects its angle and shape.
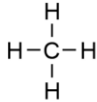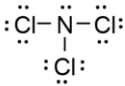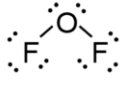
Question 2a. (i) Draw the Lewis structure (electron dot diagram) for the following molecules, and name their shapes.

Molecule	CH_4	NCl_3	OF_2
Lewis structure			
Name of shape	tetrahedral	Trigonal pyramid	bent

Question 2a. (ii) The above molecules have different shapes; however each molecule has an approximate bond angle of 109.5°.
Justify this statement by referring to the factors that determine the shape of each molecule.

Molecule	CH_4	NCl_3	OF_2
Lewis structure	H−C−H with H above and H below	:Cl−N−Cl: with :Cl: below	O with F and F
Name of shape	tetrahedral	Trigonal pyramid	bent

Bond angle is determined by the number of electron density regions around the central atom, which are arranged into a position to minimise repulsion by having maximum separation.
All molecules have 4 electron density regions / areas of negative charge around the central atom which arrange with maximum separation into a tetrahedral shape / geometry with a bond angle of (approx.) 109.5° / 109°.
In CH_4 all of the electron pairs are bonded, and so the shape of the molecule is also tetrahedral.
In NCl_3 three of the electron pairs are bonded and one is non-bonding. The observed shape of the molecule is trigonal pyramidal.
In OF_2, due to the presence of two non-bonding pairs of electrons / regions (or two bonding regions) on the central atom, OF_2 has an observed shape that is bent / v-shaped / angular.

Electronegativity is the attraction that an atom has towards electrons from another bonded atom. The greater the electronegativity the stronger the pull the nucleus of the atom has towards other electrons.

Trends in the periodic table
❑ The larger the nucleus (with the positive protons) the stronger the electronegativity, this means it increases from left to right.

❑ The further the valence electrons are from the nucleus the less the electronegativity, therefore the electronegativity decreases down a group.

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Resources

extra info
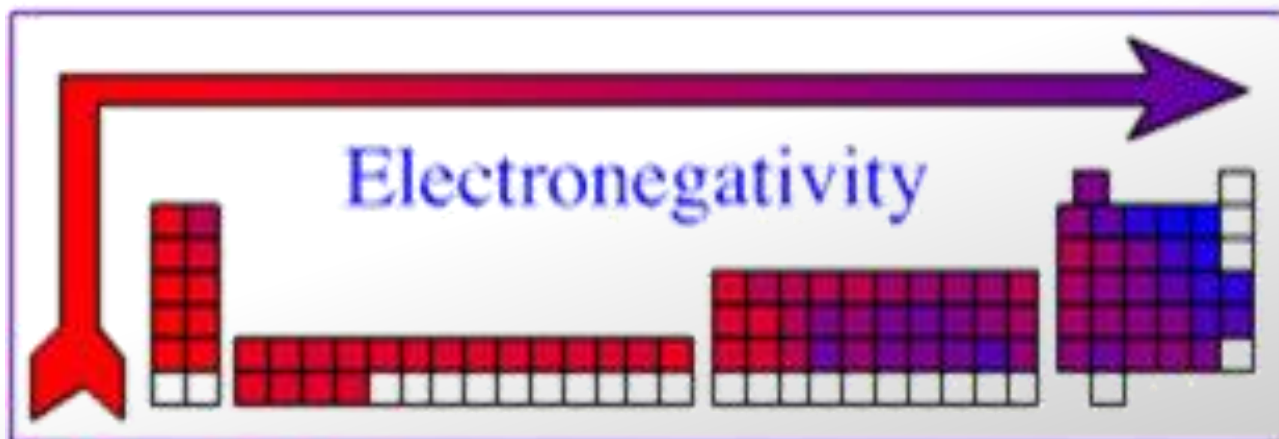
We use a **Pauling scale** to determine electronegativity. The scale starts close to 0 – with minimal electronegativity and goes up to 4 with the highest electronegativity. Most of the Inert gases do not have a value because of there no reactivity with other atoms.

Electronegativity

0.7   4

Pauling scale

1																		18
H	2												13	14	15	16	17	He
2.1																		--
Li	Be												B	C	N	O	F	Ne
1.0	1.5	3	4	5	6	7	8	9	10	11	12		2.0	2.5	3.0	3.5	4.0	--
Na	Mg												Al	Si	P	S	Cl	Ar
0.9	1.2												1.5	1.8	2.1	2.5	3.0	--
K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	Se	Br	Kr
0.8	1.0	1.3	1.5	1.6	1.6	1.5	1.8	1.8	1.8	1.9	1.6	1.6	1.8	2.0	2.4	2.8	3.0
Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Ag	Cd	In	Sn	Sb	Te	I	Xe
0.8	1.0	1.2	1.4	1.6	1.8	1.9	2.2	2.2	2.2	1.9	1.7	1.7	1.8	1.9	2.1	2.5	2.6
Cs	Ba	La	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At	Rn
0.7	0.9	1.1	1.3	1.5	1.7	1.9	2.2	2.2	2.2	2.4	1.9	1.8	1.9	1.9	2.0	2.2	--
Fr	Ra	Ac	Rf	Db	Sg	Bh	Hs	Mt	Uun	Uuu	Uub	113	Uuq	115	116	117	118
0.7	0.9	1.1	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--

Ce	Pr	Nd	Pm	Sm	Eu	Gd	Tb	Dy	Ho	Er	Tm	Yb	Lu
1.1	1.1	1.1	1.2	1.2	1.1	1.2	1.2	1.2	1.2	1.2	1.2	1.2	1.3
Th	Pa	U	Np	Pu	Am	Cm	Bk	Cf	Es	Fm	Md	No	Lr
1.3	1.5	1.7	1.3	1.3	1.3	1.3	1.3	1.3	1.3	1.3	1.3	1.5	--

GZ Science Resources

If two identical atoms are covalently bonded together then there will be exactly the same amount of attraction between the shared valence electrons to the nuclei of each atom in the bonded pair. This is because there is no electronegativity difference between atoms, and the valence electrons 'orbit' each atom evenly. This becomes a **non-polar molecule** with **non-polar bonds**.
Example - Iodine molecule I_2

← Path of electrons

Evenly shared
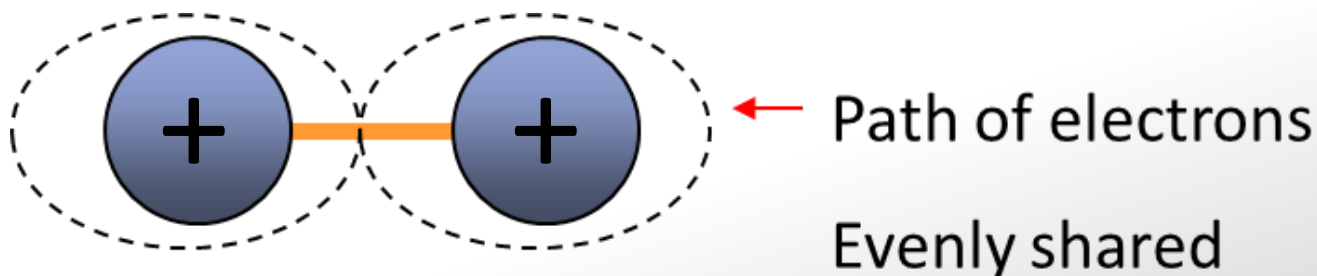
If two different types of atoms are bonded together then they will exert different levels of attraction for the orbiting electrons. That is because they may have different numbers of electron shells and different numbers of protons in their nucleus. This will cause an electronegativity difference. These bonds become **polar bonds,** creating a **dipole**, as there is an electronegativity difference and the valence electrons spend more time around the atom with the higher electronegativity value (that atom becomes slightly negative). The atom that the valence electrons spend less time around becomes slightly positive.

Example – hydrochloric acid HCl

Slightly positive

Electrons orbit less

δ+

δ-

Slightly negative

Electrons orbit more

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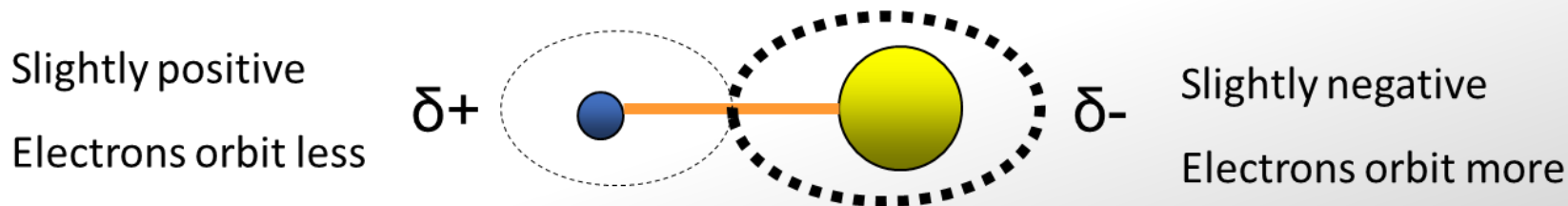
The **overall polarity** of a molecule with polar bonds depends upon whether the molecule is symmetrical or not <u>and</u> if all bonds are the same.

A symmetrical molecule (one where the centres of peripheral atoms coincide) becomes a **non-polar molecule** (<u>only</u> if all bonds are the same) – as the bond dipoles cancel out.

An unsymmetrical molecule (one where the centre of peripheral atoms do not coincide) is a **polar molecule**, – as the bond dipoles **do not** cancel out.

Non-Polar molecule

2.5
3.0

CCl_4
Nonpolar

Polar molecules

3.0

2.5 2.1

CH_3Cl
Polar

3.0

2.1

NH_3
Polar
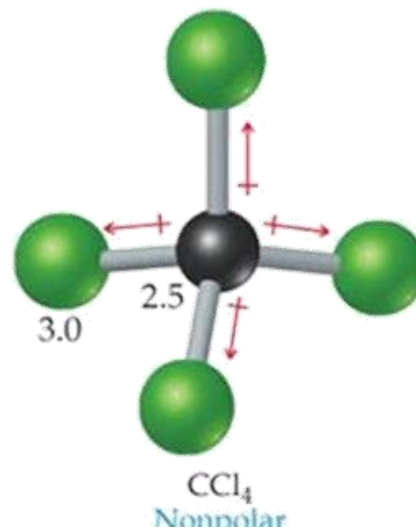
If the bonds (central to peripheral) are different, then even if molecule is symmetrical the dipoles will not cancel out and therefore the molecule is polar overall

extra info
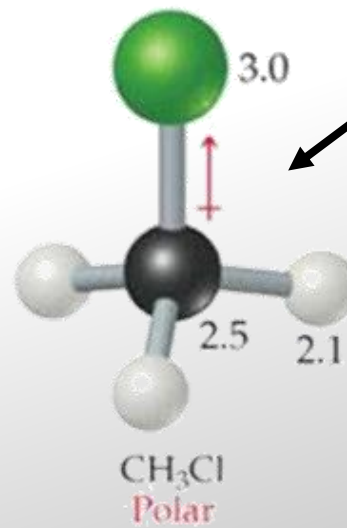
Bond types between atoms can depend on the **electronegativity** of the atoms. Rather than discrete categories, molecules fall along a continuum

Nonpolar covalent bonding
Electrons are shared equally

Cl   Cl

Polar covalent bonding
Electrons are shared unequally

H   Cl

Ionic bonding
Electrons are transferred

Na⁺   Cl⁻

Increasing ionic character

Difference in electronegativity

0.4   1.7
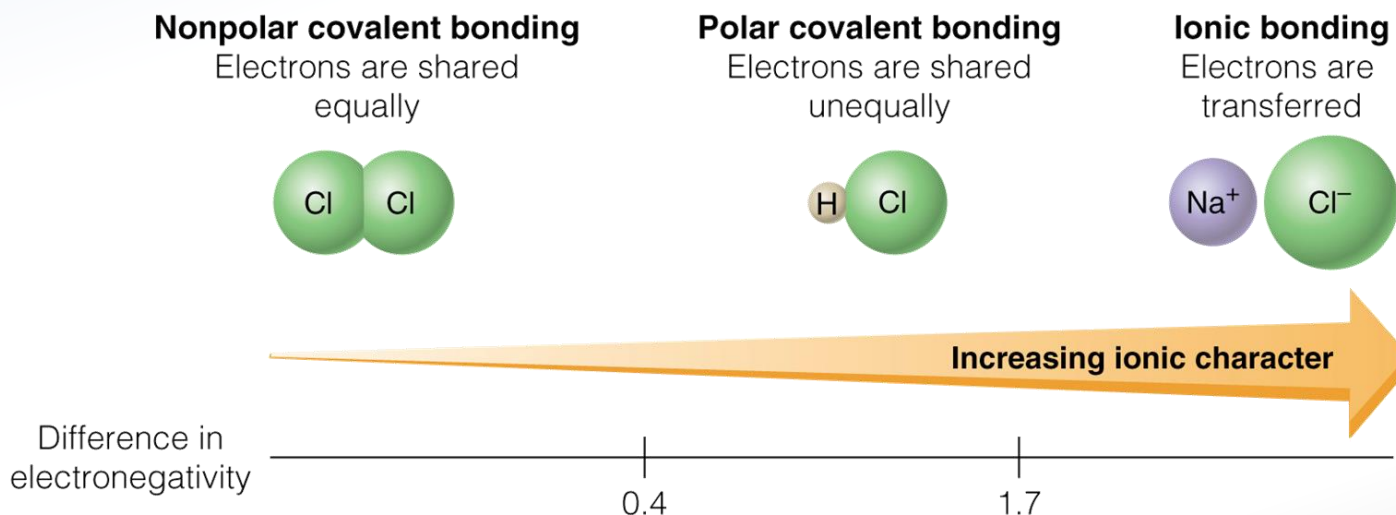
If there is little difference in electronegativity between two atoms then they tend to form a covalent bond with no polarity difference. A greater electronegativity difference creates a polar bond with uneven "sharing" of valance electrons.
If the electronegativity is even greater then there will be a complete transfer of electron from one atom (Metal) to another atom (non-metal) and ions will form that are held together with an ionic bond.

Example: Explain why molecules x (CCl_4) and y (NCl_3) are polar and non-polar?

Polar molecule	Non-polar molecule
1. molecule (NCl_3) is polar (state which one)	1. molecule (CCl_4) is non-polar (state which one)
2. (NCl_3) contains polar bonds and therefore forms dipoles due to electronegativity difference of N and Cl.	2. (CCl_4) contains polar bonds and therefore form dipoles due to electronegativity difference of C and Cl. Cl attracts more electrons than C because it has a bigger atomic number than C but with the same number of shells
3. over the whole molecule the atoms are not distributed symmetrically in 3 dimensions because its shape is (state which one) and has lone pairs of electrons	3. over the whole molecule the atoms are distributed symmetrically in 3 dimensions because its shape is (state which one)
4. Dipoles do not cancel each other out and the whole molecule is polar.	4. Dipoles cancel each other out and the whole molecule is non-polar.

Question 1c (ii): Elements M and X form a compound MX_2. Atoms of element X have a higher electronegativity value than atoms of element M, therefore the M–X bonds are polar. Depending on what elements M and X are, molecules of the compound formed will be **polar** or **non-polar**.
State the most likely shape(s) of the molecule if it is **Polar** and if it is **Non-polar**: Justify your answer and draw diagrams of the possible molecules with dipoles labelled.

If MX_2 is **polar**, this indicates that the polar M–X bonds are not spread symmetrically around the central M atom. There must be either **three or four regions of negative charge** with only two bonded atoms therefore the shape must be bent.

Polar: bent

Three regions of negative charge:

Four regions of negative charge:

Non-polar: linear
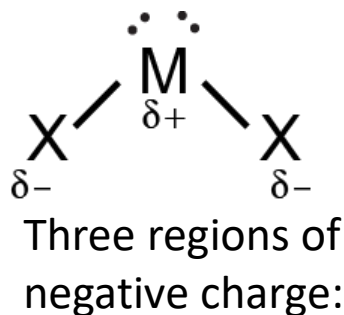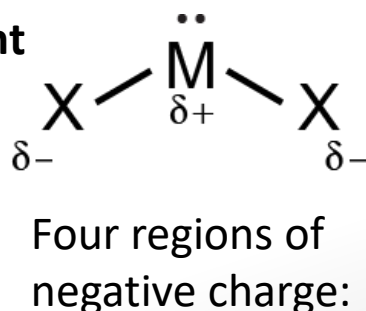
Two regions of negative charge:
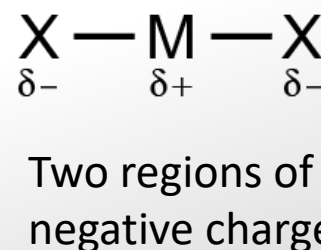
If MX_2 is **non-polar** this means that the polar M-X bonds are spread symmetrically around the central M atom. There must be only **two regions of negative charge** around the M atom, both bonded by X atoms in a linear shape.

Excellence Question

Question 1c: Molecules can be described as being polar or non-polar.
The following diagrams show the Lewis structures for two molecules, SO_2 and CO_2. Identify the polarity
- Justify your choice
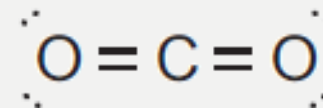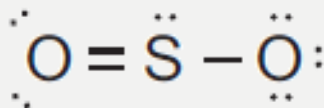
$$O = \ddot{S} - \ddot{O}:$$

$$\ddot{O} = C = \ddot{O}$$

When given a multi- choice question **never** leave it blank

State polarity of molecule first
State polarity of bonds (name atoms)
Link symmetry to dipoles cancelling out (or vice versa)

Answer 1c: SO_2 molecule is polar.
CO_2 molecule is non-polar.
The S–O / S=O bond is polar due to the **difference in electronegativity** between S and O atoms. The bonds are arranged **asymmetrically** in a bent shape around the central S atom; therefore the (bond) **dipoles do not cancel** and the molecule is polar.

The C=O bond is polar due to the **difference in electronegativity** between C and O atoms. The bonds are arranged **symmetrically** in a linear shape around the central C atom; therefore the (bond) **dipoles cancel** and the molecule is non-polar.

Question 1c: $BeCl_2$ and BF_3 are unusual molecules because there are not enough electrons for the central atoms, Be and N, to have a full valance shell. Their Lewis structures are shown below. Both Molecules have the same polarity.
- Identify the polarity
- Justify your choice

$:\ddot{C}l - Be - \ddot{C}l:$ $:\ddot{F} - B - \ddot{F}:$
                                        $:\ddot{F}:$

When given a multi- choice question **never** leave it blank

State polarity of molecule first
State polarity of bonds (name atoms)
Link symmetry to dipoles cancelling out (or vice versa)

Answer 1c: Both molecules are non-polar.
The Be-Cl bond is polar because Cl is more electronegative than Be / the atoms have different electronegativities.
Since both the bonds are the same and arranged symmetrically around the central atom, in a linear arrangement, the bond dipoles cancel out, resulting in a non-polar molecule.
The B-F bond is polar because F is more electronegative than B / the atoms have different electronegativities. Since all three bonds are the same and arranged symmetrically around the central atom, in a trigonal planar arrangement, the bond dipoles cancel out, resulting in another non-polar molecule.

Question 3b: The Lewis structures for two molecules are shown.

Molecule	H–N̈–H | H Ammonia	H–B–H | H Borane
Polarity of molecule	polar	non-polar

Ammonia, NH_3, is polar, and borane, BH_3, is non-polar. Justify this statement.

State polarity of molecule first
State polarity of bonds (name atoms)
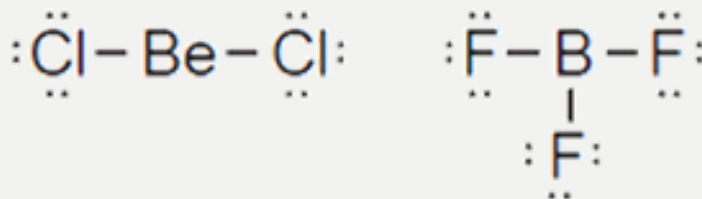Link symmetry to dipoles cancelling out (or vice versa)

Answer 3b: Each N-H bond in NH_3 is polar / forms a dipole because the N and H atoms have different **electronegativities.** The shape of the molecule (due to the presence of one non-bonding electron pair) is trigonal pyramidal which is asymmetrical, so the dipoles / bond polarities do not cancel. The **resulting NH_3 molecule is polar.**
Each B-H bond in BH_3 is polar / forms a dipole because the B and H atoms have different **electronegativities.** The shape of the molecule is trigonal planar which is symmetrical, so the dipoles / bond polarities cancel. The **resulting BH_3 molecule is non-polar.**

Excellence Question

Question 2b: Three-dimensional diagrams for two molecules are shown below.

Molecule		
Name	Dichloromethane	Tetrachloromethane
Polarity of molecule	polar	Non-polar

(i) In the boxes above, identify the polarity of each molecule, by writing either **polar** or **non-polar.**
(ii) Justify your choices.

State polarity of molecule first
State polarity of bonds (name atoms)
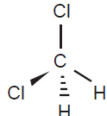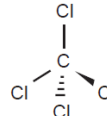Link symmetry to dipoles cancelling out (or vice versa)

In CCl_4, the four C–Cl bonds are polar, i.e. have a dipole, due to the difference in electronegativity between C and Cl. These (equally sized) dipoles are arranged in a symmetric **tetrahedral** shape, resulting in the dipoles / bond polarities cancelling each other out, so CCl_4 is non-polar.
In CH_2Cl_2, there are two types of bond, C–H and C–Cl, each polar with dipoles due to the difference in electronegativity between C and H and C and Cl. These dipoles have different polarities / sizes as H and Cl have different electronegativities. (Despite the symmetric tetrahedral arrangement) the different (sized) dipoles / bond polarities do not cancel each other out, so CH_2Cl_2 is polar.

Question 2c. The Lewis structures for two molecules are shown below.
Hydrogen cyanide, HCN, is polar, and carbon dioxide, CO_2, is nonpolar.
Both molecules are linear.
Explain why the polarities of the molecules are different, even though their shapes are the same.

Molecule	H−C≡N	O=C=O
Polarity of molecule	Polar	Nonpolar

180°

In HCN, the **two bonds are polar** due the **difference in electronegativity** between H and C, and C and N. The resulting bond dipoles are differing in size as H and N have different electronegativities, so despite the symmetric linear arrangement **the bond dipoles do not cancel** and HCN is overall polar.

The **C=O bond is also polar** due to O being more electronegative than C giving these bonds dipoles. But because both **bonds are identical and are arranged symmetrically** in a linear shape, the **bond dipoles cancel** and the molecule is non-polar overall.

Question 2b. The following table shows the Lewis structures (electron dot diagrams) for the molecules, $CHCl_3$ and NH_3.

Molecule	$CHCl_3$	NH_3
Lewis structure	H | :Cl−C−Cl: | :Cl:	H−N̈−H | H
Polarity	Polar	Polar

(i) In the boxes beside, identify the polarity of each molecule by writing either **polar** or **non-polar**.
(ii) Justify your choices.

In $CHCl_3$, there are two types of bond, C−H and C−Cl, each polar, due to the difference in electronegativity between C and H and C and Cl atoms. These dipoles have different polarities / sizes as H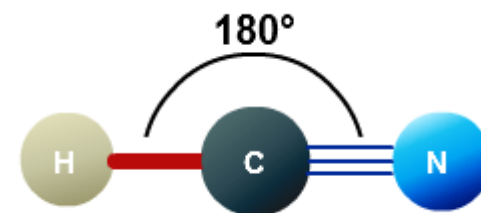 and Cl have different electronegativities. (Despite the tetrahedral arrangement appearing symmetrical) the different (sized) bond dipoles do not cancel each other out, so $CHCl_3$ is polar.
In NH_3, the three N−H bonds are polar, i.e. have a dipole, due to the difference in electronegativity between N and H atoms. These (equally sized) dipoles are arranged in a non-symmetrical trigonal pyramidal shape, resulting in the bond dipoles not cancelling each other out, so NH_3 is polar.

GZ Science Resources

When a solid mixes into a liquid and can longer be seen it has **dissolved**. The liquid is called the **solvent** and it pulls apart the bonds between the solid particles, called the **solute**, and they **diffuse**. A **solution** is then created when the solvent particles (often water) are mixed up with the broken apart solute particles.
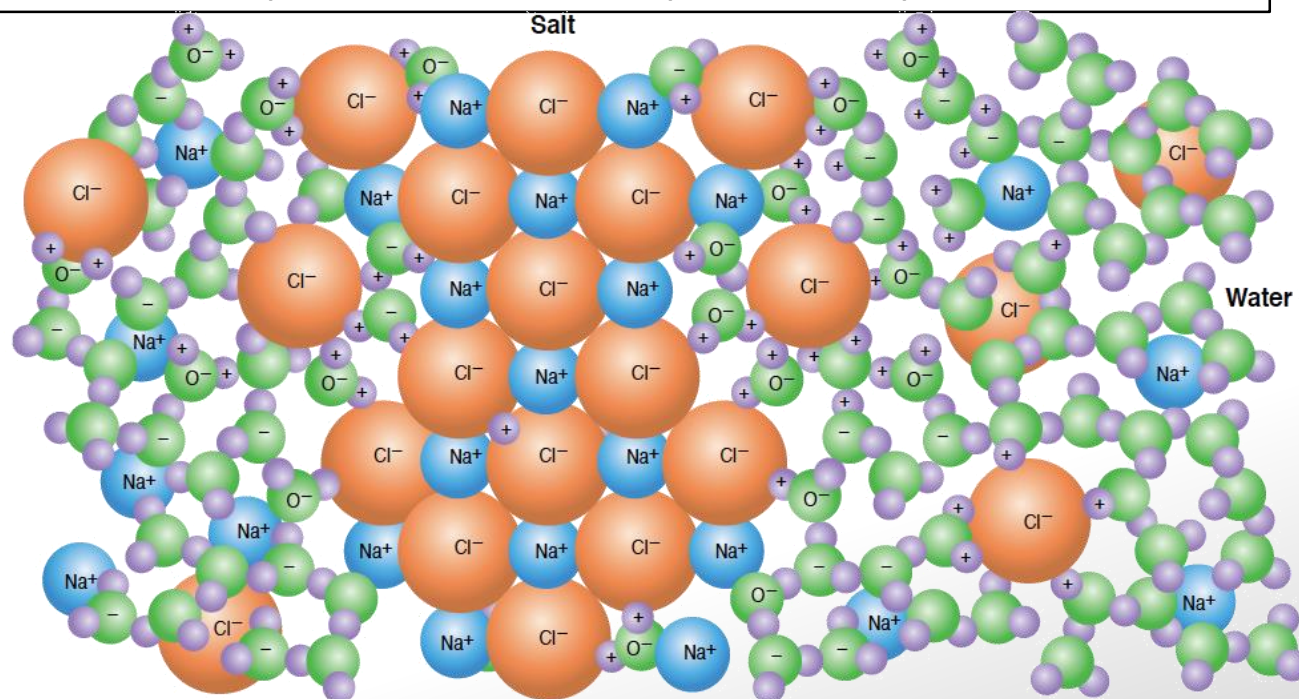
For a solute to dissolve, the solvent particles must form bonds with the solute particles. Water, being polar attracts ions because of their dipoles, and so dissolve many ionic substances.

The solubility of a substance is the **amount** of that substance that will dissolve in a given amount of solvent. Solubility is a quantitative term. Solubility will vary depending on the solvent and the solute. The terms soluble and insoluble are relative. Some substances can be sparingly soluble where only the minutest percentage dissolves. For a solute to dissolve the attraction to the solvent molecules must be stronger than the bonds holding the atoms/molecules of the solute together.
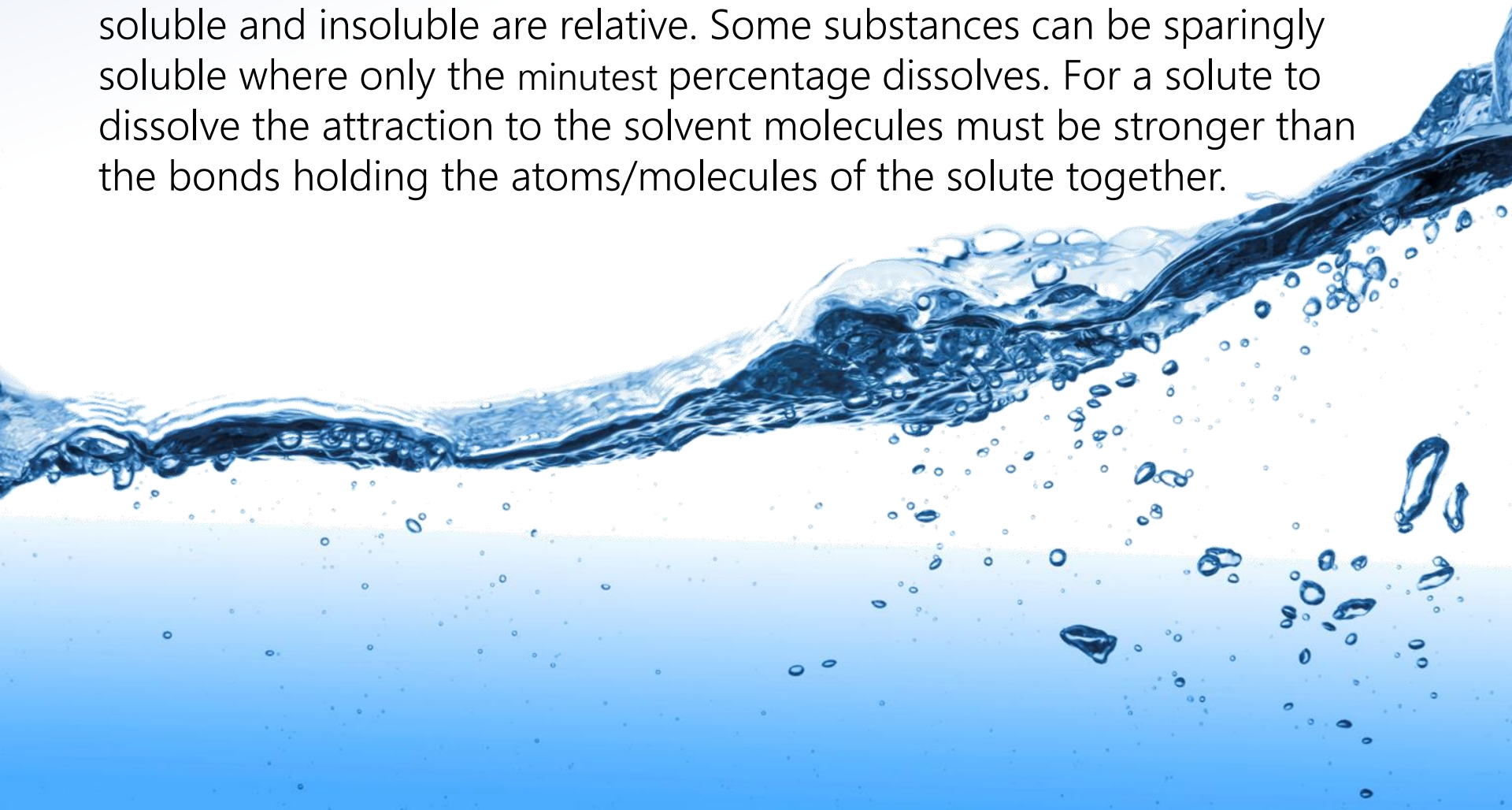
Polar substances dissolve polar substances.
e.g. Water, being polar attracts the molecules of other polar substances (e.g. HCl) and will dissolve them.

Polar substances will not dissolve non-polar substances.
e.g. Water, (polar) has a stronger attraction to itself than to non-polar molecules (e.g. cyclohexane) and will not dissolve them.

Non-polar substances dissolve non-polar substances.
e.g. Non-polar solvents (like cyclohexane) attract non-polar solutes (like napthalene) by the same weak molecular forces they attract themselves by and so will dissolve non-polar solutes.
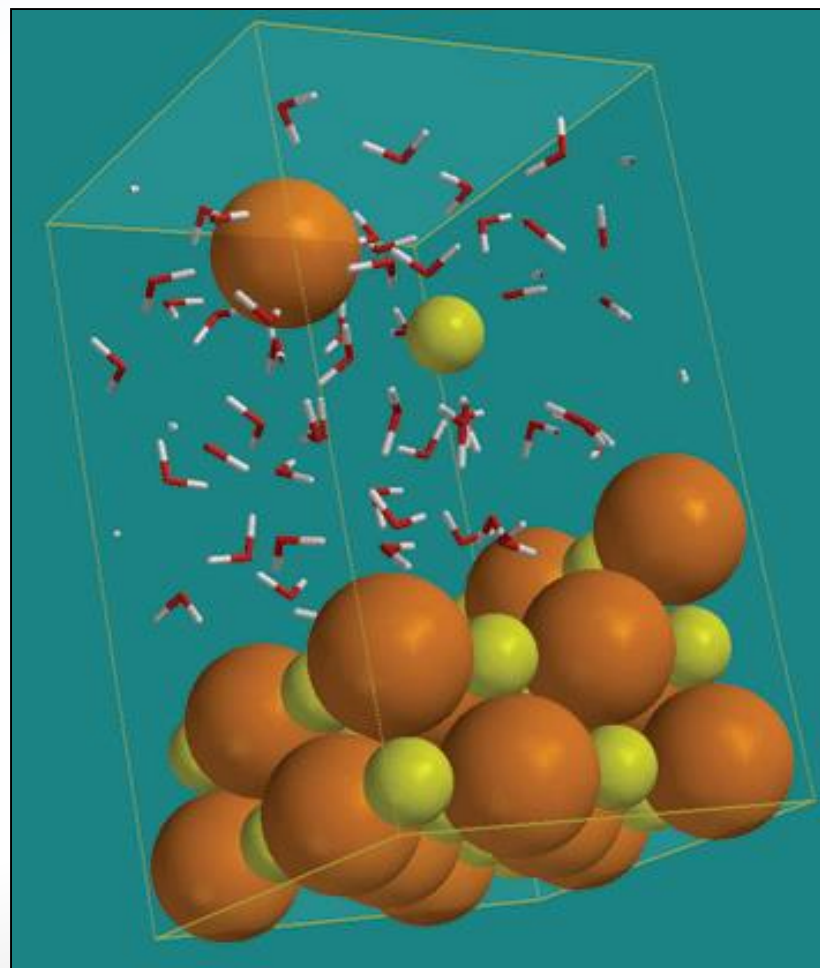
Ionic solid dissolving in water

The water molecule has dipoles, caused by the separation of charge ($\delta+$ - $\delta-$). Due to the **asymmetry** of the molecule, their dipoles reinforce making the oxygen side of the molecule partially negative ($\delta-$) and the hydrogen side partially positive ($\delta+$). Such molecules are called '**polar**'. Polarity causes a stream of water molecules to attract to a charged plastic pen.

Polarity of substances	
Polar	Non-Polar
water	cyclohexane
methanol	benzene
ethanol	hydrocarbons (e.g.
acetic acid	petrol)
hydrogen	iodine
chloride	
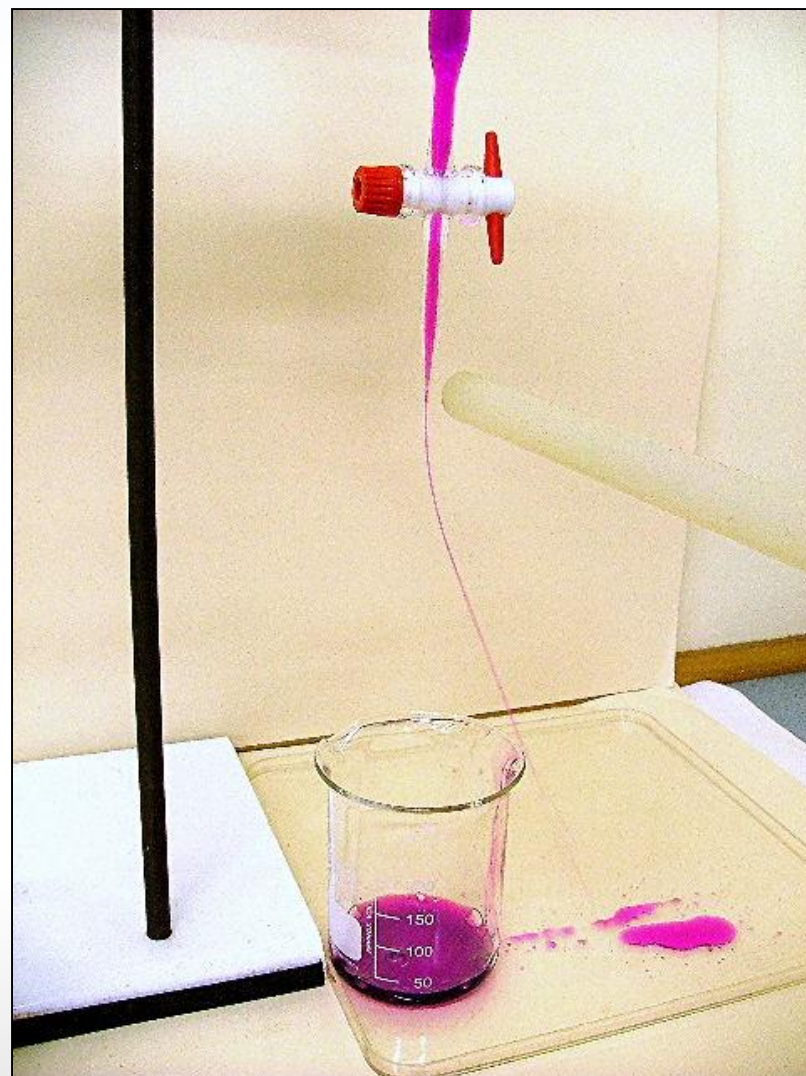
Question 3b: Use your knowledge of structure and bonding to explain the dissolving process of sodium chloride in water.
Support your answer with an annotated (labelled) diagram.

Answer 3b : Solubility

When sodium chloride is dissolved in water the **attractions** between the polar water molecules and between the ions in the salt are replaced by attractions between the water molecules and the ions. The negative charges on the oxygen ends of the water molecules are attracted to the positive Na$^+$ ions, and the positive hydrogen ends of the water molecules are attracted to the negative Cl$^-$ ions.
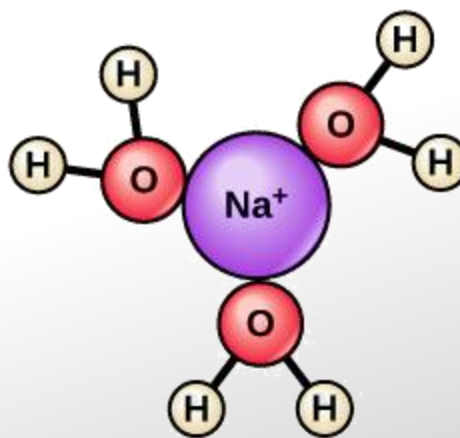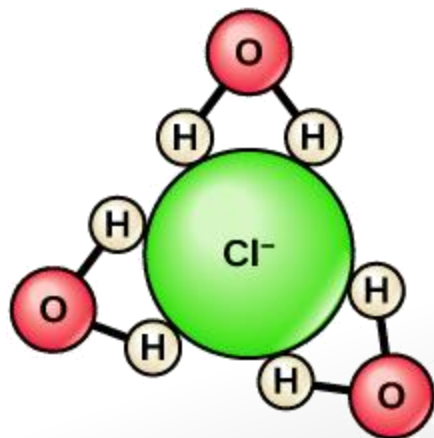
Only a few ions of each are needed in the diagram. Don't forget to label

Question 1b: (iii) Sodium chloride, NaCl, is another compound that is excreted from the body in sweat.
Use your knowledge of structure and bonding to explain the dissolving process of sodium chloride, NaCl, in water.
Support your answer with a labelled diagram.

Only a few ions of each are needed in the diagram. Don't forget to label

Answer 3b : Solubility

Sodium chloride is an ionic substance made up of Na$^+$ and Cl$^-$ ions arranged in a (3D) lattice and held together by ionic bonds. The δ- O of polar water molecules are attracted to the positive Na$^+$, while water's δ+ H is attracted to the negative Cl$^-$, this attraction is sufficiently strong to overcome the attractions between the ions in the salt / crystal /lattice (and between the water molecules in the solvent), dissolving the NaCl.

Sodium (Na)

Chlorine (Cl)

NaCl crystal structure

NaCl in water
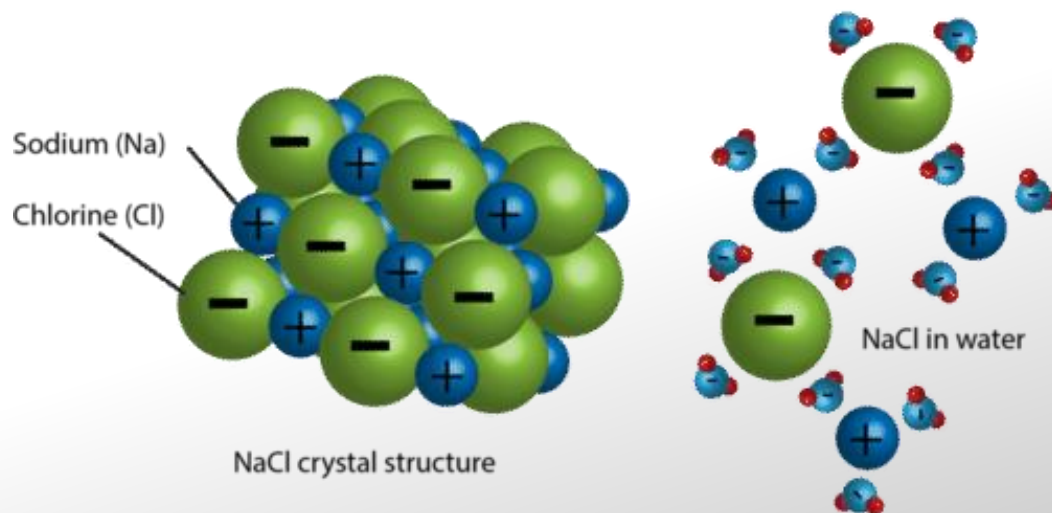
Question 3d (i) . Use an annotated diagram to show how solid **A (ionic solid)** is able to dissolve in water.
Show the solid before dissolving, and the dissolving process of the solid.
(ii) Explain the attractions that allow solid **A** to be soluble in water.
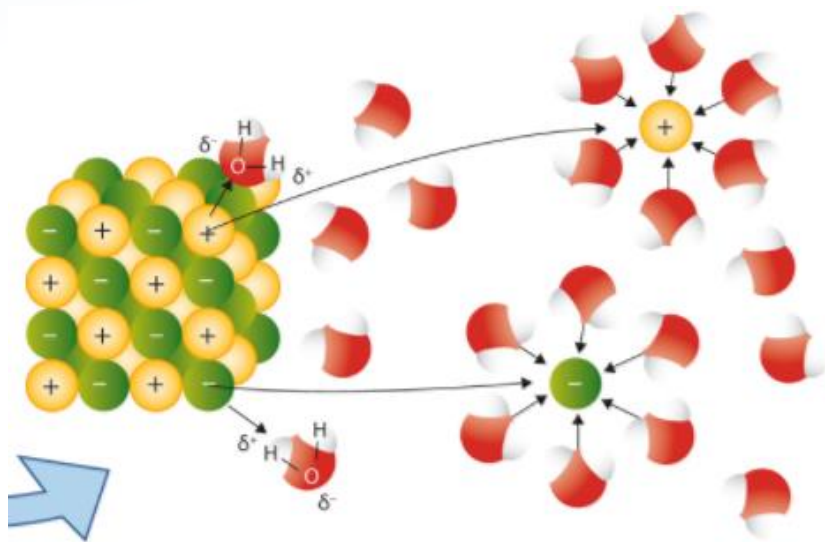
Solid A is ionic and when it dissolves in water, it **separates into its ions**. The ions are **charged** and are attracted to the charged ends of the polar water molecule. The slightly negative charges on the oxygen ends of the water molecules are attracted to the positive ion, and the slightly positive hydrogen ends of the water molecules are attracted to the negative ions. **This causes the ions to be surrounded by water molecules and it dissolves.**

This solid is soluble because the **force of attraction between the ions and water is strong enough to overcome the forces holding the ions together** along with the forces holding the water molecules together (in the solvent).

https://www.learnit.ru/476728/

Question 1c Compare the solubilities of iodine, $I_{2(s)}$, in water, $H_2O_{(l)}$ – a polar solvent, and in cyclohexane, $C_6H_{12(l)}$ – a non-polar solvent.
Use your knowledge of structure and bonding to explain the solubility of iodine in these two solvents.

Iodine is a non-polar (covalent) molecular substance made up of I_2 molecules held together by weak intermolecular forces.
Iodine is soluble in cyclohexane, but does not easily dissolve in water.
For iodine in water, the iodine-water attractions are not strong enough to overcome both the iodine-iodine / solute-solute and the strong water-water / solvent-solvent attractions.

For iodine in cyclohexane, the iodine-cyclohexane attractions are strong enough to overcome iodine-iodine / solute-solute and cyclohexane-cyclohexane / solvent-solvent attractions because all attractive forces are similar (nonpolar).

https://www.learnit.ru/476728/

Substances are grouped together according to the type of bonds they have between particles, and consequently the structure they form.

L2 will cover four groups of substances; Molecular, metallic, ionic and covalent network. The physical properties of these groups will be linked to their structure.

Molecular solids
Non-metals forming molecules

S_2
sulfur

HCl
Hydrogen chloride
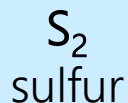
I_2
iodine

Ionic solids
Non-metals and metals together forming a ionic compound

KI
Potassium iodide

$CuSO_4$
Copper sulfate

NaCl
Sodium chloride

Metallic solids
Elements that are metals

Fe
iron

Al
aluminium

Cu
copper

Covalent network solids
Carbon and silicon dioxide

SiO_2
Silicon dioxide

C
diamond

C
graphite

non-metal + non-metal
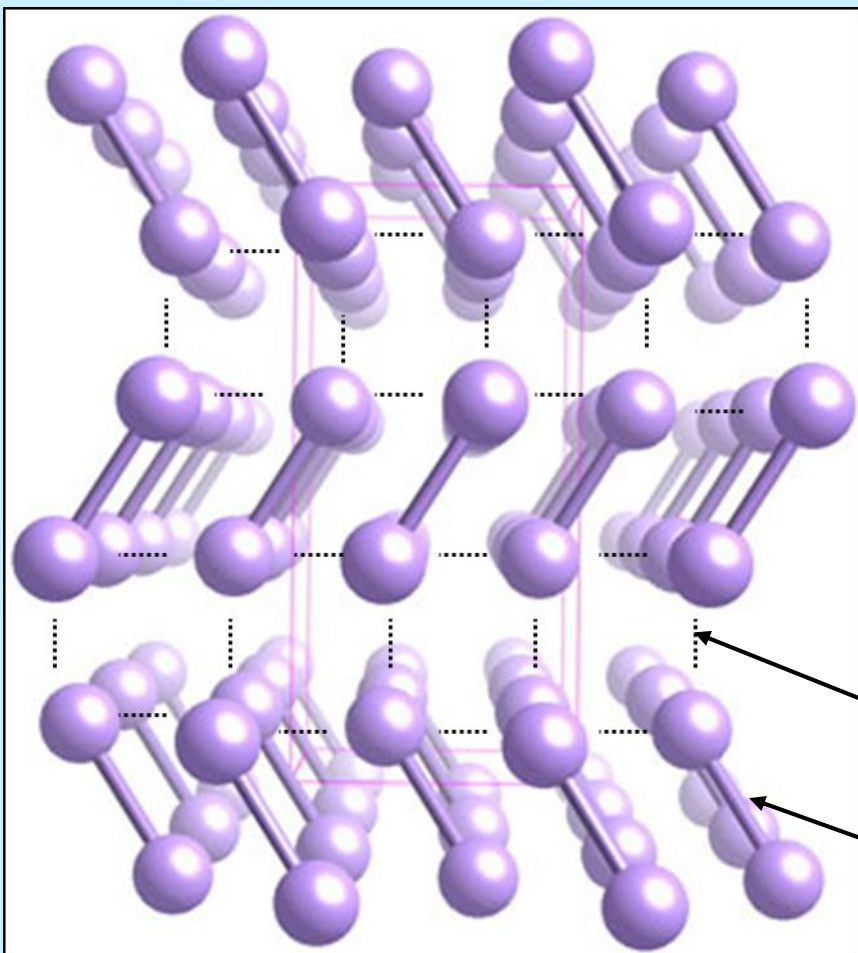
Non-polar Molecules are held together by **weak intermolecular bonding** caused by temporary dipoles only. These are induced (created) by electrons randomly spending more time around one nucleus than the other.

Within the Molecules, the atoms are held together by **strong covalent bonds**.

Weak intermolecular bond

Strong covalent bond

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Polar molecules are held together by **weak intermolecular** forces, the δ –ve end of one molecule is attracted to the δ +ve end of another.

This is due to both temporary <u>and</u> permanent dipoles (which tend to be stronger). Permanent dipoles are induced by electrons spending more time around one nucleus in the molecule that has greater electronegativity than the other.

extra info

Intermolecular Force between diploles
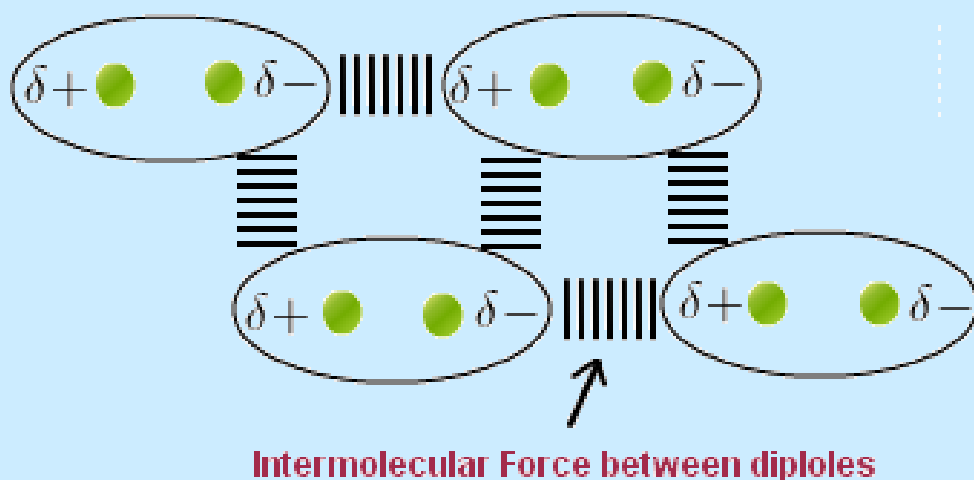
δ means slightly

Note the distinction:
Intra-molecular Forces: the strong bonding forces within a molecule. i.e. the <u>covalent bonds</u> holding the molecule together.
Inter-molecular Forces: the weak bonding forces between molecules due to the attractions between partial charges. i.e. permanent dipole

Polar molecular solids are soluble in water

1. Hydrogen chloride (HCl) is a molecular solid

2. Hydrogen chloride is composed of covalently bonded atoms to form molecules

3. These molecules are held together by weak intermolecular forces

4. These molecules are polar therefore, the electrostatic attractions of water molecules (which is stronger than the weak intermolecular forces) have sufficient strength to pull the molecules apart hence, hydrogen is soluble
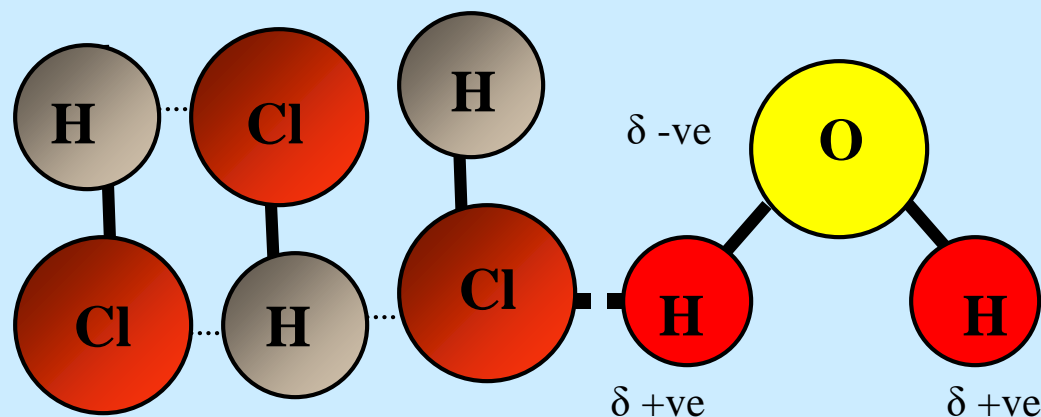
Non-polar molecular solids are insoluble in water

1. Iodine is a molecular solid

2. Iodine is composed of covalently bonded atoms to form molecules

3. These molecules are held together by weak intermolecular forces

4. Iodine is <u>non-polar</u> therefore, the electrostatic charges of the water do not have sufficient strength to overcome the weak intermolecular forces holding the molecules together hence, iodine is insoluble

δ -ve

O

H **H**

δ +ve δ +ve

δ -ve

O

H **H**

δ +ve δ +ve
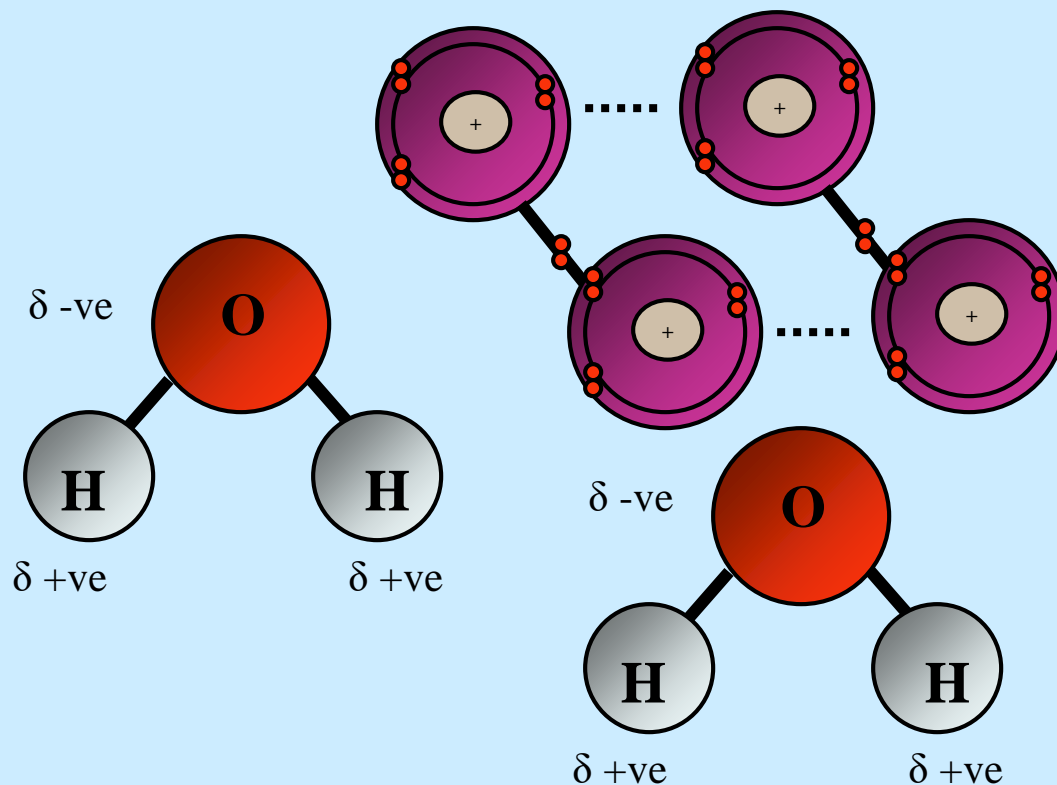
molecular solids have a low melting point

For example:

1. Carbon dioxide is a molecular solid (at low temperatures below -56∘C)

2. Carbon dioxide is composed of covalently bonded atoms to form molecules

3. These molecules are held together by weak intermolecular forces

4. These forces require small amounts of energy to break apart the solid (but not the individual molecules which are held together by strong covalent bonds) therefore, carbon dioxide has low melting point

Many molecular solids are only solid at temperatures well below 0∘C and at room temperature they are gases

Heat

For example:

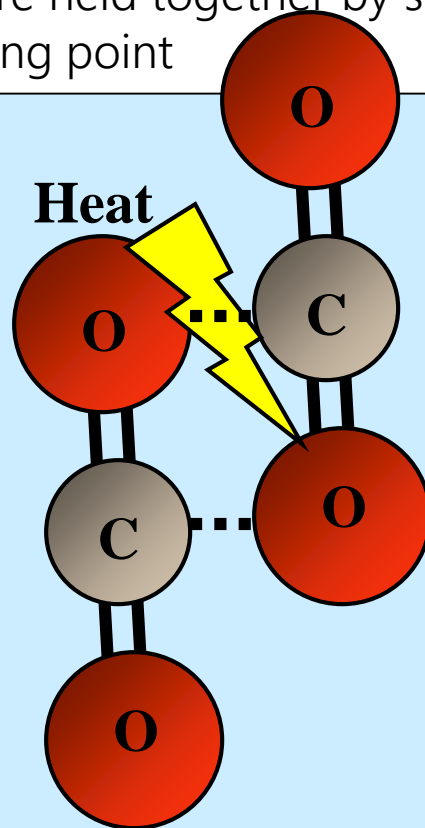1. Sulfur is a molecular solid

2. Sulfur is composed of covalently bonded atoms to form molecules

3. These molecules are held together by weak inter molecular forces

4. These forces require small amounts of energy to break apart the solid (but not the individual molecules which are held together by strong covalent bonds) therefore, sulfur is easily broken up

Force applied

Weak intermolecular force
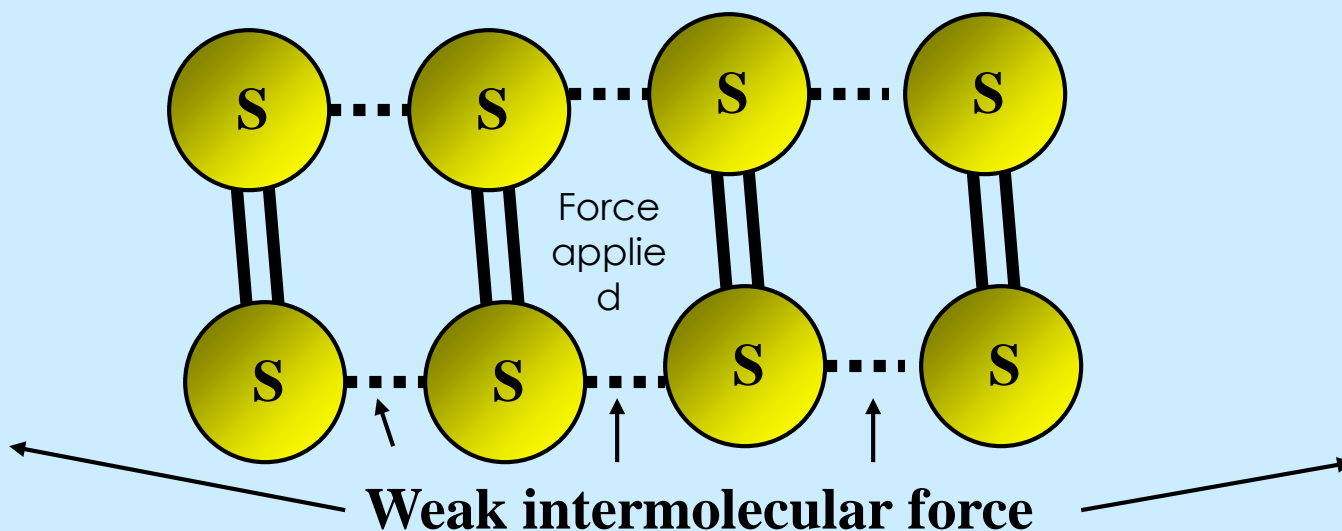
molecular solids do not conduct electricity

In order for a substance to be electrically conductive there must be free moving (mobile) charged particles

For example:
1. Iodine is a molecular solid
2. Iodine is composed of covalently bonded atoms to form molecules; weak intermolecular forces hold these molecules together
3. There are no free moving charges therefore, iodine cannot conduct electricity

nucleus

Valence electrons

Covalently shared electrons

Weak intermolecular bonding
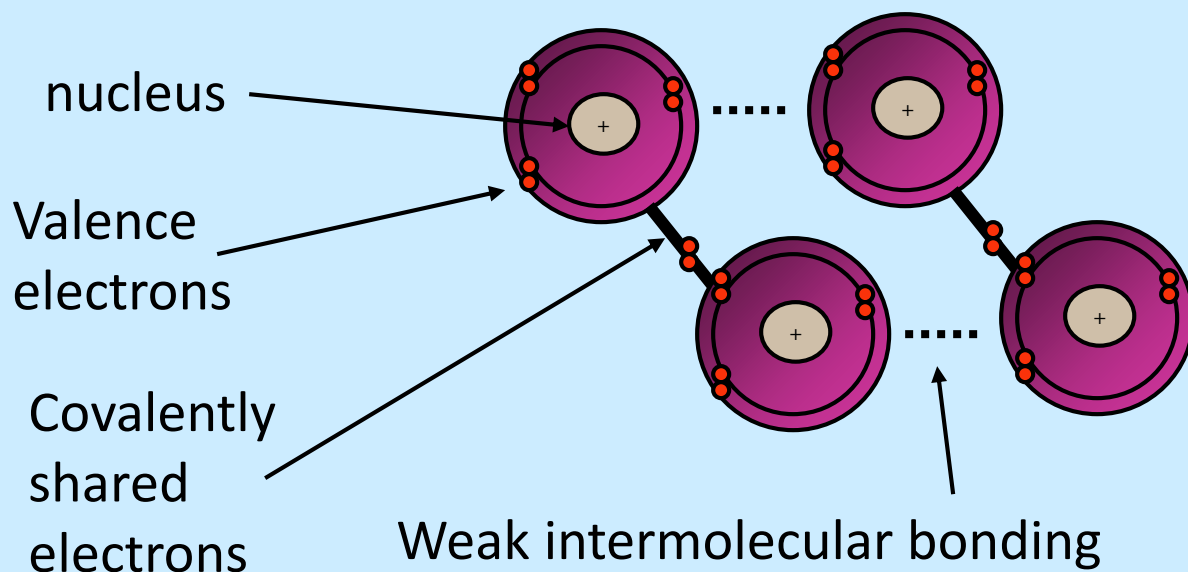
Fully occupied valence electrons remain in 'fixed orbit' around nucleus and are not available to carry charge. The molecule is neutral

Metals atoms are arranged as positive ions held in place in ordered layers by strong attractive **non-directional** bonding, forming a lattice - this gives metals strength.

fixed cation

sea of electrons
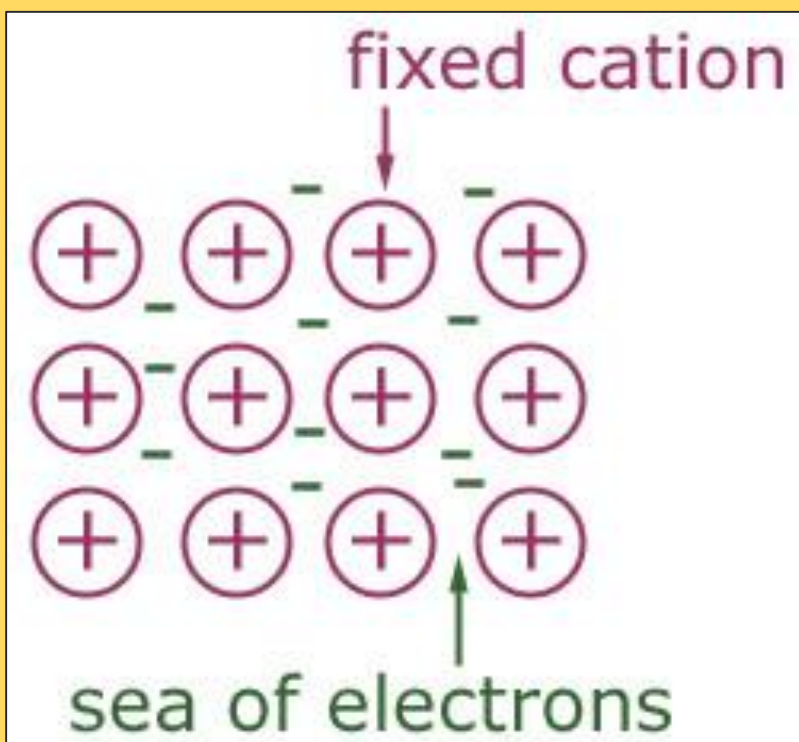
Metal atoms are held together in a 3–D lattice by metallic bonding in which valence electrons are attracted to the nuclei of neighbouring atoms. The attraction of the metal atoms for the valence electrons is not in any particular direction; therefore metal atoms can move past one another without disrupting the metallic bonding, therefore metal is ductile and malleable.

The atoms are packed tightly together - this makes metals dense

For example:

1. Copper is a metallic solid

2. Copper is arranged as positive ions held in place in ordered layers by strong attractive non-directional forces, in a sea of de-localised electrons

3. Electrons are free moving hence can carry a charge

4. Therefore, copper **can conduct electricity**

Free moving charged particles are required to carry a charge and for a substance to be electrically conductive

Electrons from the outer shells of the metal atoms move freely throughout the lattice. - this makes metals excellent conductors of heat and electricity
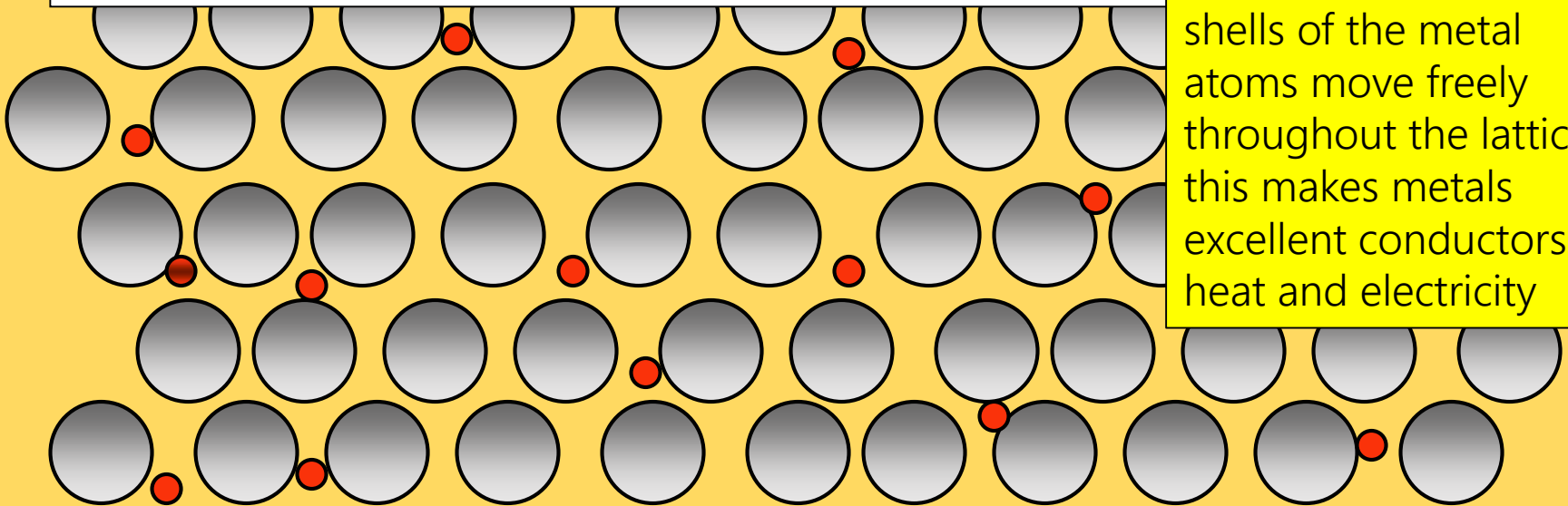
For example:

1. Lead is a metallic solid

2. Lead is arranged as positive ions held in place in ordered layers by strong attractive non-directional forces, in a sea of de-localised electrons

3. These forces require a large amount of energy to break therefore the electrostatic attractions of water molecules do not have sufficient strength to pull the atoms apart

4. Therefore, **lead is insoluble**

In order for substance to dissolve in water (a polar liquid) the attraction between the particles in a substance must be less than the attraction towards water molecules
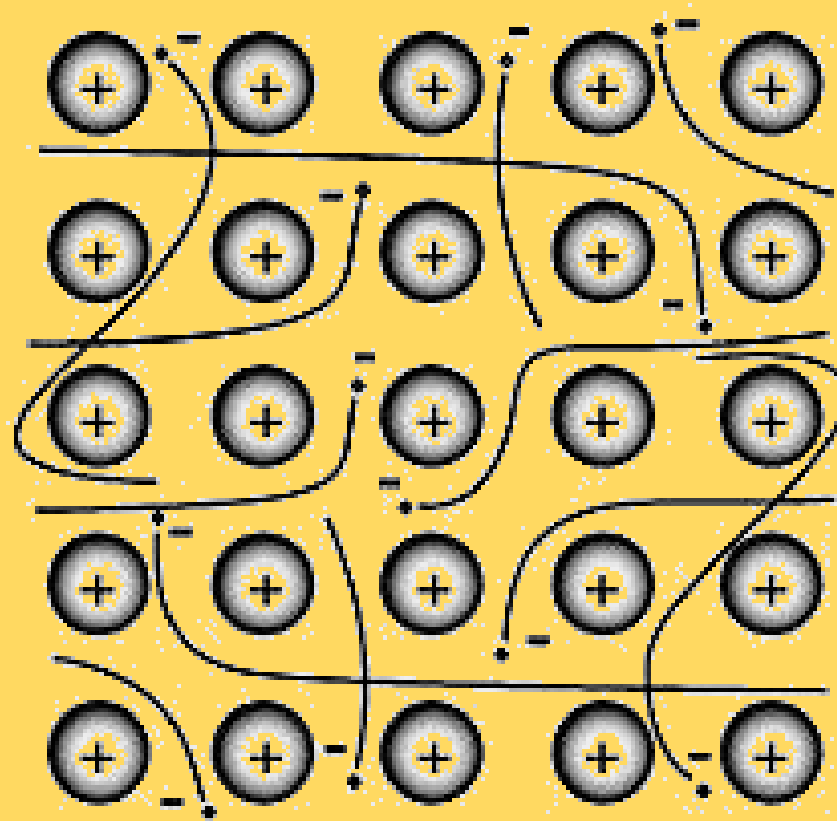
Layers of ions can slide over each other without breaking- this makes metals hard and also malleable and ductile

For example:

1. Iron is a metallic solid

2. Iron is arranged as positive ions held in place in ordered layers - **a lattice**, by strong attractive non-directional forces, in a sea of de-localised electrons

3. These forces require large amounts of energy to break apart the solid therefore aluminium is not easily broken up

4. However, when pressure is applied layers can slide over each other, and as the attractive forces are non-directional the metallic particles remain strongly bonded. – This gives the metallic solids the properties of being malleable (moulded into flat sheets) and ductile (drawn out to thin wires)

Metallic solids have a high melting point

For example:

1. Aluminium is a metallic solid

2. Aluminium is arranged as positive ions held in place in ordered layers by strong attractive non-directional forces, in a sea of de-localised electrons

3. These forces require a large amount of energy (high temperature) to break apart the metallic solid therefore the melting point is very high.

The strength of the bonds between particles determines the energy required to break them, and therefore the amount of energy to change a solid into a liquid (the melting point) where the bonds are somewhat broken. Metals in general, have very strong bonds which makes them solid at room temperature (Mercury is the exception)

Three steps to answering structure and physical properties questions.
❑ The first is state the name of the solid.
❑ The second is describe the structure of the solid.
❑ The third is link the structure of the solid to the physical property discussed.

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Metal + Non-Metal

CHLORIDE ION

SODIUM ION

Cl^-

Na^+

SODIUM CHLORIDE LATTICE STRUCTURE
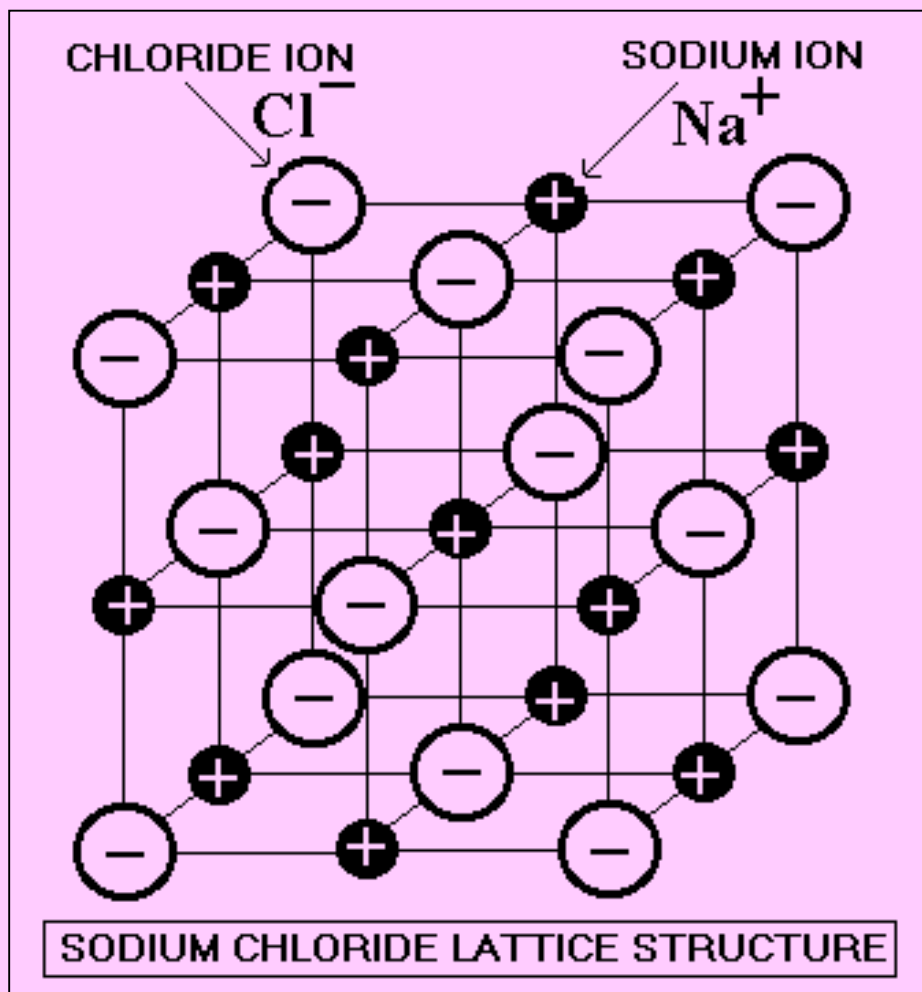
An ionic solid is composed of ions held together by strong directional electrostatic forces (ionic bonding) between +ve (cations) and –ve (anions) ions in a 3-dimensional lattice.

Ionic solids are soluble in solution

In order for substance to dissolve in water (a polar liquid), the attraction between the particles in a substance must be less than the attraction towards water molecules

For example:

1. Sodium chloride (NaCl) is an ionic solid

2. Sodium chloride is composed of ions held together by strong directional electrostatic attractions between +ve and –ve ions in a lattice

3. The electrostatic attractions of water molecules have sufficient strength to pull the ions apart

4. Therefore, the solid will dissolve and is soluble

NaCl first placed in water
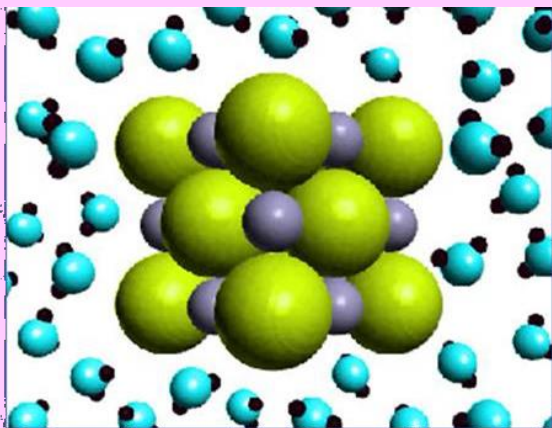
Na$^+$ and Cl$^-$ ions breaking apart

The positive hydrogen end of water is attracted to the anions and the negative oxygen end of water is attracted to the cations

Ionic solids are conductive when in solution or molten only

For example:

1. Sodium chloride is an ionic solid

2. Sodium chloride is composed of ions held together by strong directional electrostatic forces between +ve and –ve ions in a 3-d lattice

3. When solid the ions are not free to move therefore it doesn't conduct electricity

4. However, when molten, or dissolved in solution, the bonds are broken and the ions are free to move therefore sodium chloride can conduct electricity

Free moving charged particles are required to carry a charge and for a substance to be electrically conductive

Na+ Cl- Na+ Cl-

Cl- Na+ Cl- Na+

For Example:

1. Sodium chloride is an ionic solid

2. Sodium chloride is composed of ions held together by strong directional electrostatic attractions between +ve and –ve ions in a 3-d lattice so requires a lot of energy to break the bonds

3. However, because the ionic bonding is directional, if sideways force is applied and a sheet of the lattice slides then ions of the same charge may come in close contact with each other and repel hence the ionic solid is brittle (and can break into pieces)

force

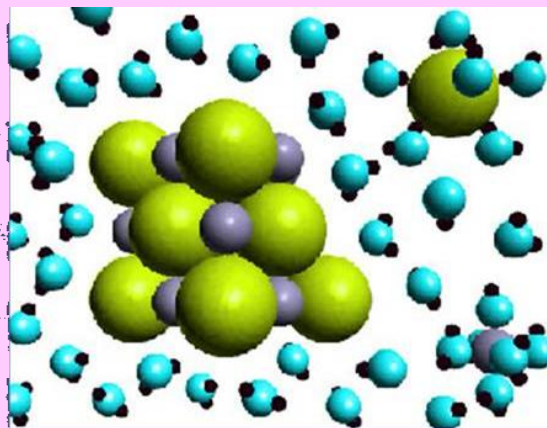GZ Science Resources

Ionic solids have a high melting point

For example:

1. sodium chloride is an ionic solid

2. Sodium chloride is composed of ions held together by strong directional electrostatic attractions between +ve and –ve ions in a 3-d lattice

3. Because these strong bonds require a large amount of energy to break the ionic solids have a high melting point.

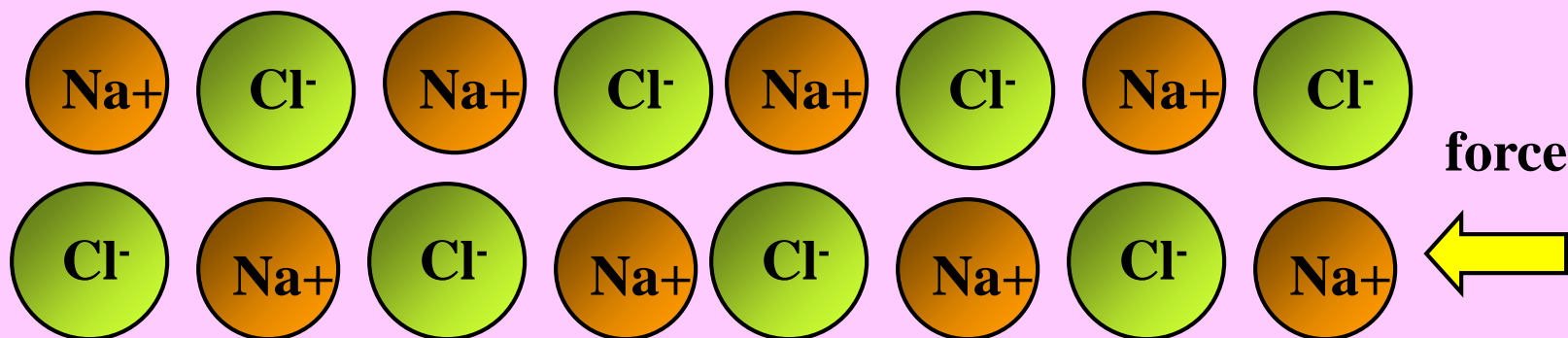Na+ Cl⁻ Na+ Cl⁻

Cl⁻ Na+ Cl⁻ Na+

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Resources
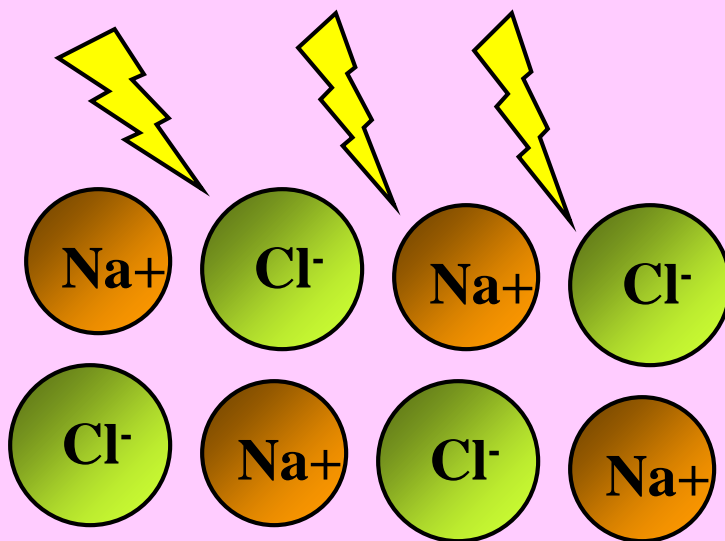
diamond

graphite

Silicon dioxide

All atoms are held together by strong covalent bonds

Diamond is a 3-dimensional covalent network structure where atoms are held together by strong covalent bonds in all planes

Graphite is a covalent network structure that is in 2 dimensional sheets (graphite). Between the layers are free moving electrons from the valance electrons of the carbon atoms.

Silicon dioxide (SiO_2) is a 3-dimensional covalent network structure

For example:

1. Diamond is a 3-dimensional covalent network structure (diamond)

2. All atoms are held together by strong covalent bonds

3. There is no free moving charged particles

4. Therefore, diamond **cannot conduct electricity**
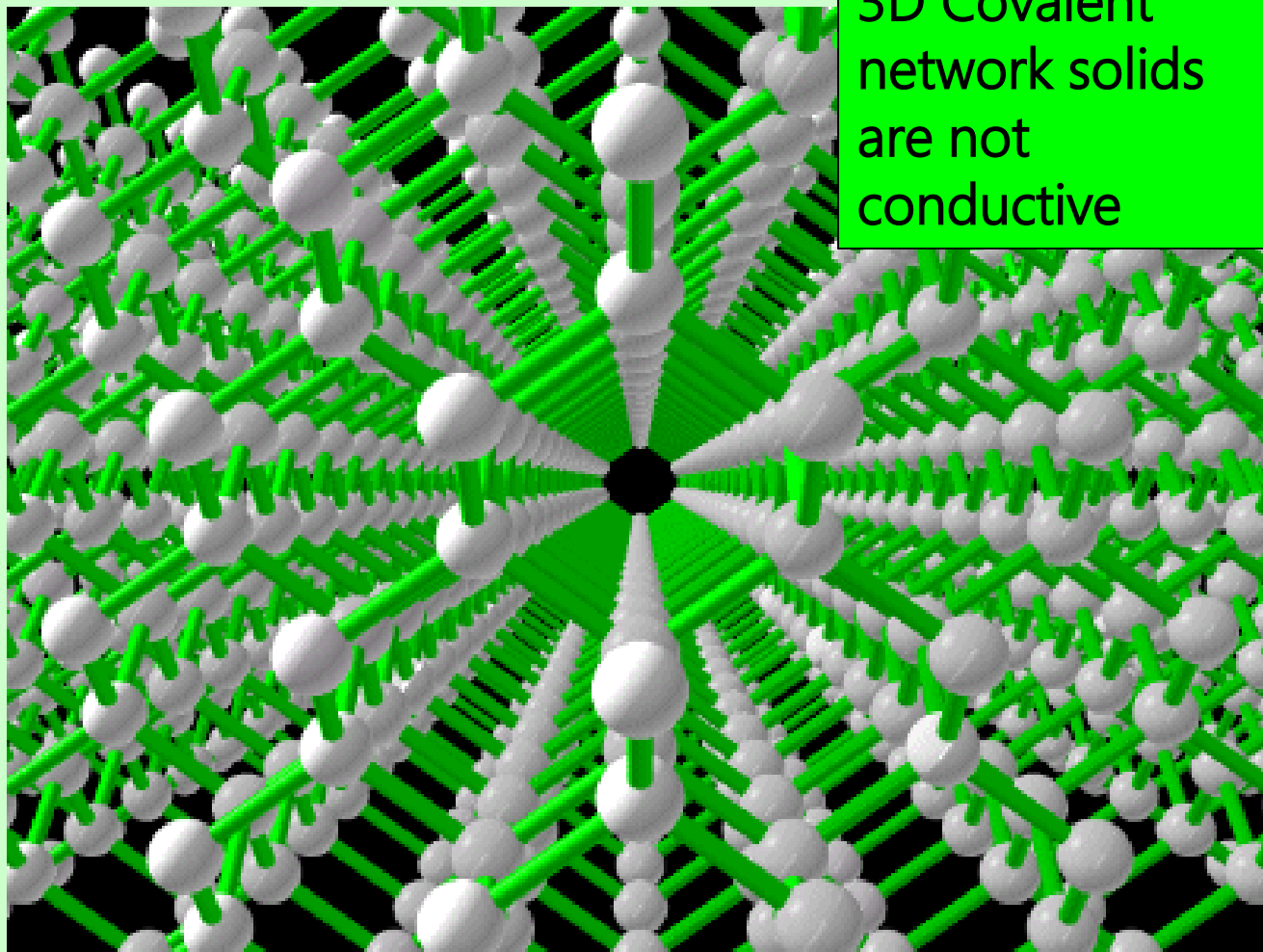
3D Covalent network solids are not conductive

2D Covalent network solids are conductive

For example:

1. Graphite is a covalent network that is in 2-dimensional sheets

2. Between the layers are free moving electrons from the valance electrons of the carbon atoms.

3. The free moving electrons can carry a current

4. Therefore, graphite **can conduct electricity**
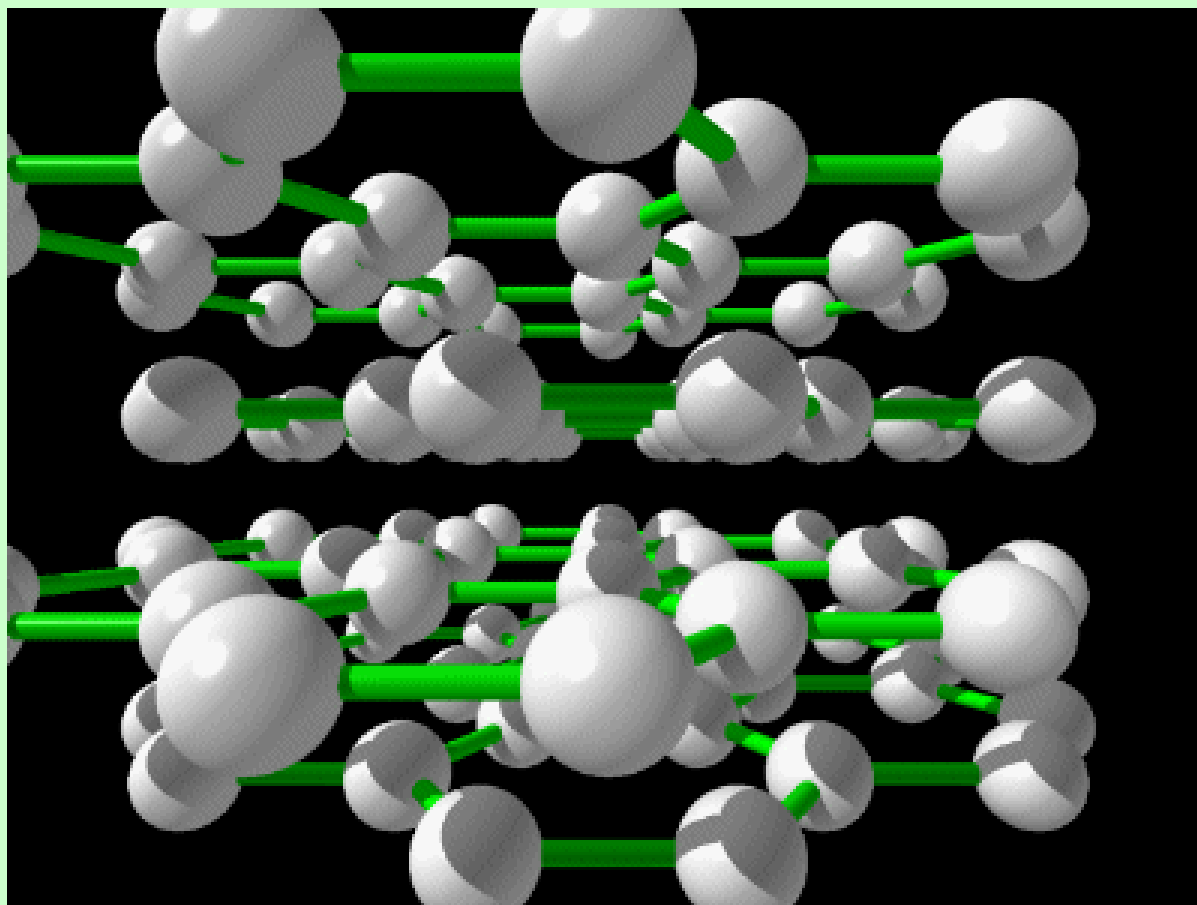
Covalent network solids are not soluble

For example:

1. Silicon Dioxide is a 3-dimensional (or 2-dimensional) covalent network structure

2. All atoms are held together by strong covalent bonds

3. These forces require a large amount of energy to break therefore, the electrostatic attractions of water molecules do not have sufficient strength to pull the ions apart

4. Hence, silicon dioxide will not dissolve in water and is insoluble
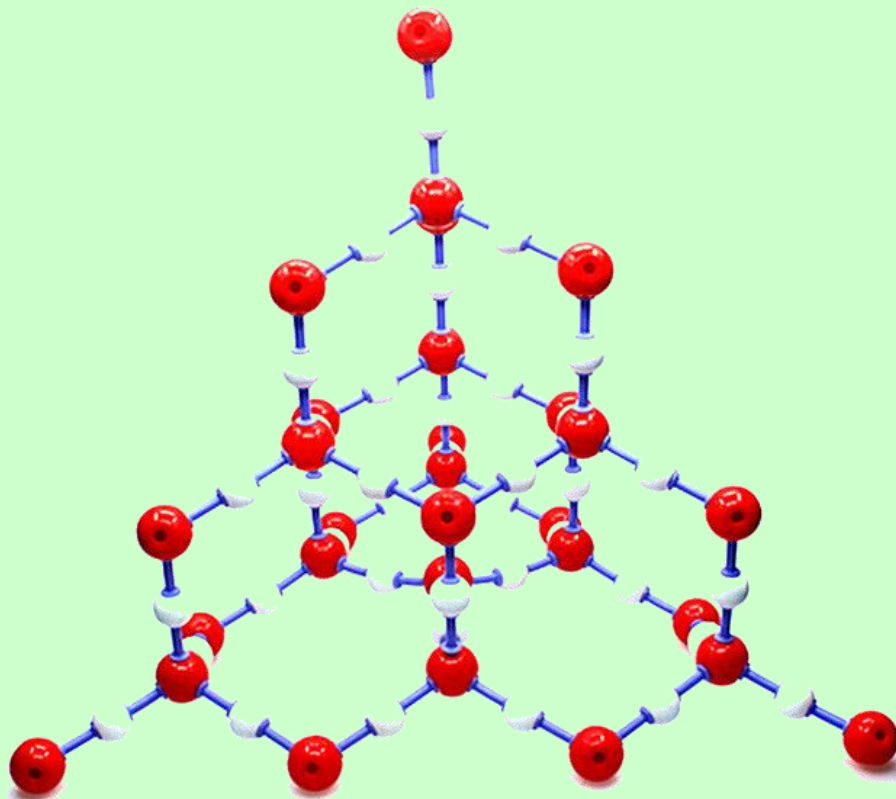
Covalent network solids have a high melting point

For example:

1 Diamond is a 3-dimensional (or 2-dimensional) covalent network structure

2. All atoms are held together by strong covalent bonds

3. These forces require a large amount of energy to break

4. Therefore, diamond has a very high melting point.
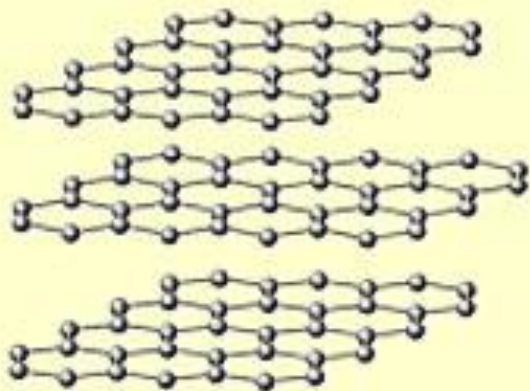
graphite   diamond   silica

For example:

1. Diamond is a 3-dimensional covalent network structure

2. All atoms are held together by strong covalent bonds

3. These forces require a large amount of energy to break

4. Therefore diamond is very hard.

3-dimensional Covalent network solids are hard

Covalent bond   Carbon atom

Carbon atoms

Weak binding forces

Covalent bond

Structure of Diamond   **Structure of Graphite**

https://www.diamond-jewelry-pedia.com/hardness-of-diamond.html
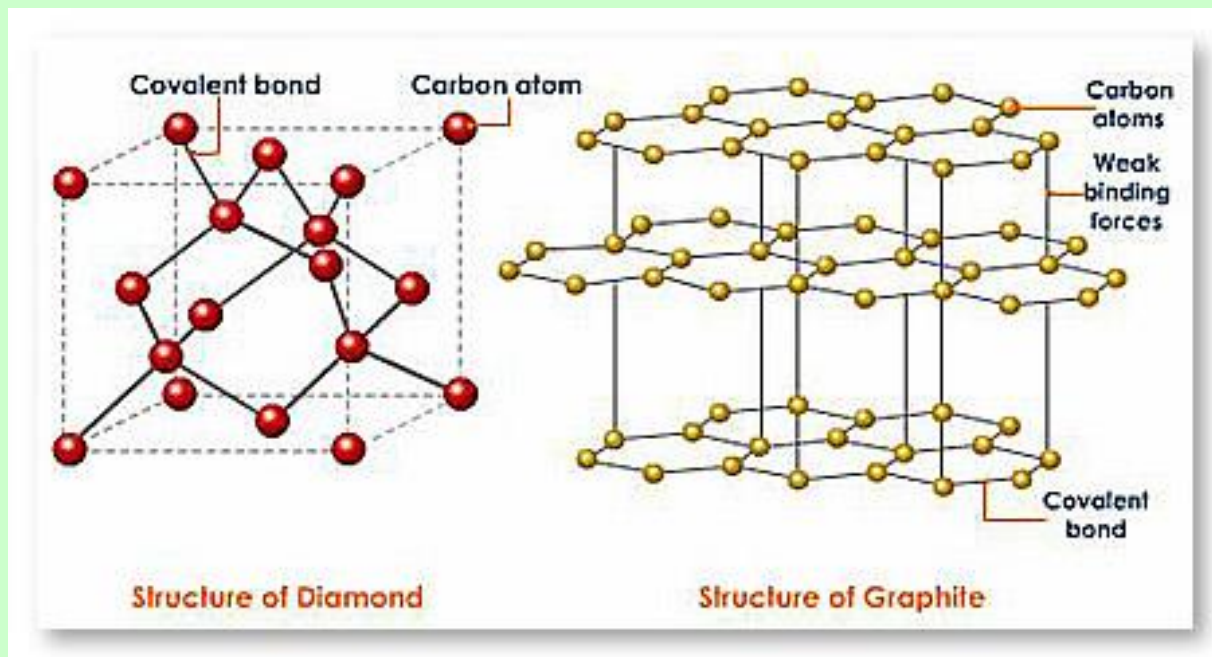
For example:

1. Graphite is a 2-dimensional covalent network structure

2. Atoms are held together by strong covalent bonds in 2-dimensional layers

3. However, the attractive forces holding the layers together are very weak and are broken easily, so the layers easily slide over one another, but the attraction is not strong enough to hold the layers together

4. Therefore, graphite is considered soft.
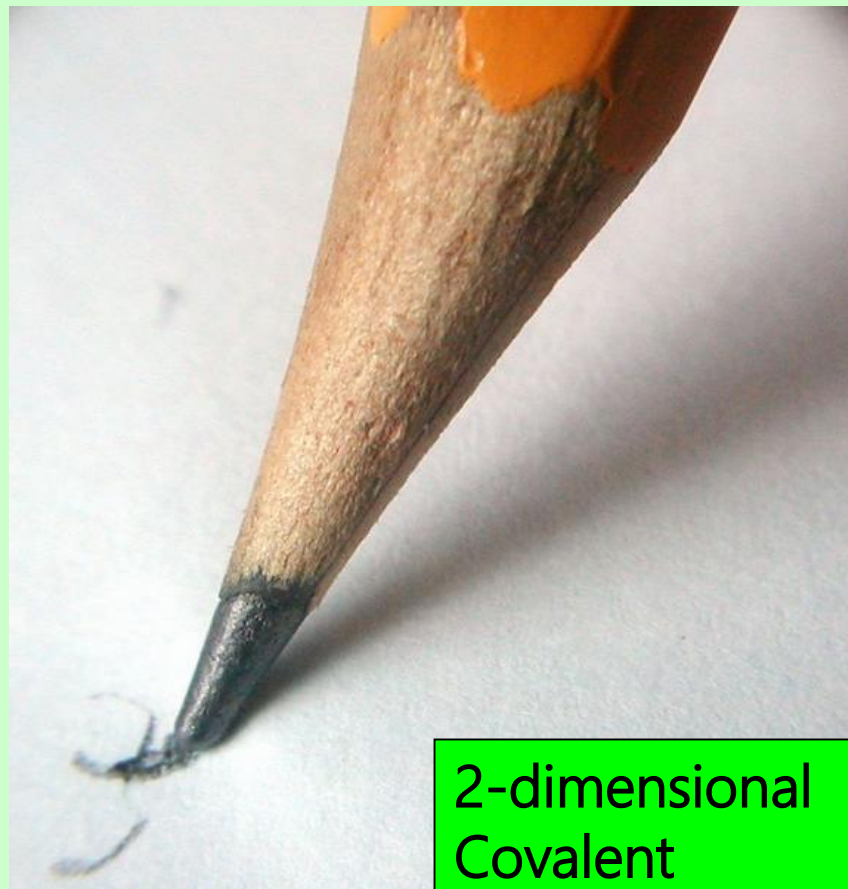
2-dimensional Covalent network solids are soft

Solids Summary

Name of solid substance	Type of particle in solid	Attractive force broken when solid melts	Attractive force between particle – weak or strong (hardness)	Relative melting point	solubility	Electrical conductivity	Malleable
Molecular	molecules	Weak inter molecular	weak	low	Yes if polar No if non-polar	no	no
Metallic	atoms	Metallic bonding	strong	high	no	yes	yes
Ionic	ions	Electrostatic Ionic bonding	strong	high	Yes	Only if molten or in solution	No - brittle
Covalent Network 3-D	atoms	Covalent bonding	strong	high	no	no	no
Covalent Network 2-D	atoms	Covalent bonding	Strong (but weak between layers)	high	no	yes	no

Question 2a: Complete the table below by stating the type of substance, the type of particle, and the bonding (attractive forces) between the particles for each of the substances.

Substance (for example)	Type of substance	Type of particle	Attractive forces between particles
$C_{(s)}$ Graphite	Covalent network	Atom	Covalent (and weak intermolecular forces)
$Cl_{2\ (s)}$ chlorine	Molecular	Molecules	Weak intermolecular forces
$CuCl_{2(s)}$ copper chloride	Ionic	Ion	Ionic bonds / electrostatic attraction
$Cu_{(s)}$ copper	Metal	Atom / cations and electrons	Metallic bonds / electrostatic attraction

This chart needs to be learnt

There will not necessarily be one example for each group but information from this chart MUST be used in following questions about solids

Question 2b: Explain why chlorine is a gas at room temperature, but copper chloride is a solid at room temperature.
In your answer, you should refer to the particles and the forces between the particles in **both** substances.
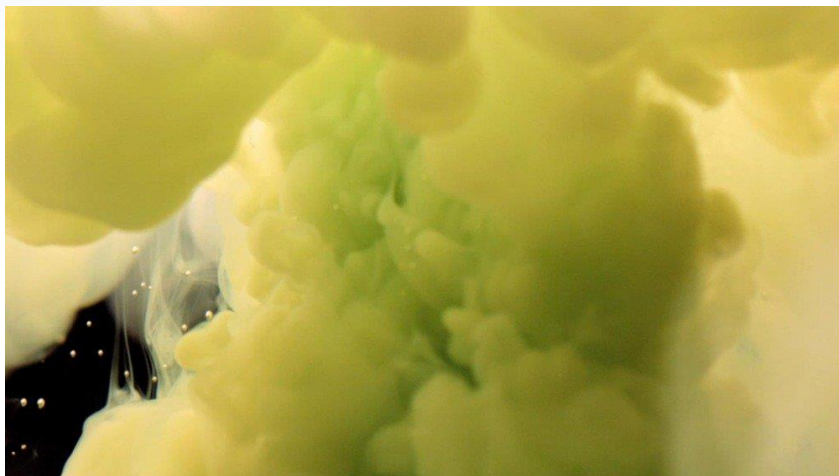
Chlorine is a molecular substance composed of chlorine molecules held together by weak intermolecular forces. The weak intermolecular forces do not require much heat energy to break, so the boiling point is low (lower than room temperature); therefore chlorine is a gas at room temperature.

Copper chloride is an ionic substance. It is composed of a lattice of positive copper ions and negative chloride ions held together by electrostatic attraction between these positive and negative ions. These are strong forces, therefore they require considerable energy to disrupt them and melt the copper chloride; hence copper chloride is a solid at room temperature.

Question 2b (ii) : Using your knowledge of structure and bonding, explain why, although both graphite and copper are good conductors of electricity, copper is suitable for electrical wires, but graphite is not.

For a substance to conduct electricity, it must have charged particles which are free to move.

Graphite is a covalent network solid composed of layers of C atoms covalently bonded to three other C atoms. The remaining valence electron is delocalised (ie free to move) between layers; therefore these delocalised electrons are **able to conduct** electricity.

Copper is a metallic substance composed of copper atoms packed together. Valence electrons are loosely held and are attracted to the nuclei of the neighbouring Cu atoms; ie the bonding is non-directional. These delocalised valence electrons are **able to conduct** an electrical current.
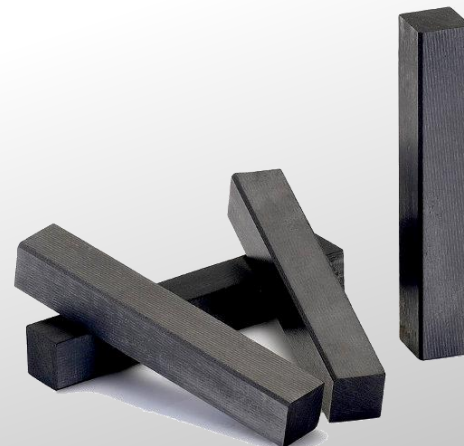
Excellence Question

Question 2b (ii) : Using your knowledge of structure and bonding, explain why, although both graphite and copper are good conductors of electricity, copper is suitable for electrical wires, but graphite is not.

For a substance to be made into wires, it needs to be stretched or drawn out without breaking.

In **graphite**, the attractive forces holding the layers together are very weak and are broken easily, so the layers easily slide over one another, but the attraction is not strong enough to hold the layers together and allow it to be drawn into wires or although the layers can slide due to weak forces, if graphite was to be made into a wire the very strong covalent bonds within the layers would have to be broken

Copper metal is malleable and can easily be drawn into wires since, as it is stretched out, the non-directional metallic bonding holds the layers together, allowing it to be stretched without breaking.

Question 2a: Complete the table below by stating the type of substance, the type of particle, and the type of bonding (attractive forces) between the particles for each of the two substances. Mg (magnesium) and I_2 (iodine)

This chart needs to be learnt

Answer 2a:

	Type of substance	Type of particle	Attractive forces between particles
Mg	Metallic	Atoms / cations and electrons	Metallic bonds / electrostatic attraction between positive ion (cation) and electron
I_2	Molecular	Molecules	Intermolecular forces

There will not necessarily be one example for each group but information from this chart MUST be used in following questions about solids

Identify Metal first then Ionic (metal + non-metal), next Covalent network (C or SiO_2) and then molecular (non-metal + non-metal)

Question 2b: Graphene is a new 2-dimensional material made of carbon atoms. Graphene can be described as a 'one-atom-thick' layer of graphite. A diagram of graphene and two of its properties is shown below. Use your knowledge of structure and bonding to explain the two properties of graphene given above.
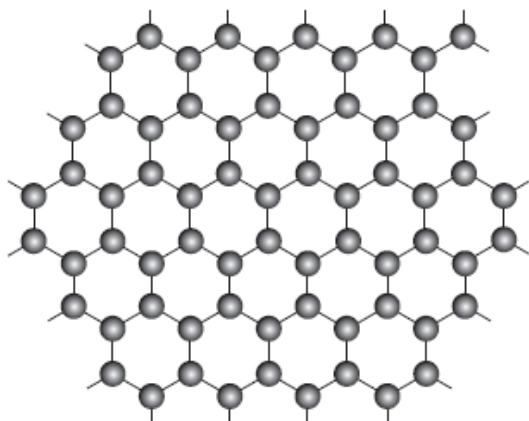
Properties of graphene:

Melting point: very high

Electrical conductivity: excellent

Answer 2b: Graphene has strong covalent bonds. Because the covalent bonds are strong / there are a large number of covalent bonds, it requires a lot of energy to break these bonds, and therefore the melting point is high.

Each carbon atom is bonded to only three other carbon atoms. Therefore each carbon atom has free / delocalised /valence electron(s), to conduct electricity.

.

Question 2c: Solid Mg and I_2 were tested for three physical properties. The table below shows the results of the tests. Use your knowledge of structure and bonding to explain the results of the tests.

Refer back to particle chart

Substance tested	Physical property		
	Ductile	Soluble in cyclohexane (non-polar solvent)	Conducts electricity
Mg	yes	no	yes
I_2	no	yes	no

Answer 2c: Magnesium atoms are held together in a 3–D lattice by metallic bonding in which valence electrons are attracted to the nuclei of neighbouring atoms.
Iodine molecules are held together by weak intermolecular forces.

REMEMBER: (Properties are limited to hardness (including malleability and ductility), electrical conductivity, melting and boiling points and solubility)

Question 2c: Solid Mg and I_2 were tested for three physical properties. The table below shows the results of the tests. Use your knowledge of structure and bonding to explain the results of the tests.

Refer back to particle chart

Substance tested	Physical property		
	Ductile	Soluble in cyclohexane (non-polar solvent)	Conducts electricity
Mg	yes	no	yes
I_2	no	yes	no

Answer 2c: Ductility
The attraction of the Mg atoms for the valence electrons is not in any particular direction (non-directional) ; therefore Mg atoms can move past one another without disrupting the metallic bonding, therefore Mg is ductile.

The attractions between iodine molecules are directional. If pressure is applied the repulsion between like-charged ions will break the solid (brittle) , therefore I_2 is not ductile.

Substance tested	Physical property		
	Ductile	Soluble in cyclohexane (non-polar solvent)	Conducts electricity
Mg	yes	no	yes
I_2	no	yes	no

Refer back to particle chart

Answer 2c: Dissolving in cyclohexane
Magnesium **does not dissolve in cyclohexane** because cyclohexane molecules are not attracted to the magnesium atoms in the metallic lattice.

Iodine **is soluble**, as iodine is a non-polar molecule. The iodine molecules and cyclohexane molecules form weak intermolecular attractions.

Electrical conductivity
Valence electrons of Mg atoms are free to move throughout the structure. This means that magnesium can conduct electricity. There are free moving charge particles.

Iodine does not conduct electricity as it does not contain delocalised electrons (free moving charged particles). Molecules are neutral compounds.

Question 3a: Complete the table below by stating the type of solid, the type of particle, and the attractive forces between the particles in each solid.

This chart needs to be learnt

Answer 3a:

There will not necessarily be one example for each group but information from this chart MUST be used in following questions about solids

Substance	Type of Substance	Type of particle	Attractive forces between particles
Cu(s)	metal / metallic	Atom (or cation and delocalised electrons)	metallic bond
PCl₃(s)	molecular	molecule	intermolecular (forces)
SiO₂(s)	covalent network	atom	covalent bond
KCl(s)	ionic	ion	ionic bond

Identify Metal first then Ionic (metal + non-metal), next Covalent network (C or SiO$_2$) and then molecular (non-metal + non-metal)

Question 3b: Phosphorus trichloride, PCl_3, is a liquid at room temperature, and does not conduct electricity.
Explain these two observations in terms of the particles, structure, and bonding of PCl_3.

Refer back to particle chart

Answer 3b: Phosphorus trichloride, PCl_3, is a **molecular solid**, made up of non-metal phosphorus and chlorine atoms **covalently** bonded together. The molecules are held together by **weak intermolecular forces**. Since these forces are weak, not much energy is required to overcome them, resulting in low melting / boiling points. (In the case of PCl_3, its melting point is lower than, and its boiling point is higher than room temperature, so it is liquid.)

PCl_3 does not contain free moving ions nor any delocalised / free moving valence electrons, meaning PCl_3 **does not contain any charged particles**. Since free moving charged particles are required to carry electrical current, PCl_3 is unable to conduct electricity.

REMEMBER: (Properties are limited to hardness (including malleability and ductility), electrical conductivity, melting and boiling points and solubility)

Question 3c: Consider each of the solids copper, Cu, silicon dioxide, SiO_2, and potassium chloride, KCl.
Complete the table below by identifying which of these solids have the listed physical properties:

Refer back to particle chart

Physical properties	Solid
The solid is insoluble in water and is malleable.	METAL - copper
The solid is soluble in water and is not malleable.	IONIC – potassium chloride
The solid is insoluble in water and is not malleable.	COVALENT NETWORK – silicon dioxide

REMEMBER: Draw up a quick chart for each physical property and each solid

Excellence Question

Answer 3c: Justify TWO of your choices in terms of the particles, structure, and bonding of these solids. You may use diagrams in your justification.

State the properties and solid type

Refer back to particle chart

Cu is **insoluble in water and malleable**.

1. Copper is a metal made up of an array of atoms (or ions) held together by **non-directional forces** between the positive nuclei of the atoms and the delocalised / free moving valence electrons.

2. There is no attraction between the copper atoms and the (polar) water molecules, therefore Cu is **insoluble in water**.

3. Since the attractive forces are non-directional, when pressure is applied, the Cu atoms can move past each other to change shape without the bonds breaking, so Cu **is malleable**. (Note – labelled diagrams can provide replacement evidence).

REMEMBER: discuss solid structure first then each of the properties. You need both properties for full marks.

Excellence Question

Question 3c: Justify TWO of your choices in terms of the particles, structure, and bonding of these solids. You may use diagrams in your justification.

Refer back to particle chart

KCl is **soluble in water and not malleable**.

1. KCl is made up of positive K^+ ions, and negative Cl^- ions, ionically bonded in a 3D lattice.

2. When added to water, polar water molecules form electrostatic attractions with the K^+ and Cl^- ions. The partial negative charge, δ^-, on oxygen atoms in water are attracted to the K^+ ions and the partial positive, δ^+, charges on the H's in water are attracted to the Cl^- ions, causing KCl to **dissolve in water**.

3. KCl is **not malleable** because if pressure is applied to an ionic lattice, it forces ions with the same charge next to each other; they repel each other and break the structure.
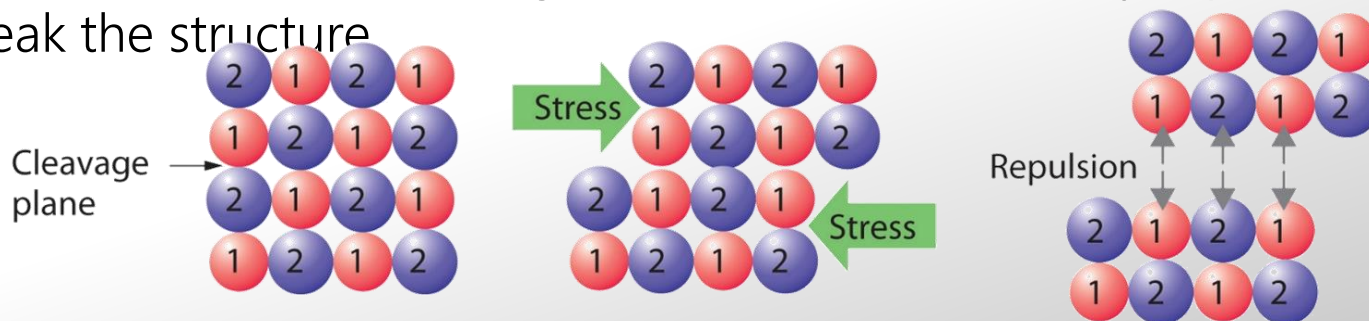
Excellence Question

Question 3c: Justify TWO of your choices in terms of the particles, structure, and bonding of these solids. You may use diagrams in your justification.

Refer back to particle chart

SiO_2 is insoluble in water and not malleable.

1. SiO_2 is a **covalent network** made up of atoms covalently bonded together in a 3D lattice structure.

2. (Covalent bonds are strong), Polar water molecules are not strong / insufficiently attracted to the Si and O atoms, therefore **SiO_2 is insoluble** in water.

3. SiO_2 is not malleable because if pressure is applied, the directional / strong covalent bonds have to be broken before the atoms can move.

(Note - labelled diagrams can provide replacement evidence).
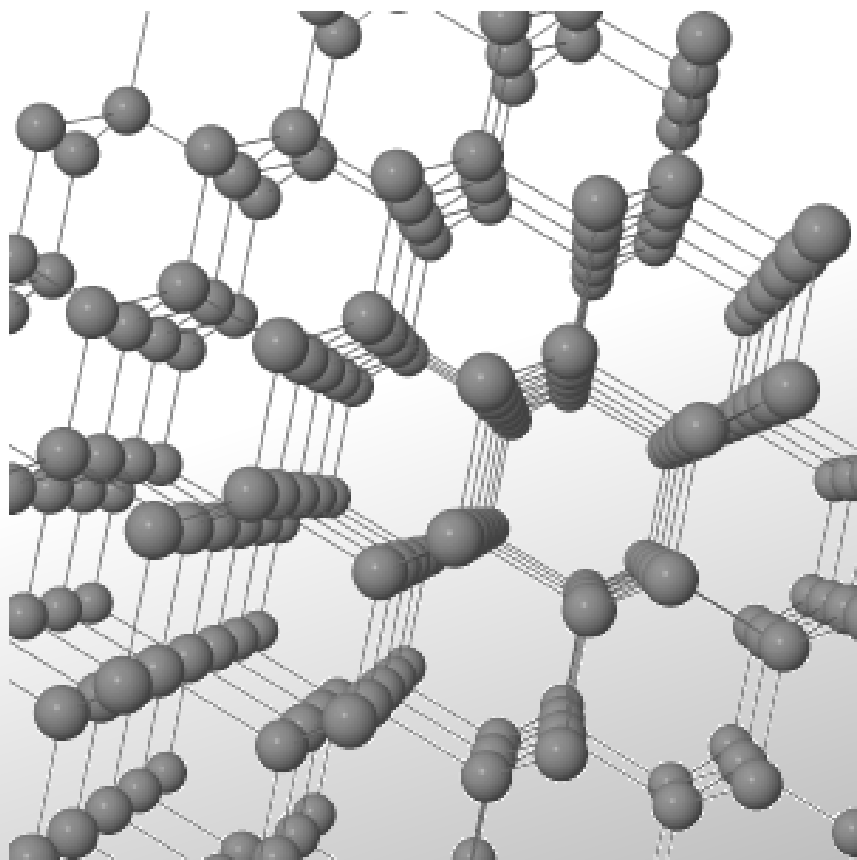
Question 2a: Complete the table below by stating the type of substance, the type of particle, and the attractive forces between the particles in the solid for each substance.

This chart needs to be memorised

Answer 2a:

Substance	Type of substance	Type of particle	Attractive forces between particles
$ZnCl_2(s)$ (zinc chloride)	ionic	ions	ionic
$C(s)$ (graphite)	covalent network	atoms	covalent
$CO_2(s)$ (carbon dioxide / dry ice)	molecular	molecules	intermolecular

There will not necessarily be one example for each group but information from this chart MUST be used in following questions about solids

Identify Metal first then Ionic (metal + non-metal), next Covalent network (C or SiO_2) and then molecular (non-metal + non-metal)
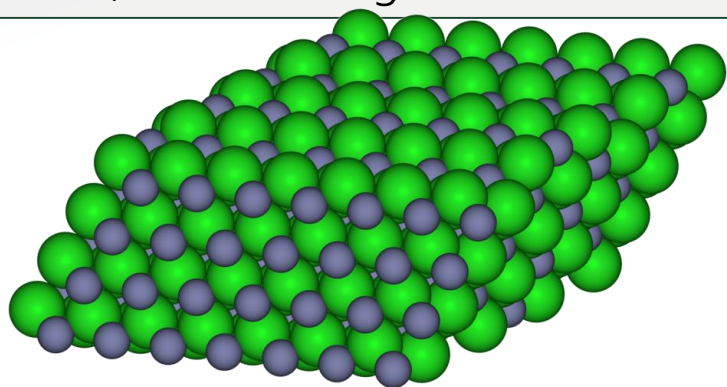
Question 2b : Carbon (graphite) conducts electricity when it is solid, whereas zinc chloride, $ZnCl_2$, will not conduct electricity when solid, but will conduct when molten.
Justify this statement in terms of the particles, structure, and bonding for both substances.

For a substance to conduct electricity, it must have charged particles which are free to move.

Graphite is a covalent network solid composed of layers of C atoms covalently bonded to three other C atoms. The remaining valence electron is delocalised (ie free to move) between layers; therefore these delocalised electrons are **able to conduct** electricity.

$ZnCl_2$ is an ionic compound that cannot conduct electricity when solid because the ions (charged particles) are fixed in place in a 3D lattice structure and unable to move. When molten, the ionic bonds between the ions break, so the ions are free to move in the molten liquid. With charged particles / ions free to move, $ZnCl_2$ can then conduct electricity.

Excellence
Question

Question 2c : Solid zinc chloride, $ZnCl_{2(s)}$, is soluble in water. Dry ice, $CO_{2(s)}$, is not readily soluble in water. Justify these statements in terms of the particles, structure, and bonding of these substances.

For a substance to be soluble Polar water molecules need to be insufficiently attracted to the particles of the substance placed in

Polar water molecules attract the ions in **zinc chloride's** 3-D lattice strongly enough to separate and dissolve them. The negative charges on the oxygen ends of the water molecules are attracted to the positive Zn^{2+} ions, and the positive hydrogen ends of the water molecules are attracted to the negative Cl^- ions, forming hydrated ions that can spread out through the solution. Therefore it is soluble.

The polar water molecules are unable to interact with the non-polar **carbon dioxide** molecules strongly enough to break the intermolecular forces between the carbon dioxide molecules. Therefore it is not soluble.

Question 3a : Complete the table below by stating the type of solid, the type of particle, and the type of bonding (attractive forces) between the particles in each solid.

Solid	Type of solid	Type of particle	Attractive forces between particles
$Al(s)$ (Aluminium)	*metal / metallic*	*atoms (or cations and delocalised valence electrons)*	*metallic (bonds)*
$MgCl_2(s)$ (Magnesium chloride)	*ionic compound*	*ions*	*ionic (bonds)*
$S_8(s)$ (Sulfur)	*molecular*	*molecules*	*intermolecular (bonds)*

Excellence Question

Question 3b: Circle the substance which has the lowest melting point.
$Al_{(s)}$ $MgCl_{2(s)}$ $S_{8(s)}$
Justify your choice, referring to the attractive forces between the particles of ALL three substances.

Sulfur has the lowest melting point.
Sulfur is a molecular substance with weak intermolecular forces between the molecules.
These forces do not require much energy to overcome, so they will break at lower temperatures, giving sulfur a lower melting point.
Al is a metal with strong metallic bonds. These attractions require a lot of energy to overcome, so the melting point is higher than sulfur's melting point.
$MgCl_2$ is an ionic compound with strong ionic bonds between the cations and anions.
These bonds also require a lot of energy to overcome, so the melting point is also higher than sulfur's melting point.

Question 3c: Circle the substance which is malleable.

$Al_{(s)}$ $MgCl_{2(s)}$ $S_{8(s)}$

Justify your choice by referring to the structure and bonding of your chosen substance.

You may include a diagram or diagrams in your answer.

Aluminium is malleable.

Aluminium is a metal made up of atoms / cations in a sea of electrons which are held together by non-directional metallic bonds in a (3D) lattice. The metallic bonds are non-directional as the (bonding) electrons are delocalised across the lattice / shared by many atoms. When a force (or pressure) is applied the atoms / layers can move without breaking / disrupting these non-directional bonds thus the structure can change shape without breaking the lattice.

Achieved Question

Question 3a. Complete the table below by choosing the appropriate type of solid that matches the properties shown in the table.
Types of solid: Ionic, Metallic, Covalent Network, Molecular.

Solid	Melting point (°C)	Boiling point (°C)	Conducts electricity?	Soluble in water?	Type of solid
A	290	732	solid – no molten – yes	Yes, solution conducts electricity	Ionic
B	44	280	No	No	Molecular
C	1710	2230	No	No	Covalent Network
D	660	2470	Solid and molten – yes	No	Metallic

https://www.learnit.ru/476728/

Question 3b. Explain why Solid **A** does not conduct electricity in the solid state, but will conduct when molten or when dissolved in water.
Refer to the particles, structure, and bonding of this substance.

Electrical conductivity requires a substance to have **mobile (free moving) charged particles**.
Solid A is an ionic solid made up of a 3-D lattice of positive and negative ions (cations and anions) that are attracted to each other.
❑ In the **solid** state, these ions are **rigidly held in a lattice by strong ionic bonds**, so cannot move around.
❑ When **molten**, the ions are able to move freely so it can conduct electricity.
❑ In **aqueous solution**, the ions are also free to move so the solution can also conduct electricity.

Solid	Melting point (ºC)	Boiling point (ºC)	Conducts electricity?	Soluble in water?	Type of solid
A	290	732	solid – no molten – yes	Yes, solution conducts electricity	Ionic

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Excellence
Question

Question 3c. Elaborate on the differences in the melting points of solids **B** and **D** with reference to their particles, structure, and bonding.

Solid B is composed of discrete covalent molecules which are held together by weak intermolecular forces. These weak intermolecular forces are easily broken, so the molecules can be separated with little energy, therefore the melting point is low.

Solid D is a metal made of a 3-D lattice of metal atoms surrounded by a sea of delocalised valence electrons, which are strongly attracted to all the nuclei in the lattice. This forms a strong metallic bond which requires a large amount of energy to break, therefore the melting point is high at 660°C.

Solid	Melting point (°C)	Boiling point (°C)	Conducts electricity?	Soluble in water?	Type of solid
A	290	732	solid – no molten – yes	Yes, solution conducts electricity	Ionic
B	44	280	No	No	Molecular
C	1710	2230	No	No	Covalent Network
D	660	2470	Solid and molten – yes	No	Metallic

Question 1a. Complete the table below by stating the type of solid, the type of particle, and the type of bonding (attractive forces) between the particles in each solid.

Solid	Type of solid	Type of particle	Attractive forces between particles
$Na(s)$ (sodium)	Metal / metallic	Atoms/cations in a sea of delocalised electrons	Metallic bond
$NaI(s)$ (sodium iodide)	ionic	ions	Ionic bond / electrostatic attraction between (oppositely charged) ions
$I_2(s)$ (iodine)	(covalent) molecular	molecules	(weak) intermolecular (forces)

Question 1b. Sodium, Na(*s*), is malleable, whereas sodium iodide, NaI(*s*), is brittle. Explain these observations by referring to the structure and bonding of each substance.

Sodium is a metallic solid made up of atoms in 3D lattice held together by nondirectional metallic bonds (or cations non-directionally electrostatically attracted to a surrounding sea of electrons). When a force (or pressure) is applied, the atoms / layers can move without breaking / disrupting these non-directional bonds; thus the structure can change shape.

NaI is made up of alternating positive ions / Na$^+$ ions, and negative ions / I$^-$ ions, ionically bonded in a 3D lattice. NaI is not malleable because if pressure is applied to an ionic lattice, it forces ions with the same charge next to each other; they repel each other and break the structure.

Question 3a. Classify the following chemical process as exothermic or endothermic and give a reason for your choice.

(i) $C_{(s)} + O_{2(g)} \rightarrow CO_{2(g)}$ $\Delta_rH = -394$ kJ mol^{-1}
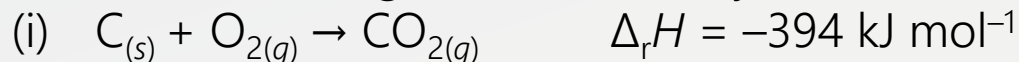
(ii) In the reaction above, $C_{(s)}$ in the form of graphite can conduct electricity. The product, carbon dioxide, $CO_{2(g)}$, does not conduct electricity.

Use your knowledge of structure and bonding to explain this observation.

Exothermic as Δ_rH is negative.

To be able to conduct electricity, there needs to be mobile/free moving charged particles. Graphite, $C_{(s)}$, is an extended covalent network solid. Each carbon atom is covalently bonded to 3 other carbon atoms in hexagonal layers. This leaves one delocalised electron per carbon atom that is mobile and able to carry a charge, so graphite conducts electricity.

Carbon dioxide is a covalent molecule. The molecules are held together by weak intermolecular forces, so it is a gas at room temperature. There are no free moving ions or electrons in their structure. Therefore, it can't conduct electricity.

Enthalpy (or Heat Content) is the energy in a substance due to kinetic energy of particles and potential energy in chemical bonds.

Enthalpy change ΔH is the difference in enthalpy of products H_P and reactants H_R

Enthalpy increases

Heat energy IN

Enthalpy decreases

Heat energy OUT
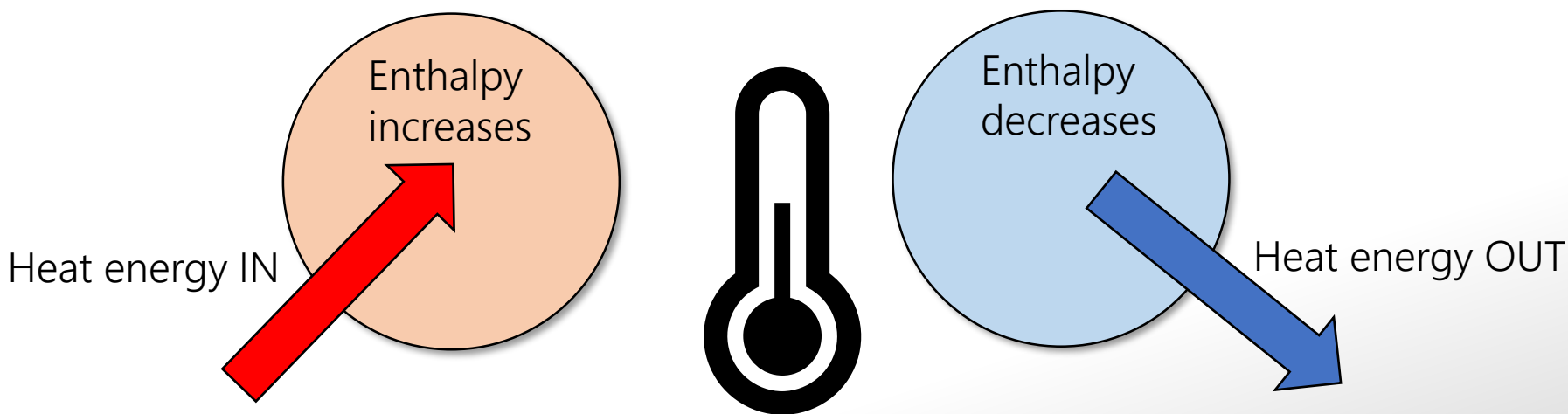
$$\Delta H = H_P - H_R$$

The unit for Enthalpy is kilojoules (kJ)

H_P (products) and H_R (reactants) cannot be measured directly.

We can measure Enthalpy change (ΔH) by measuring energy;

Released to surroundings (Exothermic Reactions)	Absorbed from surroundings (Endothermic Reactions)
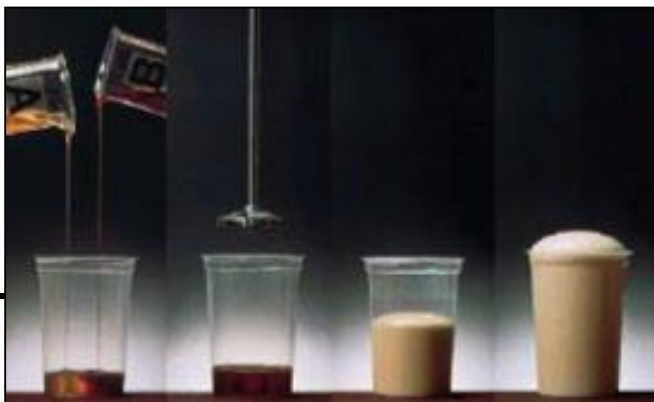
Exothermic Reactions	Endothermic Reactions
These are reactions where **heat energy** is **released** into the surroundings.	These are reactions where **heat energy** is **absorbed** from the surroundings.
Surroundings gain heat energy. (increase in temperature)	Surroundings lose heat energy. (Decrease in temperature)
Products will have less energy than reactants.	Products will have more energy than reactants.
ΔH is NEGATIVE (-)	ΔH is POSITIVE (+)
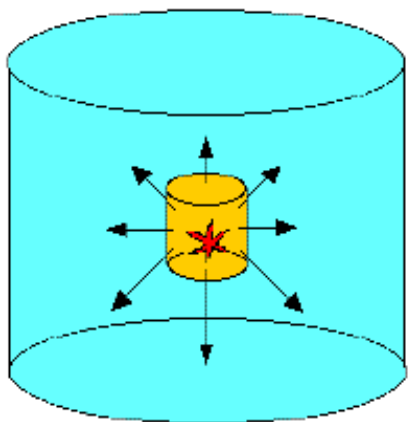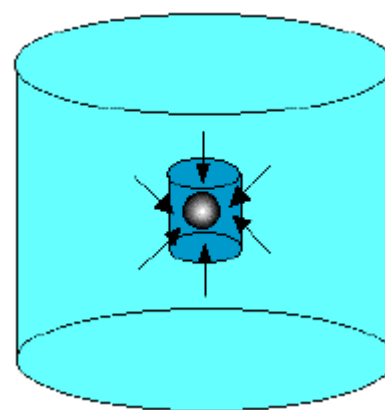
Exothermic reactions

Any combustion reaction is **exothermic**. The bonds holding the atoms of fuel molecules together (usually consisting of carbon and hydrogen atoms) release a lot of energy in the form of light and heat when they are broken. The total energy holding the bonds together in the products are less than the total energy in the reactions and the difference is released.
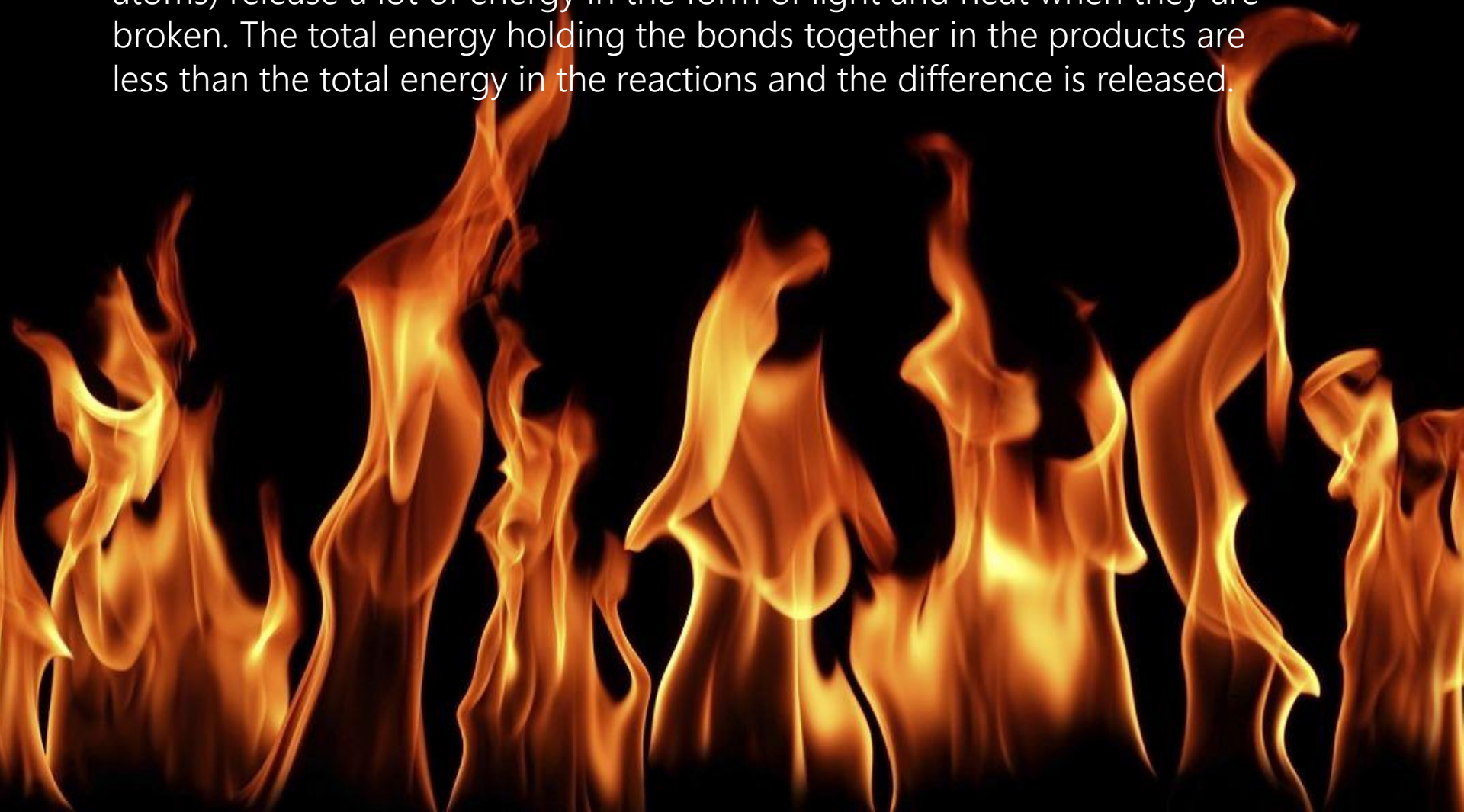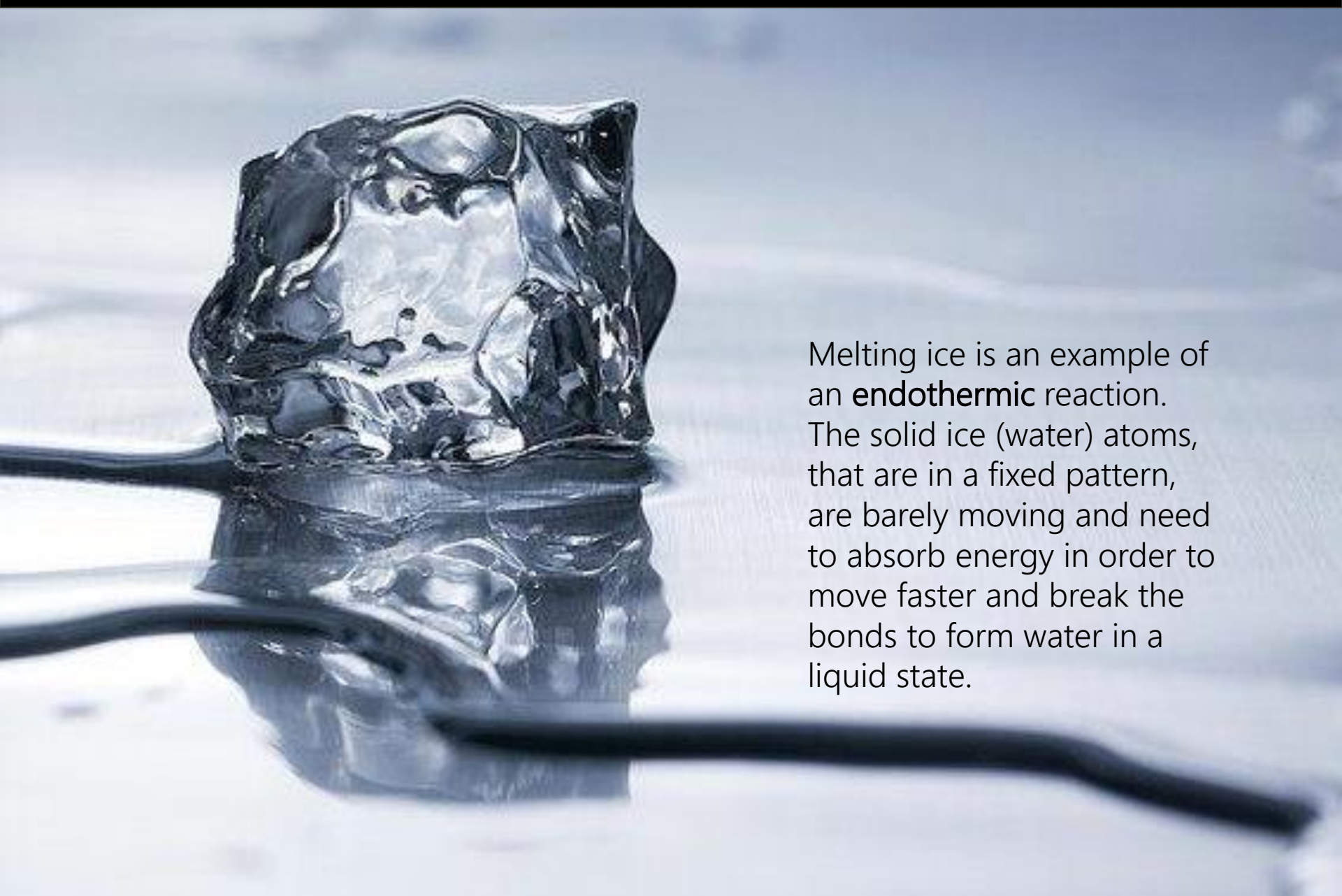
Melting ice is an example of an **endothermic** reaction. The solid ice (water) atoms, that are in a fixed pattern, are barely moving and need to absorb energy in order to move faster and break the bonds to form water in a liquid state.

Bonds forming between atoms and molecules release energy therefore **bond forming is an exothermic reaction.** Bonds are formed to form a stable molecule.

Bonds breaking between atoms and molecules require energy therefore **bond breaking is an endothermic reaction.** The input of energy (usually light or heat energy) cause the atoms and molecules to move faster and 'pull away' from each other. Each type of bond has its own specific amount of energy, called *bond enthalpy* measured in kJ, required to break its bond.

Gas

condensation

evaporating
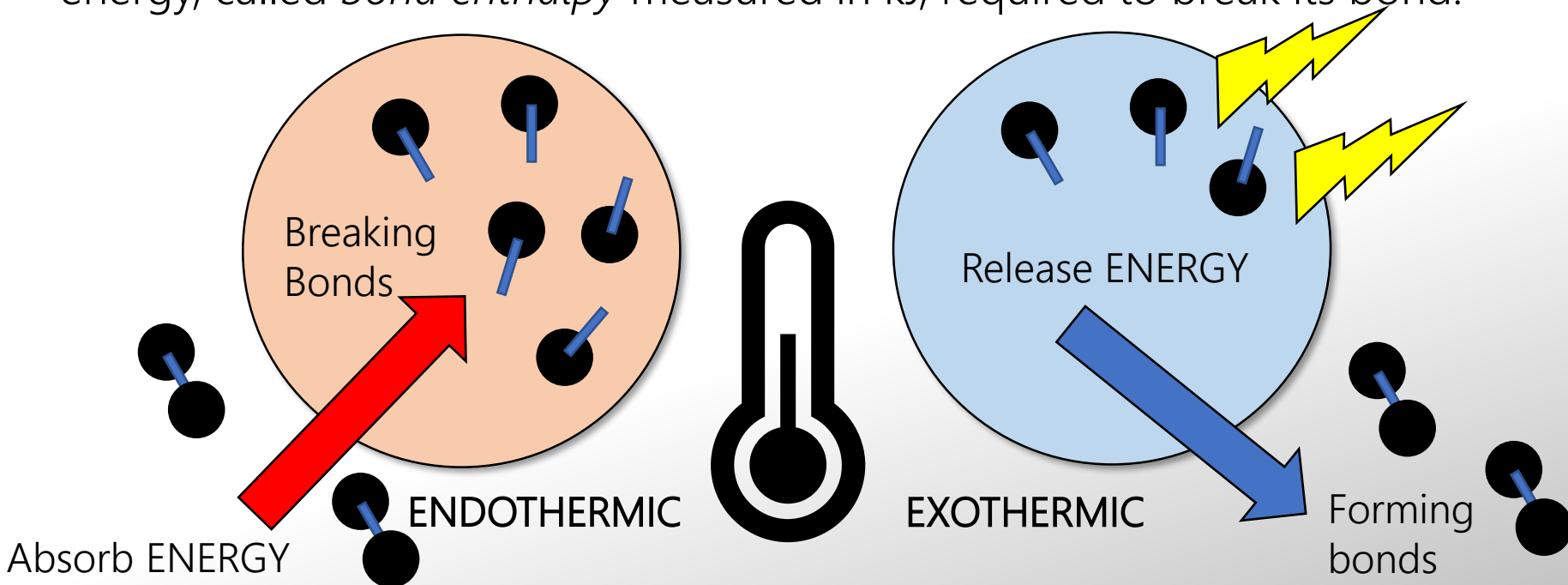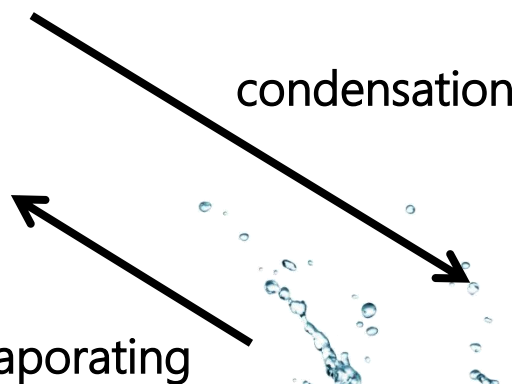
sublimation

deposition

Liquid

Freezing

Solid

Melting

If energy is absorbed or released the particles which make up the matter can change state. A change of state is a physical reaction and it is reversible.

Solid particles are packed closely and only vibrate in a fixed position. Liquid particles are also packed closely but the particles move around more. Gas particles have a lot of space between them and move around quickly.
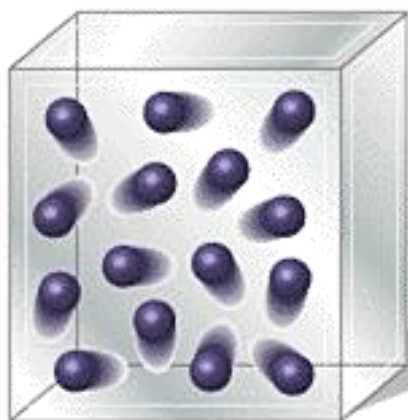
Gas   Liquid   Solid

Cool or compress

Heat or reduce pressure

Cool

Heat

Total disorder; much empty space; particles have complete freedom of motion; particles far apart.

Disorder; particles or clusters of particles are free to move relative to each other; particles close together.

Ordered arrangement; particles are essentially in fixed positions; particles close together.

Particles of different states have different kinetic energy levels

Kinetic energy causes particles to move. The more kinetic energy a particle has the faster it moves. Kinetic energy can be added to a particle by adding heat energy (and heats up). The heat energy is then **transformed** into the kinetic energy.
Kinetic energy can also be lost from a particle, which slows it down, when it changes back into heat energy and is lost (and cools).
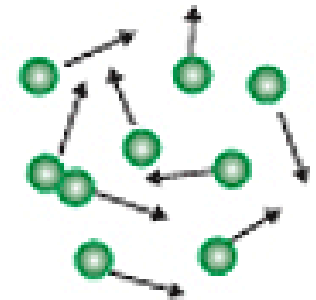
Energy, Heat   Energy, Heat
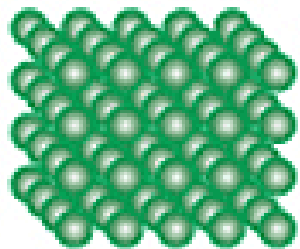
The first state: Solid
(i.e. Ice)

Maintains formation with
a little vibration only
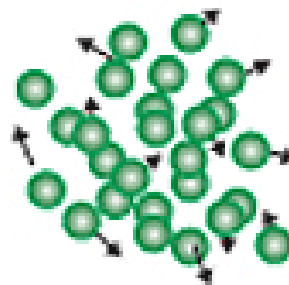
The second state: Liquid
(i.e. water)

Movement of molecules or
atoms become free and
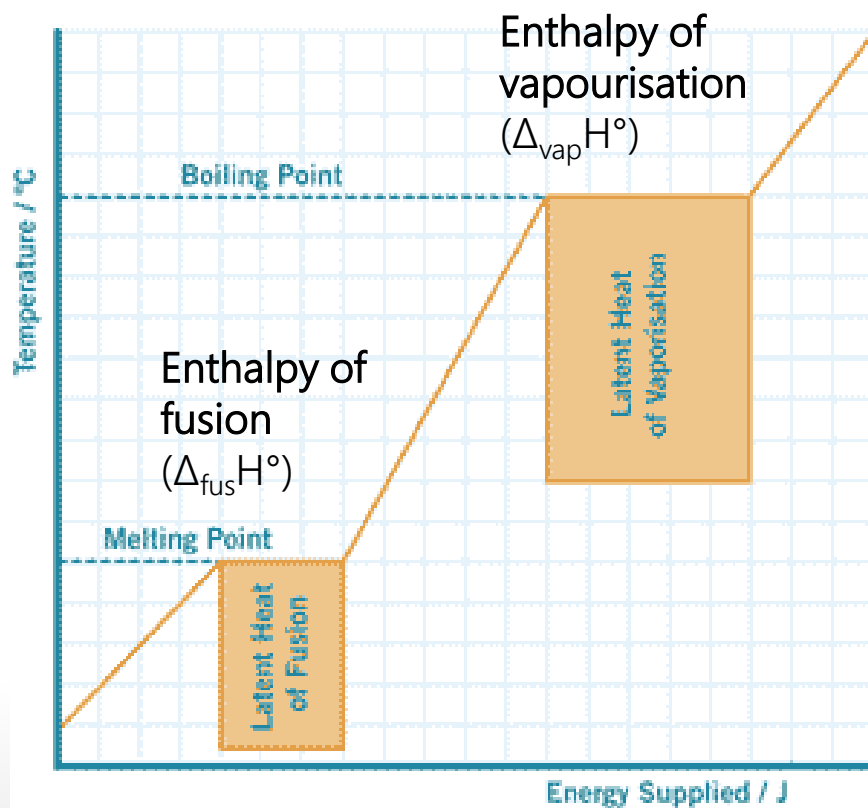disorderly

The third state: Gas
(i.e. vapor)

Movement of molecules
and electrons become
further freer

When heat energy is added to a solid substance at a particular temperature called the melting point, it will change state into a liquid.

Prior to this point a rise in heat energy will also show a rise in temperature. At the melting point the heat energy will be used to break the bonds in a solid – an **endothermic** reaction called **latent heat of fusion-** rather than show a temperature increase.

The same occurs at the boiling point from a liquid to a gas – an endothermic reaction called **latent heat of vaporisation.**

Enthalpy of vapourisation ($\Delta_{vap}H°$)

Boiling Point

Temperature / °C

Enthalpy of fusion ($\Delta_{fus}H°$)

Latent Heat of Vaporisation

Melting Point

Latent Heat of Fusion

Energy Supplied / J

If **more energy is released** when water bonds to the solute than it takes to separate the solute, the dissolving is **exothermic** and the temperature increases. An example is adding a strong acid (such as sulfuric acid) or base (such as sodium hydroxide)

However, for some substances dissolving the reaction is **endothermic**, for example Potassium chloride in water. More energy is needed to break the bonds between this ionic salt than released when new bonds are formed with the ions and water molecules.
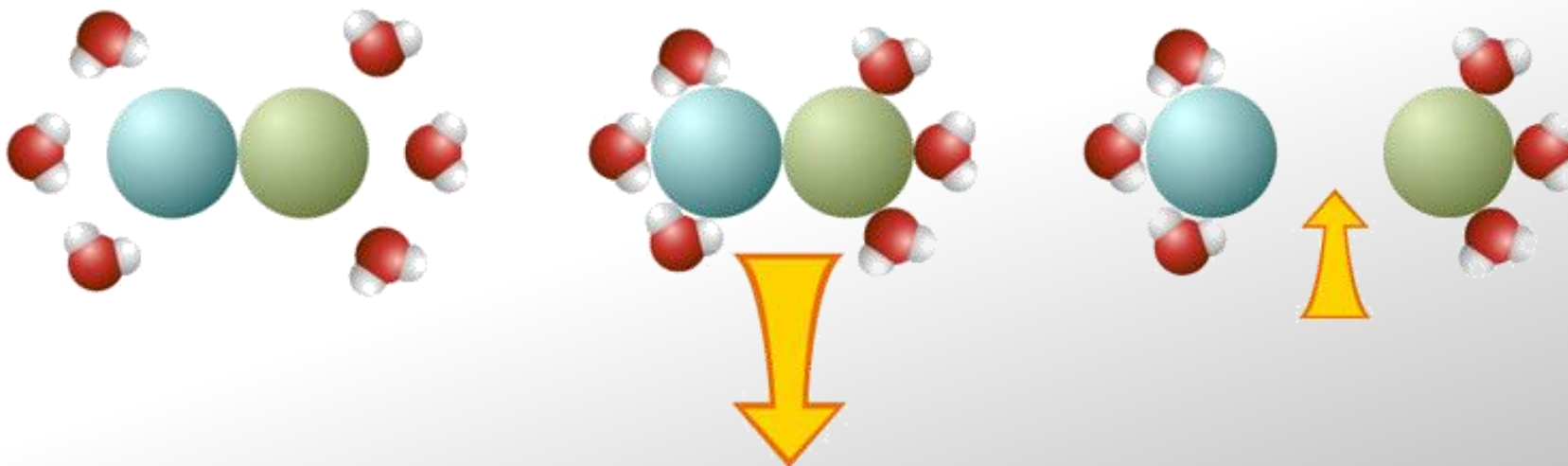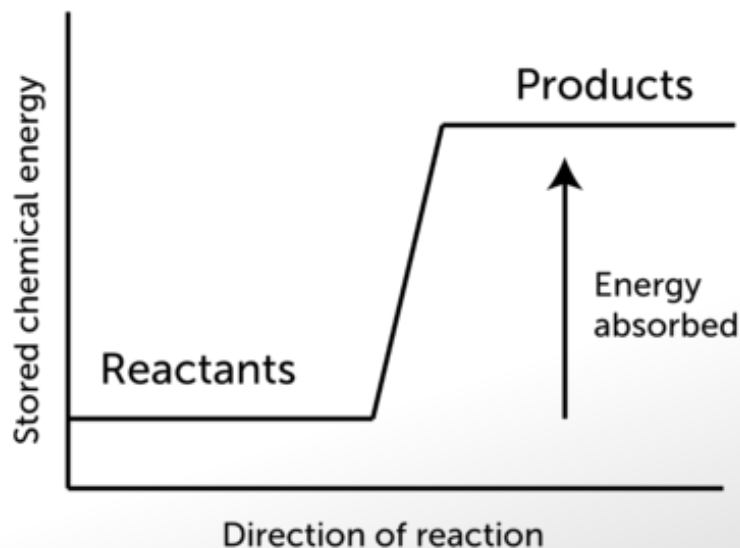
Enthalpy Diagrams can be used to show the **relative amounts of enthalpy** of the reactants and products in a reaction, as well as the direction and relative size of enthalpy change. In endothermic reactions, the enthalpy change will be positive and in exothermic reactions, the enthalpy change will be negative.
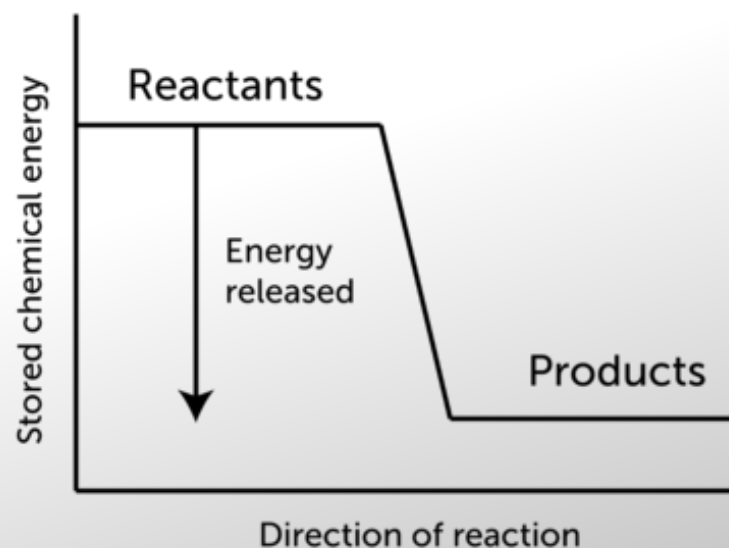
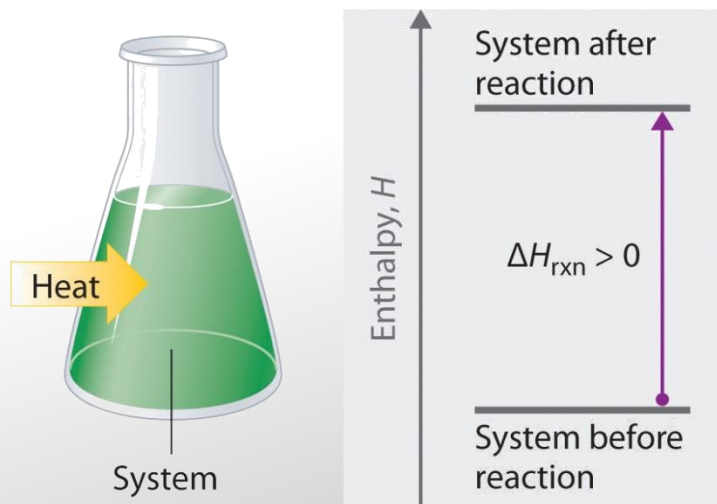| Endothermic Reactions	Exothermic Reactions
e.g. Reacting methane with steam at high pressure and temp. Energy is absorbed	e.g. Burning of methane in air. Energy is released

An exothermic reaction will release energy and the products will be at a lower enthalpy level than the reactants. The reaction system will feel will feel hot to the touch as the energy is released as heat energy.

An endothermic reaction will absorb energy and the products will be at a higher enthalpy than the reactants. The reaction system will feel cool to the touch as heat energy is taken from the surroundings, including your skin, and used to break bonds in the molecules.

$\Delta_r H$ is the enthalpy of a reaction and is measured in $kJmol^{-1}$
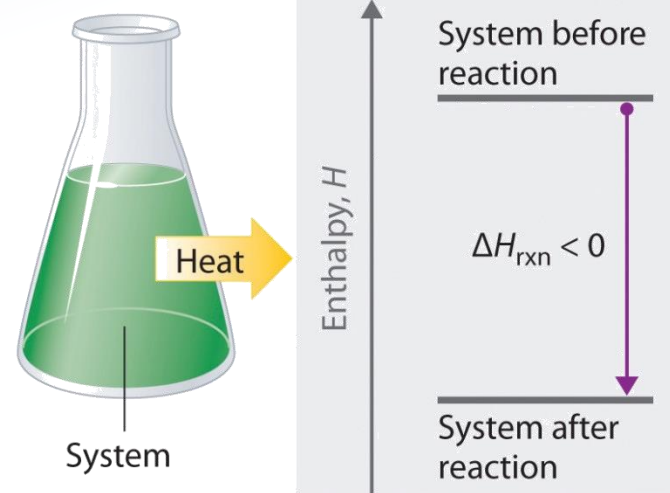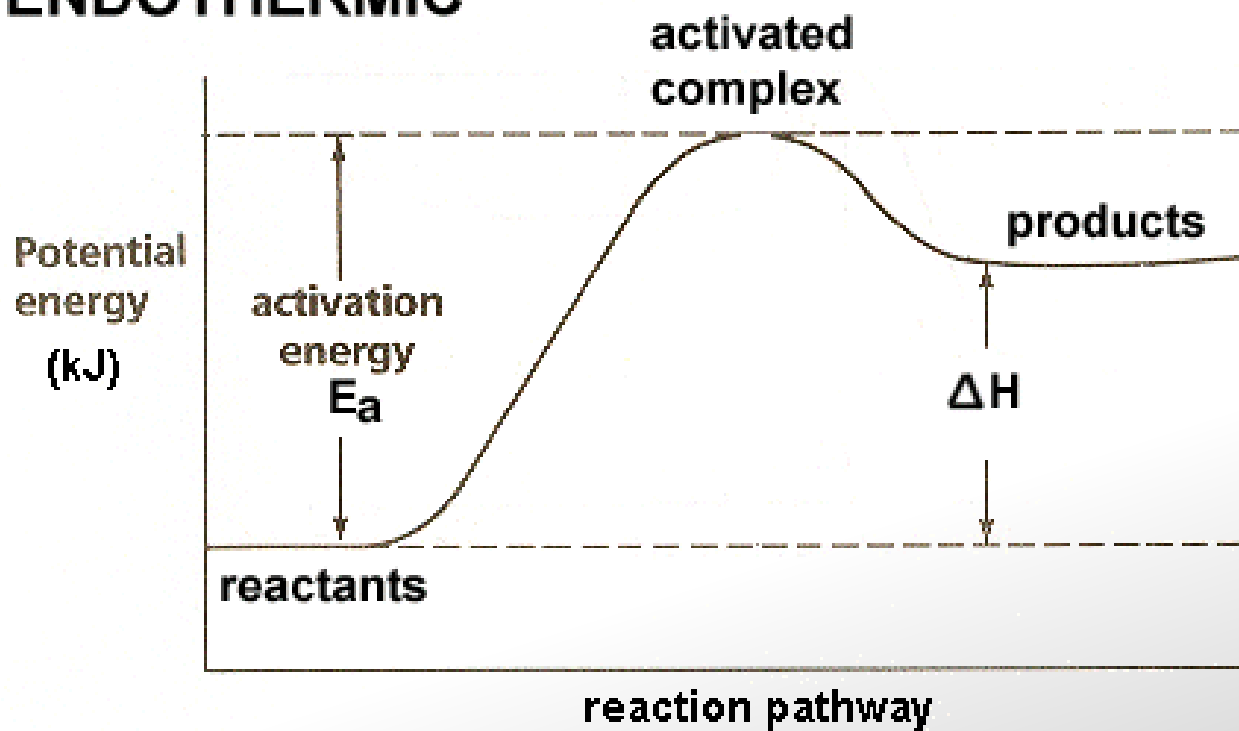

(a) Exothermic reaction

(b) Endothermic reaction

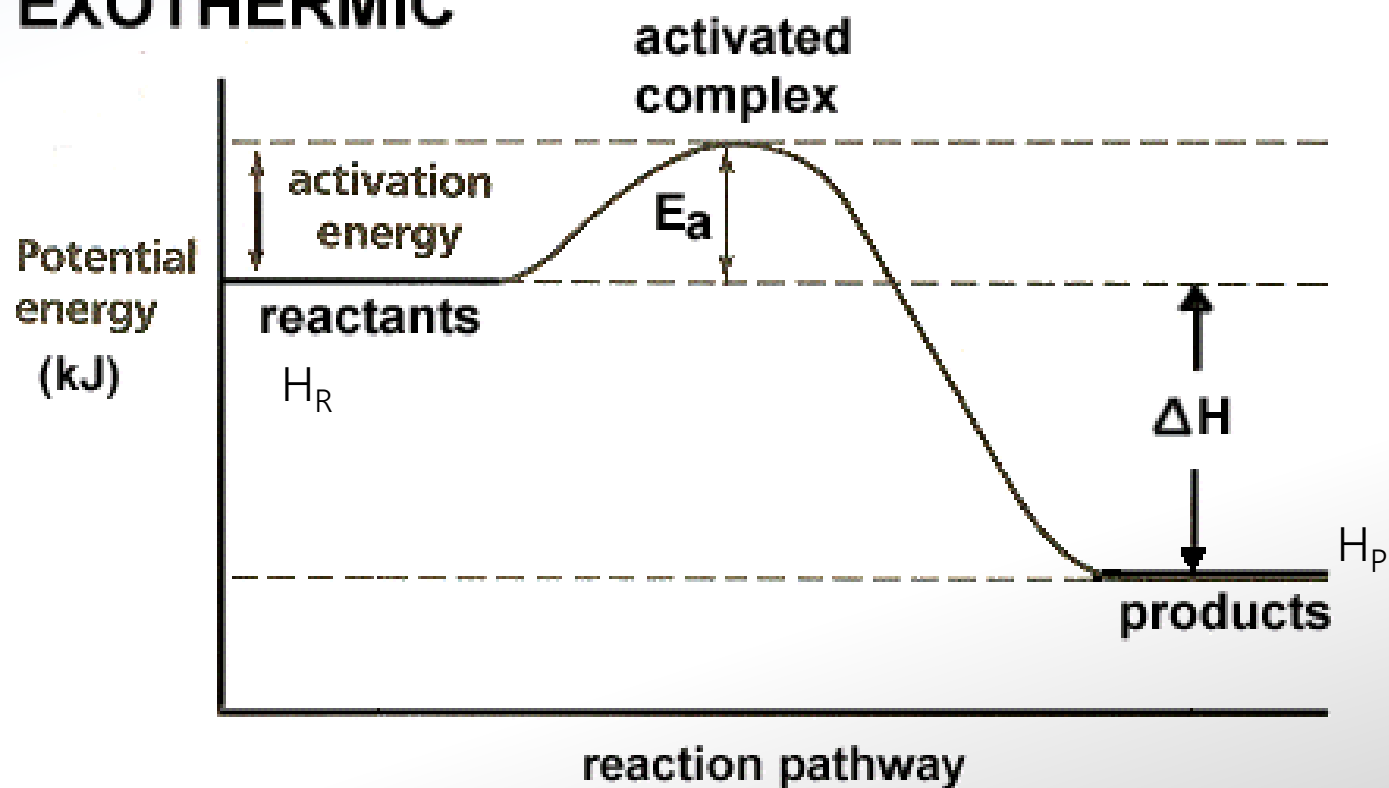Endothermic Reaction e.g. Reacting methane with steam at high pressure and temp. Energy is absorbed

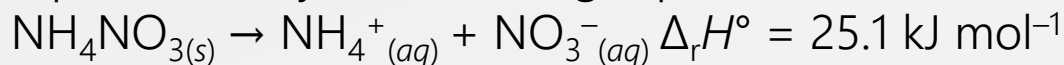Exothermic Reaction e.g. concentrated Hydrochloric acid reacting with zinc metal

Merit Question

Question 3a: Dissolving ammonium nitrate in a beaker containing water can be represented by the following equation:

$$NH_4NO_{3(s)} \rightarrow NH_4^+{}_{(aq)} + NO_3^-{}_{(aq)} \quad \Delta_rH° = 25.1 \text{ kJ mol}^{-1}$$
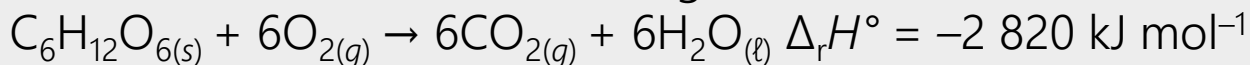
Give the term below that best describes this process and give the description that best describes what you would observe happening to the beaker during this process.

Answer 3a: Endothermic - Gets colder
The process is endothermic since the enthalpy change ($\Delta rH°$) is positive, which indicates that energy is absorbed by the system as the ammonium nitrate dissolves. Since heat energy is absorbed by the system from the surroundings (water & beaker), the water or beaker will get cooler as they lose heat energy.

Heat absorbed is always endothermic

Question 3b: Glucose is an important source of energy in our diet. The equation below shows the combustion of glucose to form carbon dioxide and water.

$$C_6H_{12}O_{6(s)} + 6O_{2(g)} \rightarrow 6CO_{2(g)} + 6H_2O_{(\ell)} \quad \Delta_rH° = -2\ 820 \text{ kJ mol}^{-1}$$

Give the term below that best describes this process and give a reason

Answer 3b: Exothermic
The reaction is exothermic because the enthalpy change ($\Delta rH°$) is negative; indicating that heat energy is produced during the reaction.
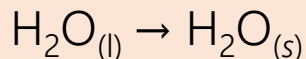
Question 3a (i): When solid sodium hydroxide is added to water, the temperature increases.
- Identify the term that best describes this reaction
- Give a reason for your choice

Answer 3a (i) : Exothermic, as the temperature increases, which shows energy is being released.

Heat released is always exothermic

Question 3a(ii): The freezing of water to form ice can be represented by the following equation.
$H_2O_{(l)} \rightarrow H_2O_{(s)}$
- Identify the term that best describes this reaction
- Give a reason for your choice

Answer 3a(ii): Exothermic, weak intermolecular attractions form between the water molecules, this releases energy.

Achieved Question

Question 2a: Hand warmers contain a supersaturated solution of sodium ethanoate which, when activated, crystallises and releases heat.
- Identify the term that best describes this reaction
- Give a reason for your choice

Heat released is always exothermic

Answer 2a: Exothermic because the temperature of the solution increases / heat is released / particles slow down / bonds are formed

Question 2b(i): Glucose is made in plants during photosynthesis when carbon dioxide gas, $CO_{2(g)}$, and water, $H_2O_{(l)}$, react to produce glucose, $C_6H_{12}O_{6(aq)}$, and oxygen gas, $O_{2(g)}$. The photosynthesis reaction can be represented by the following equation:
$$6CO_{2(g)} + 6H_2O_{(l)} \rightarrow C_6H_{12}O_{6(aq)} + 6O_{2(g)} \quad \Delta_rH° = +2803 \text{ kJ mol}^{-1}$$
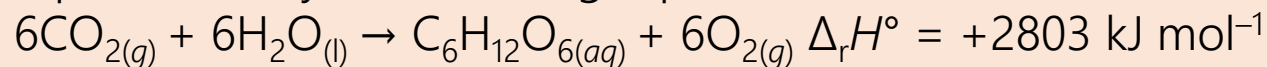
Answer 2b(i): Endothermic because the Δ_rH^o value is positive / it uses the sun's energy

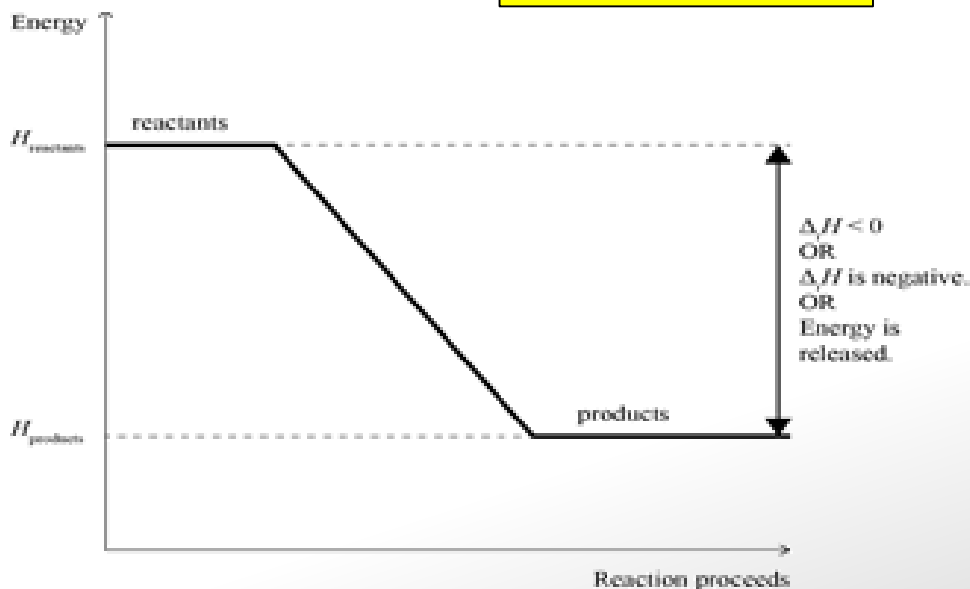Always attempt these questions as there is only 2 possible answers
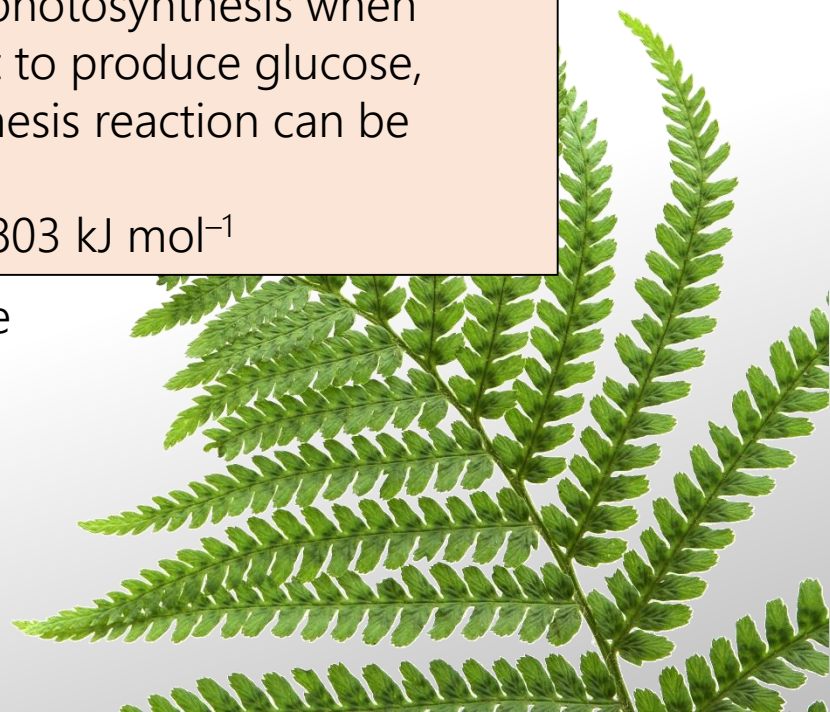
Excellence Question

Question 2c (iii): Complete, including labels, the energy diagram for the combustion of butane gas showing reactants, products, and the change in enthalpy.

Question 2c (iv): Butane gas is a useful fuel because when it undergoes combustion, energy is released. Explain why energy is released in this reaction, in terms of making and breaking bonds.

Answer 2c (iii):

Both required to be correct for Excellence

Energy

$H_{reactants}$ reactants

$\Delta H < 0$
OR
ΔH is negative.
OR
Energy is released.

$H_{products}$ products

Reaction proceeds

Answer 2c (iv): When butane undergoes combustion, heat is released, so it is an exothermic reaction.

Bond-making is an exothermic process / releases energy and bond-breaking is endothermic / requires energy. For the overall reaction in the combustion of butane to release energy, more energy is given out as bonds are made (when the products, CO_2 and H_2O are formed) than the energy being used to break the bonds (in the reactants, C_4H_{10} and O_2).

Question 1a Instant cold packs are useful for treating sports injuries on the field. They contain salts such as ammonium nitrate, NH_4NO_3. When the packs are activated, the salt dissolves in water, causing the temperature to decrease.
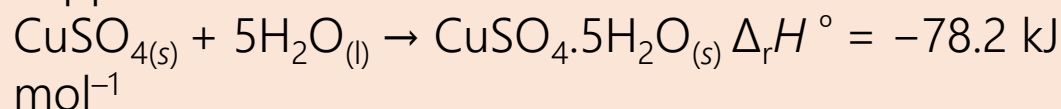
- Identify the term that best describes this reaction
- Give a reason for your choice

Heat absorbed is always endothermic

Answer 1a : Endothermic
The temperature decreased OR heat / energy has been absorbed.

Question 1b: The equation for hydrating anhydrous copper sulfate is as follows:
$CuSO_{4(s)} + 5H_2O_{(l)} \rightarrow CuSO_4.5H_2O_{(s)}$ $\Delta_r H\degree = -78.2$ kJ mol^{-1}

- Identify the term that best describes this reaction
- Give a reason for your choice

Answer 1b: Exothermic.
The enthalpy of the reaction is negative / energy has been released.

Question 1c (i): Pentane, C_5H_{12}, is a liquid at room temperature. It evaporates at 36.1°C in an endothermic process.
(i) Explain why the evaporation of pentane is an endothermic process.

Answer 1c(i) : Energy is required to change pentane from a liquid to a gas. The energy / heat is used to break weak intermolecular forces / bonds / attraction between pentane molecules.
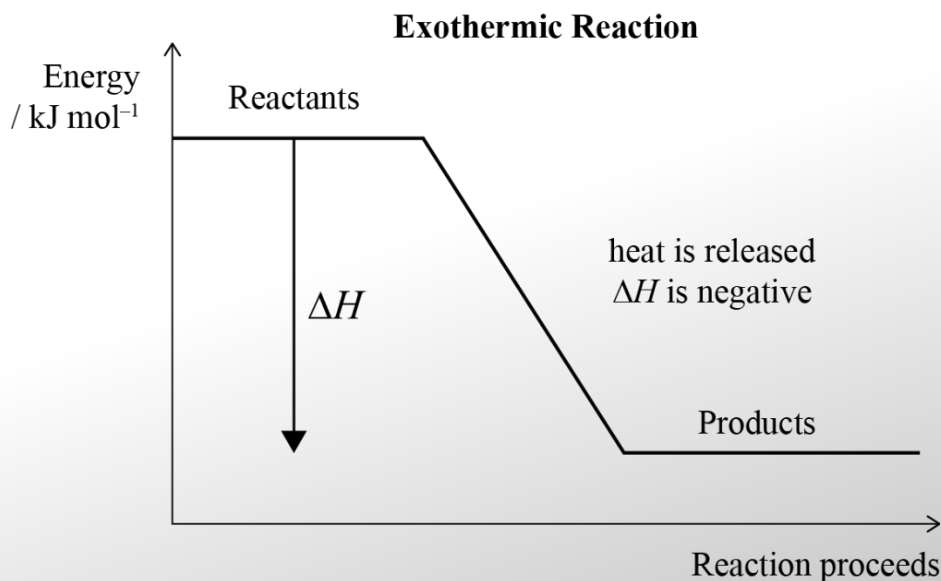
Solid →liquid →gas is always endothermic

Answer 1c(ii)

Question 1c(ii) : Draw, including labels, the energy diagram for the combustion of pentane, $C_5H_{12(l)}$.
Pentane combustion: $C_5H_{12(l)} + 8O_{2(g)} \rightarrow 5CO_{2(g)} + 6H_2O_{(l)}$
$\Delta_rH° = -3509$ kJ mol^{-1}
Include in your diagram the reactants, products, and change in enthalpy.

Exothermic Reaction

Energy / kJ mol^{-1}

Reactants

ΔH

heat is released
ΔH is negative

Products

Reaction proceeds

Question 1a: When solid calcium chloride, $CaCl_{2(s)}$, reacts with water, the temperature increases.
Which term that best describes this reaction.

Temperature increase is always exothermic

Exothermic
The temperature increased / energy or heat has been released into the surroundings /
energy is lost from the substance $(CaCl_2)$

Question 1b (i): When a person sweats, water is lost from the body by evaporation. This is an endothermic process. This evaporation speeds up when a person exercises.
(i) Explain why the evaporation of water in sweat from the body is endothermic, and why exercise increases this evaporation.

The water in sweat is changing state from liquid to gas. It needs to absorb energy to break the forces / bonds between liquid water molecules. It absorbs this from the heat of the body. The temperature of the body increases when exercising, so more water can be evaporated.
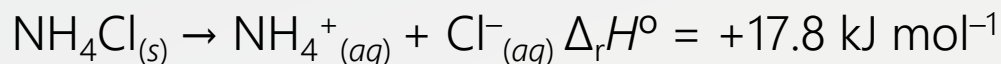
Excellence Question with b(i)

Question 1b (ii): Draw a labelled enthalpy diagram for the evaporation of water, $H_2O_{(\ell)}$.
$H_2O_{(\ell)} \rightarrow H_2O_{(g)}$ $\Delta_r H° = 40.7$ kJ mol^{-1}
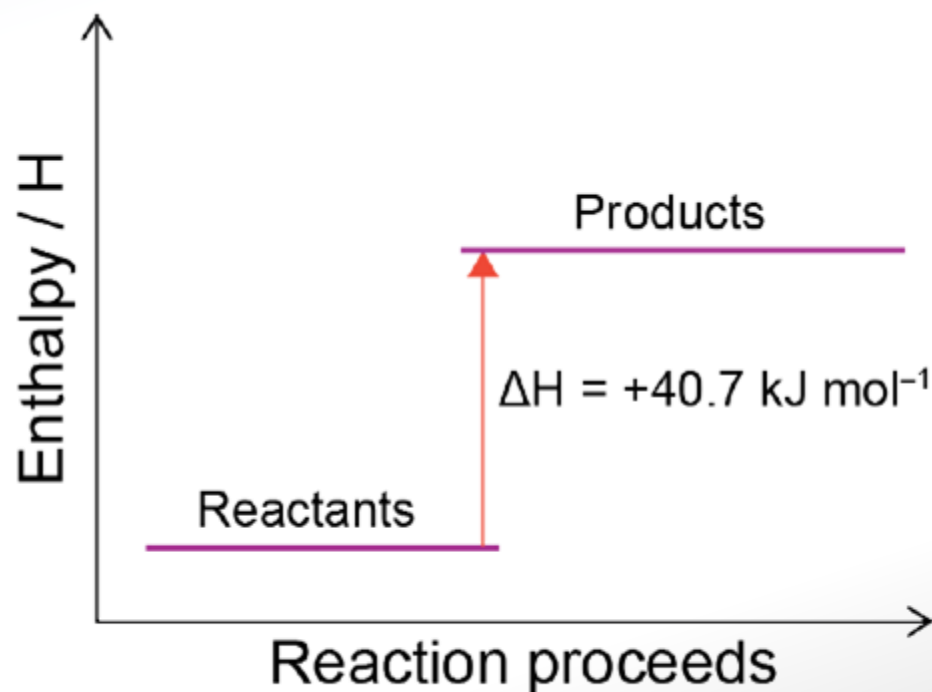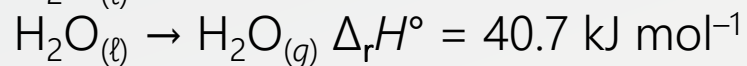
Can show activation energy but not required.

Question 1a. The equation for the dissolving of ammonium chloride, NH_4Cl, in water is shown below.

$NH_4Cl_{(s)} \rightarrow NH_4^+{}_{(aq)} + Cl^-{}_{(aq)}$ $\Delta_rH^o = +17.8$ kJ mol^{-1}

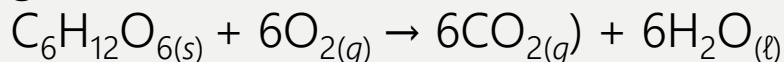Circle the term that best describes this reaction:

Endothermic   exothermic

Give a reason for your choice.

Endothermic because the enthalpy change is positive.

Question 1b (i) Respiration is the process by which energy is released from glucose.

$C_6H_{12}O_{6(s)} + 6O_{2(g)} \rightarrow 6CO_{2(g)} + 6H_2O_{(\ell)}$
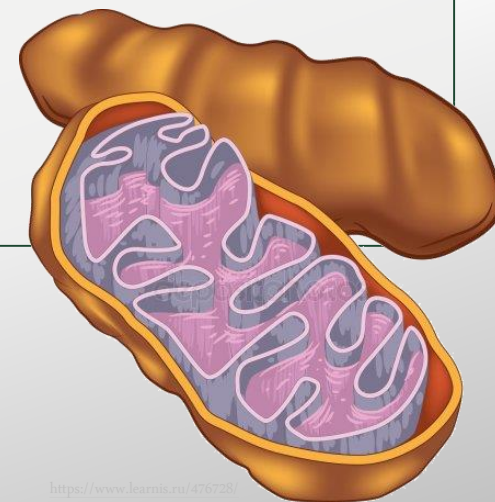
Circle the term that best describes this reaction:

endothermic   exothermic

Give a reason for your choice.

Exothermic because energy is released.

Question 1b (ii) . Water formed in the respiration reaction evaporates.

$H_2O_{(\ell)} \rightarrow H_2O_{(g)}$

Explain whether this process is endothermic or exothermic

The evaporation of water is endothermic because energy is absorbed to break the attractive forces between water molecules.

Merit Question

Question 1c. (i) Butane is used to fuel a camping stove. Butane burns readily in oxygen.
The following is an energy profile diagram for the combustion of butane.
Explain how the diagram shows that the enthalpy change for this reaction is negative.
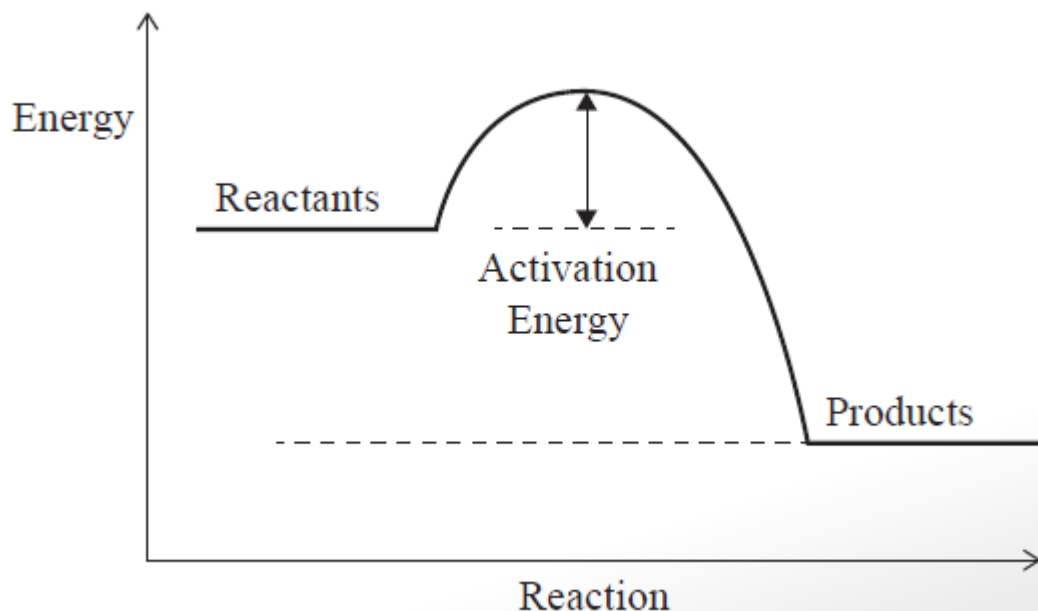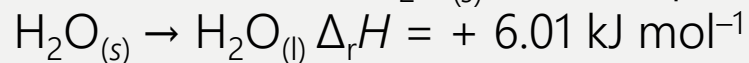

This is an exothermic reaction; the total energy of the products is less than the total energy of the reactants. Therefore, as the change in enthalpy is the difference in energy between products and reactants, the change is negative and the difference in energy is released as heat to the surroundings

Question 1d. Ice, $H_2O_{(s)}$, is often placed into drinks. As the ice melts, the drink cools.
$H_2O_{(s)} \rightarrow H_2O_{(l)}\ \Delta_r H = +\ 6.01\ kJ\ mol^{-1}$
Use your knowledge of enthalpy changes associated with changes of state to elaborate on the reason why the drink cools.

The melting of ice is endothermic, as (intermolecular) bonds are being broken as water changes from solid to liquid. This requires energy to be absorbed from the surroundings/the drink. This causes the temperature of the drink to decrease.

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The number of particles in a substance is measured in moles (unit n)

1 mole of particles = 6.02214 x 10^{23} particles for any substance!

6.02x10²³

AVOGADRO'S NUMBER

*

The mass of 1 mole of each element in grams, is equivalent to the relative Atomic Mass (Ar), which is an average of the isotopes mass in the proportions of isotopes found on Earth, and approximates to the number of protons and neutrons in the atom.

For example 1 mole of Carbon (Ar = 12.01) is 12.01g

If the actual number of particles is known (atoms, ions or molecules) then the amount, n, is easily calculated since

$$n = \frac{\text{number of particles}}{6.02 \times 10^{23}}$$

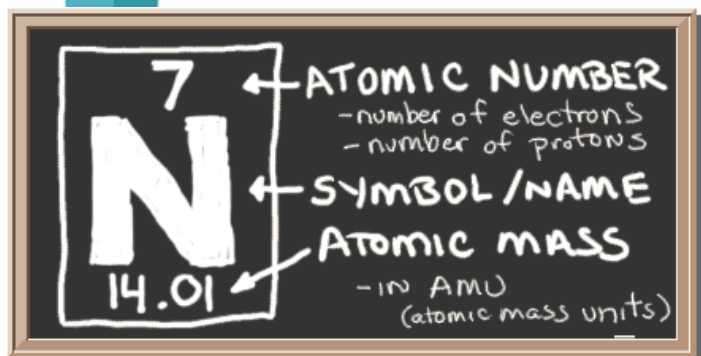For example 12×10^{24} sodium ions is equivalent to

$$\frac{12 \times 10^{24}}{6.02 \times 10^{23}}$$

= 20 mol of sodium ions.

6.02×10²³

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The *molar mass* is given as the *mass* of a given element or compound in grams (g), and divided by the amount, in moles, of the substance (mol). The unit for Molar Mass is, therefore, gmol^{-1}

Molar mass is calculated by adding up the **atomic mass** (or mass number) or each individual atom

The atomic mass used at L2 Chemistry is provided on a standardised periodic Table, also provided in all Examinations. This will be the correct value to use for calculations.

Calculating Molar Mass of a compound

1. Identify how many of each atom	2. Write a sum using the molar mass for each atom	3. Calculate the total
e.g. H_2SO_4 (2 H) (1 S) (4 O)	$2 \times M(H) + M(S) + 4 \times M(O)$ $(2 \times 1) + 32 + (4 \times 16)$	$M(H_2SO_4)$ $= 98$ gmol^{-1}
e.g. CuSO4.5H2O (1 Cu) (1 S) (9 O) (10 H)	$M(Cu) + M(S) + 9 \times M(O) + 10 \times M(H)$ $63.5 + 32 + (9 \times 16) + (10 \times 1)$	$M(CuSO_4.5H_2O)$ $= 249.5$ gmol^{-1}

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$$n = m / \mathscr{m}$$

We use the formula above to calculate the number of moles present of a substance when given the **mass of the substance** and the **molar mass** (atomic mass) from the Periodic Table.

m - Mass is measured in grams (g)

μ - Molar Mass is measured in gmol^{-1}

n   - Moles are measures in mols

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EXAMPLE
Calculate the amount (in moles) of methane, CH_4, in 12.5 g of the gas?
$M(CH_4) = 16.0$ g mol^{-1}

$$n = m / \mathcal{M}$$

1) Write down the formula and rearrange if necessary

$n(CH_4) = m/ \mu$

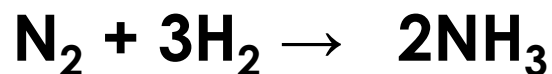2) Write in the values (with units) underneath and calculate

$n(CH_4) = 12.5g/16.0g$ mol^{-1}

$n(CH_4) = 0.781mol$

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A **balanced equation** will be given, which shows each substance with its lowest common number of moles in relationship to the other substances.

$$N_2 + 3H_2 \rightarrow 2NH_3$$

In the equation above **1 mole** of nitrogen gas (N_2) will react with **3 moles** of hydrogen gas (H_2) to form **2 moles** of ammonia (NH_3)

Of course the actual amount of moles you have in any reaction will depend on the mass you start with but the ratios of the other substances **will remain the same as in the equation**

For example: the number of moles of NH_3 produced will always be 2 x the number of moles of N_2 and 2/3 x the number of H_2

What is Stoichiometry? It is an area of chemistry that involves using relationships between reactants and/or products in a chemical reaction to determine quantitative data.

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In a question that requires you to calculate either mass or moles of substance – that substance is referred to as the **unknown** (U)

The substance that the mass is given for is referred to as the **known** (K)

When you need to calculate the mass or moles of an unknown you will need to multiply the moles of the known by a **mole ratio** from the equation.

The mole ratio is **x U / K** and the values of these are from the **balanced equation.**

$$N_2 + 3H_2 \rightarrow 2NH_3$$
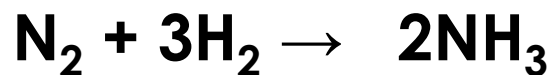
For example in the equation above if H_2 was the unknown (3 moles in equation) and NH_3 was the known (2 moles in equation) then:
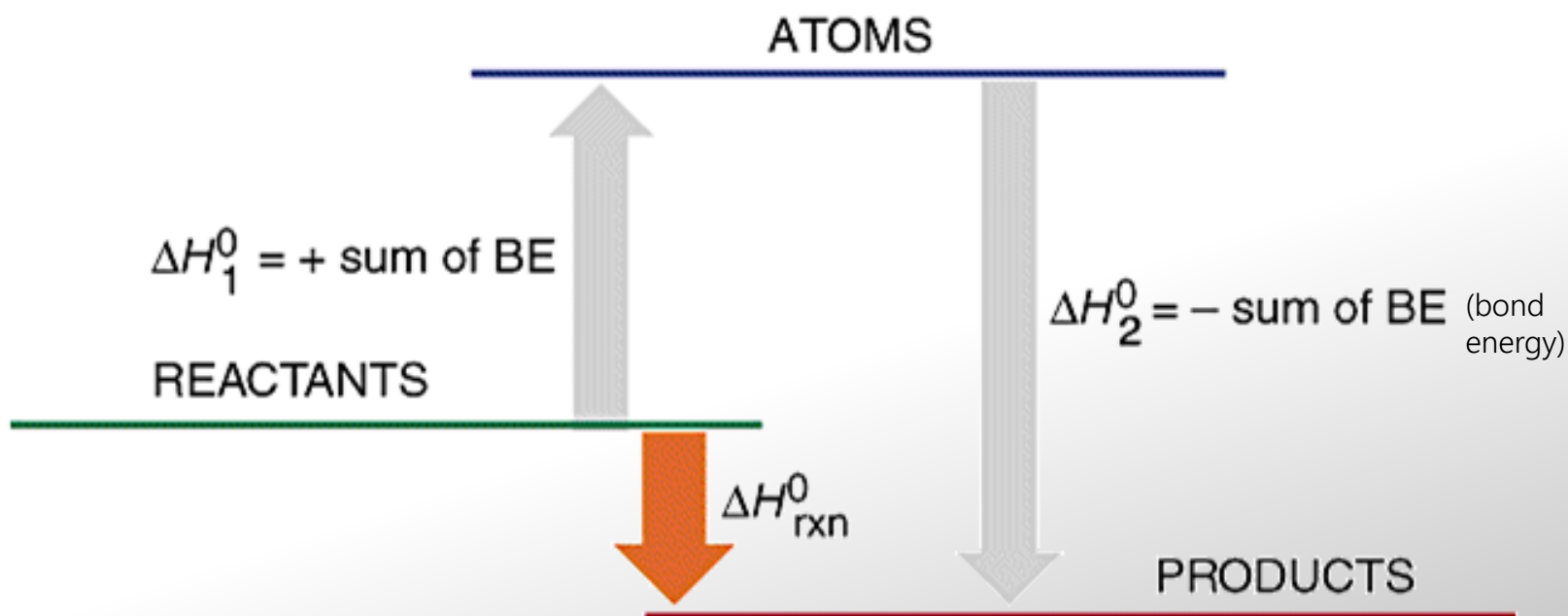
moles of hydrogen = moles of ammonia x 3/2

Exothermic - energy released   Δ H negative

Endothermic - energy absorbed   Δ H positive
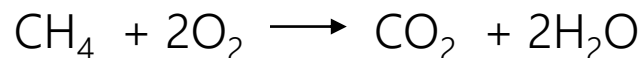
Standard Enthalpy of Reaction $\Delta_r H^{\ominus}$
"The enthalpy change when products are formed from their constituent reactants under standard conditions."

ATOMS

$\Delta H^0_1 = +$ sum of BE

$\Delta H^0_2 = -$ sum of BE (bond energy)

REACTANTS

ΔH^0_{rxn}

PRODUCTS

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You can perform stoichiometry calculations using energy changes from thermochemical equations. Using a balanced chemical equation to calculate amounts of reactants and products is called *stoichiometry*. Energy released or absorbed can be calculated per amount of substance.
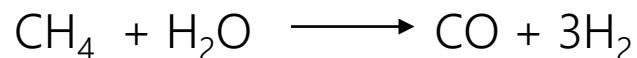
Exothermic	Endothermic
$CH_4 + 2O_2 \longrightarrow CO_2 + 2H_2O$	$CH_4 + H_2O \longrightarrow CO + 3H_2$
$\Delta H = -888 kJmol^{-1}$	$\Delta H = 206 \ kJmol^{-1}$
This thermochemical equation reads; 888kJ of heat is released when 1 mole of CH_4 reacts with 2 moles of O_2 to produce 1 mole of CO_2 and 2 moles of H_2O	This thermochemical reaction reads; 206kJ of heat is absorbed when 1 mole of CH_4 reacts with 1 mole of H_2O to produce 1 mole of CO and 3 moles of H_2.

Use thermochemical equations to find ΔH, n and m.   n = m/M
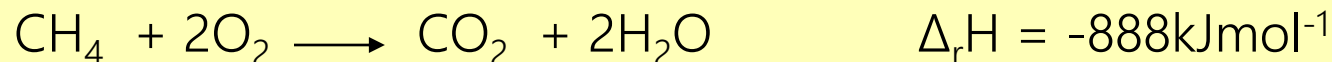
n = moles (6.02 x 10^{23} particles)   m = mass (grams)   M = Molar Mass ($gmol^{-1}$)

The **mole**, abbreviated **mol**, is an SI unit which measures the number of particles in a specific substance. **One mole** is equal to $6.02214179 \times 10^{23}$ atoms

$$CH_4 + 2O_2 \longrightarrow CO_2 + 2H_2O \qquad \Delta_rH = -888 \text{kJmol}^{-1}$$

Use the equation above to find heat released if 2.5 moles of CH_4 burns.

1 mole of CH_4 releases 888kJ
2.5 moles CH_4 releases **x** kJ
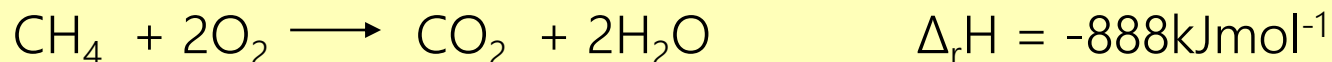x = 2.5 x 888 = **2220kJ**

An equation is a mole ratio – the number in front of each substance tells you how many moles of that there is to any other substance.
For example there is 1 mole of CH_4 to every 2 moles of O_2
The enthalpy of the equation shows you the amount of energy per unit of substance.
888 = 1CH_4 888 = 2O_2 (444 = 1O_2)

$$CH_4 + 2O_2 \longrightarrow CO_2 + 2H_2O \qquad \Delta_rH = -888 \text{kJmol}^{-1}$$

Calculate the **amount** (in moles) of H_2O produced when the reaction above releases **10,000kJ**.

Amount of mols in equation

2 moles H_2O when 888kJ released

x moles H_2O when 10000kJ released

$x = \dfrac{2 \times 10000}{888} = 22.5$ moles

Amount energy per unit of substance

Total energy released

An **alternative method** is to find out how much energy is released per mole first
2 moles H_2O = 888kJ
Therefore 1 mole H_2O = 444kJ
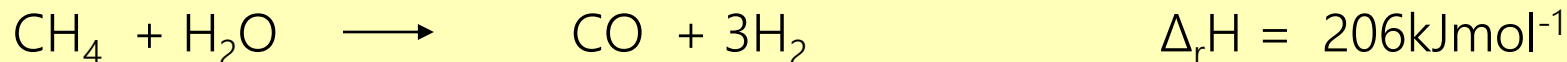
10,000kJ/444kJ = 22.5

So 22.5 moles of water are produced at 444kJ to reach 10,000kJ
(22.5 x 444 = 10,000)

$$CH_4 + H_2O \longrightarrow CO + 3H_2 \qquad \Delta_rH = 206kJmol^{-1}$$

$$M(C) = 12gmol^{-1} \qquad M(O) = 16gmol^{-1}$$

Calculate the energy required to produce 1kg of CO gas from the reaction above

Step one
moles of CO produced
M = 1000g $M(CO) = 28gmol^{-1}$
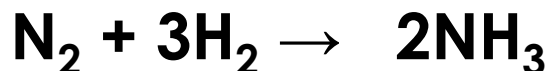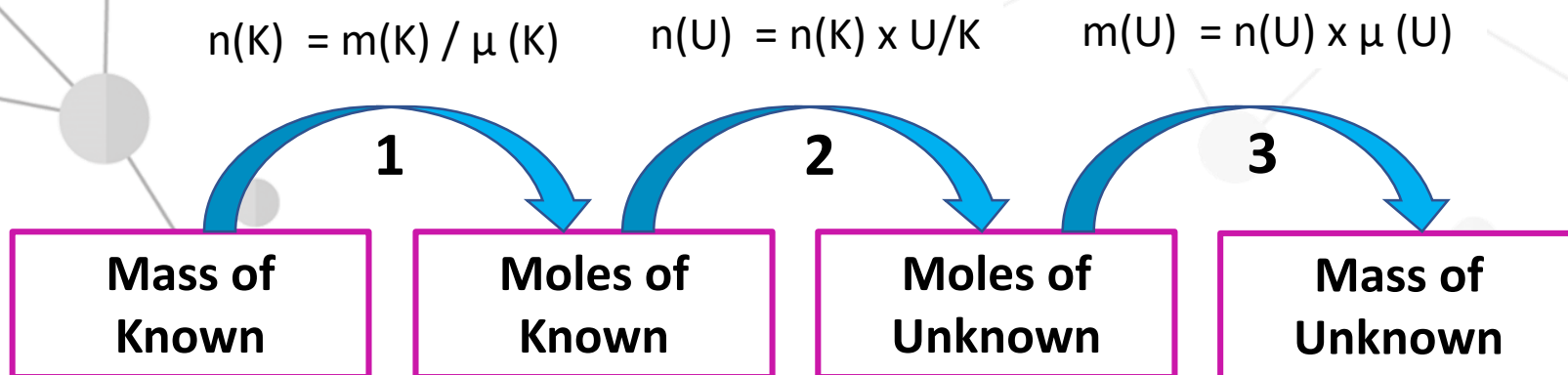
n = m/M
n = 1000/28 = 35.7 moles

1kg = 1000g. Must be converted to grams first

If Molar mass is not given then use the periodic table

The units are kJ **not** $kJmol^{-1}$ as it is total amount not amount per mole.

Step two
1 mole CO produced requires 206kJ (as per the equation above)
35.7 mols CO produced so.....
enthalpy = 35.7 x 206 = **7354kJ**

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The step that you begin and stop at depends on what information the question provides and what they want you to calculate. For example if the question provides the number of moles of a known and asks for the moles of the unknown only step 2 is required.

$n(K) = m(K) / \mu(K)$ $n(U) = n(K) \times U/K$ $m(U) = n(U) \times \mu(U)$

1   **2**   **3**

| **Mass of Known** | **Moles of Known** | **Moles of Unknown** | **Mass of Unknown** |

$N_2 + 3H_2 \rightarrow 2NH_3$

❑ Question asks for <u>moles of H_2</u> if moles of N_2 is given as 0.250mol . N_2 is Known (K) and H_2 is Unknown (U) so **only step 2 is required**. $n(H_2) = 0.025 \times U/K = 0.025 \times 3/1$

❑ Question asks for <u>mass of H_2</u> if moles of N_2 is given as 0.025mol. **Step 2 and 3 are required**. (note step one is not required as moles of known already given)

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	Known _____	Unknown _____
	x U / K	**2**
n (mols)		
m (g)	**1** ÷	x **3**
μ (gmol^{-1})		

1. n(K) = m(K) / μ (K)

2. n(U) = n(K) x U/K

3. m(U) = n(U) x μ (U)

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Step One:

> You need to establish which chemical /compound in an equation is the Known (K) – this will be the one that has information about it's mass. Write K above this in the equation.
> Establish which is the unknown (U). This will be the compound/chemical in the equation that the question is asking you to find the mass for. Write a U above this in the equation.
> Calculate the n(mols) of **Known**

$$n(K) = m(K) / \mu(K)$$

mass is from the question

molar mass added up from periodic table or given

Moles   mass   molar mass   add up each atom's molar mass in molecule

Step two:

Calculate mols of Unknown:

$$n(Unknown) = n(known) \times \frac{U}{K}$$

← Unknown mols from equation e.g **6** CO_2 mols = 6

← Known mols from equation

From Step One

Step three:

Calculate mass of Unknown (answer in g)

$$m(U) = n(U) \times \mu(U)$$

← molar mass calculated by adding up molar mass from each atom

From Step Two

$$c = n / v$$

We use the formula above to calculate the concentration of a unknown substance when given the **concentration of a known solution** when a titration has reached end point.

The **moles of each** substance is found in a balanced equation.

c - concentration is measured in $molL^{-1}$

v - volume is measured in L

n   - moles are measures in mols

$n(K) = c(K) \times v(K)$ $n(U) = n(K) \times U/K$ $c(U) = n(U) / v(U)$

1   **2**   **3**

| **Concentration and volume of Known** | **Moles of Known** | **Moles of Unknown** | **Concentration of Unknown** |

All steps must be followed with each, calculation (and equation) written down as **3 significant figures** consistently.

The entire calculation can done continuously on a calculator with the final answer of the **concentration of the unknown written down as 3sgf and units of molL^{-1}**

Highlight your final answer and make sure it appears to be a sensible answer. If you have time go through the calculation again quickly to check.

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Measurements depend on conditions
When measuring an enthalpy change you will get different values under different conditions. For example, the enthalpy change of a particular reaction will be different at different temperatures, different pressures or different concentrations of reactants. The values for enthalpy are given for standard conditions, indicated by the superscript θ

Standard conditions include:
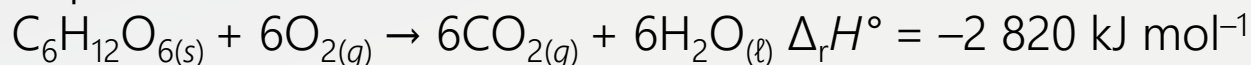Temperature of 25°C
Atmospheric pressure conditions of 1ATM
Concentration of 1mol per Litre

Question 3b(ii) : Females who are moderately active need 9 800 kJ of energy per day. Calculate the number of moles of glucose that would provide this daily energy requirement.
$C_6H_{12}O_{6(s)} + 6O_{2(g)} \rightarrow 6CO_{2(g)} + 6H_2O_{(\ell)}$ $\Delta_rH° = -2\ 820$ kJ mol^{-1}

Answer 3b(ii):
9800 kJ / 2820 kJ mol–1 = 3.48 mol

Question 3c(ii) : The equation below shows the combustion of butane.
$C_4H_{10(g)} + 13/2\ O_{2(g)} \rightarrow 4CO_{2(g)} + 5H_2O_{(g)}$ $M(C_4H_{10}) = 58.1$ g mol^{-1}.
When 100 g of butane undergoes combustion, 4 960 kJ of energy is released. Calculate the enthalpy change when 1 mole of butane undergoes combustion.
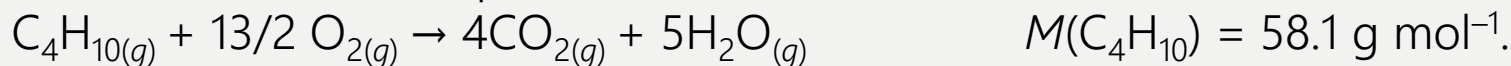
Answer 3c(ii):
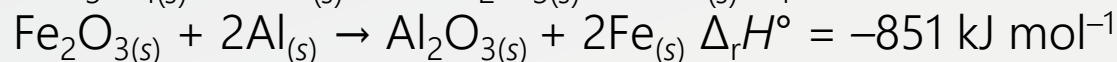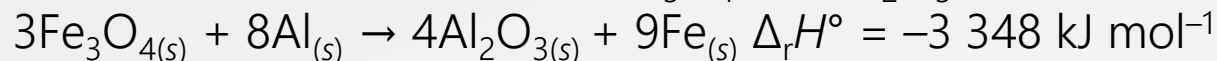$n(C_4H_{10})$ = 100 g / 58.1 g mol^{-1}
= 1.7212 mol
–4960 kJ / 1.7212 mol = –2882 kJ mol^{-1}

Calculating the amount of energy per mole

An equation and n=m/M are required for this type of thermochemical calculation

Question 3d: The iron oxides Fe_3O_4 and Fe_2O_3 react with aluminium as shown below.

$3Fe_3O_{4(s)} + 8Al_{(s)} \rightarrow 4Al_2O_{3(s)} + 9Fe_{(s)}$ $\Delta_rH° = -3\ 348$ kJ mol^{-1}

$Fe_2O_{3(s)} + 2Al_{(s)} \rightarrow Al_2O_{3(s)} + 2Fe_{(s)}$ $\Delta_rH° = -851$ kJ mol^{-1}

Justify which iron oxide, Fe_3O_4 or Fe_2O_3, will produce more heat energy when 2.00 kg of iron is formed during the reaction with aluminium.

Your answer should include calculations of the heat energy produced for the given mass of iron formed.

M(Fe) = 55.9 g mol^{-1}.

n(Fe) = 2000 g / 55.9 g mol–1 = 35.78 mol

Fe_3O_4:

3348 kJ / 9 = 372 kJ mol–1

372 kJ mol–1 × 35.78 mol = 13 310.16 kJ

= (–)1.33 ×10^4 kJ

Fe_2O_3:

851 kJ / 2 = 425.5 kJ mol–1

425.5 kJ mol–1 × 35.78 mol = 15 224.4 kJ

= (–)1.52 × 10^4 kJ

Complete each calculation one after the other

Don't forget to make a comparison statement

Therefore Fe_2O_3 produces more heat energy when 2 kg iron is formed.

Question 3c: Methanol and ethanol can both be used as fuels. Their combustion reactions can be represented by the following equations:

Methanol combustion: $2CH_3OH + 3O_2 \rightarrow 2CO_2 + 4H_2O$ $\Delta_r H^o = -1450$ kJ mol^{-1}

Ethanol combustion: $C_2H_5OH + 3O_2 \rightarrow 2CO_2 + 3H_2O$ $\Delta_r H^o = -1370$ kJ mol^{-1}

Justify which fuel, methanol or ethanol, will produce more heat energy when 345 g of each fuel is combusted in excess oxygen.

$M(CH_3OH) = 32.0$ g mol^{-1}

$M(C_2H_5OH) = 46.0$ g mol^{-1}

Answer 3c :

$n(CH_3OH) = m / M = 345 / 32 = 10.78$ mol

$n(C_2H_5OH) = m / M = 345 / 46 = 7.50$ mol

2 mol CH_3OH release 1 450 kJ of energy

1 mol CH_3OH releases 725 kJ of energy

10.78 mol CH_3OH releases 725 kJ x 10.78mol

= 7 816 kJ of energy

Complete each calculation one after the other

Don't forget to make a comparison statement

1 mol C_2H_5OH releases 1 370 kJ of energy

7.5 mol C_2H_5OH releases 1 370 kJ x 7.50mol = 10 275 kJ of energy

Therefore C_2H_5OH releases more energy when 345 g of fuel are combusted.

Excellence Question

Question 2b(ii) : Calculate how much energy is absorbed or released in the photosynthesis reaction if 19.8 g of carbon dioxide gas, $CO_{2(g)}$, reacts completely with excess water, $H_2O_{(l)}$, to form glucose, $C_6H_{12}O_{6(aq)}$, and oxygen gas, $O_{2(g)}$. Show your working and include appropriate units in your answer.
$M(CO_2) = 44.0$ g mol^{-1}
$6CO_{2(g)} + 6H_2O_{(l)} \rightarrow C_6H_{12}O_{6(aq)} + 6O_{2(g)}$ $\Delta_r H° = +2803$ kJ mol^{-1}
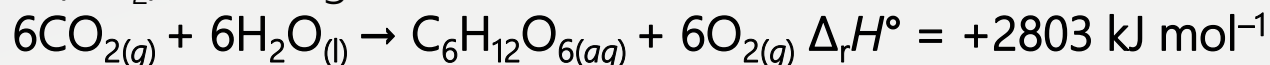
An equation and n=m/M are required for this type of thermochemical calculation

Answer 2b(ii):

$n(CO_2) = m/M$ $= 19.8/ 44.0$ $= 0.450$ mol

Since 6 moles of CO_2 reacting requires 2803 kJ of energy
then 1 mole of CO_2 reacting requires $2803/6 = 467.2$ kJ of energy
and 0.450 moles of CO_2 requires $467.2 \times 0.450 = 210$ kJ of energy absorbed.

Convert mass into mols

Units are in kJ and the + sign indicates energy is absorbed

Enthalpy change ($\Delta_r H$) for a reaction, can be collected from a thermochemical investigation. Values required are masses of water in which the reaction takes place, temperature change (in °C) and the specific heat capacity value for water. (These calculations will be done in Level 3)

Often questions will ask why a particular investigation set up does not provide the same thermochemical data as the accepted enthalpy change.

Reasons can include errors such as:

1. Some energy is used to heat the metal can and the air surrounding the experiment / the experiment was not conducted in a closed system
2. Incomplete combustion of butane.
3. Some butane may have escaped before being ignited.
4. The butane in the gas canister was impure.
5. Some water evaporated
6. Some energy was converted to light and sound
7. Not carried out under standard conditions

Therefore, not all of the energy released by the combustion of butane was transferred to heating the water

Thermometer

Metal can

Water

Tap

GAS — Small camping gas 'bottle'

Excellence Question
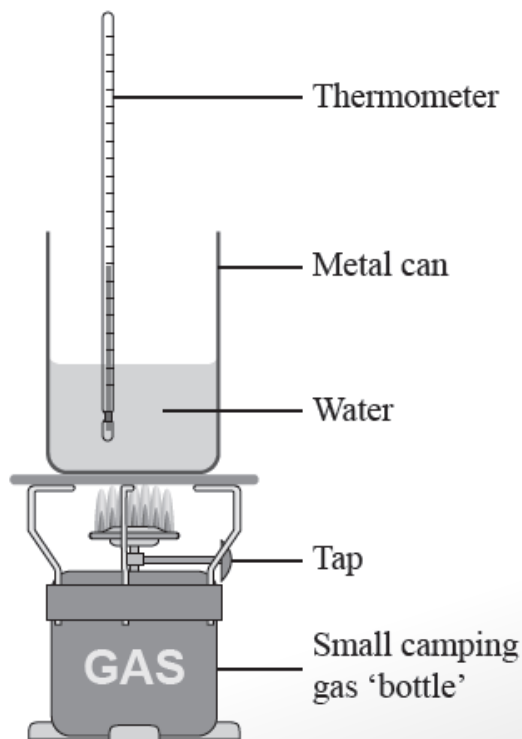
Question 2c: A small camp stove containing butane gas, $C_4H_{10(g)}$, is used to heat some water, as shown in the diagram below. A student measures the temperature change in the water and calculates that when 3.65 g of butane is combusted, 106 kJ of heat is released.

The reaction for the combustion of butane is shown in the equation below.

$2C_4H_{10(g)} + 13O_{2(g)} \rightarrow 8CO_{2(g)} + 10H_2O_{(l)}$
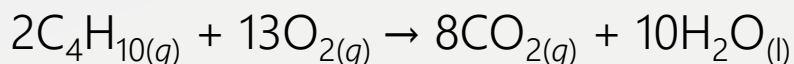
(i) Calculate the enthalpy change ($\Delta r\,H$) for this reaction, based on the above measurements. $M(C_4H_{10}) = 58.0$ g mol^{-1}

Values are all given to 3 sgf

Thermometer

Metal can

Answer 2c:

$n(C_4H_{10}) = m/M \quad = 3.65/\,58.0 \quad = 0.0629$mol

Water

If 0.0629 moles of C_4H_{10} releases 106 kJ of energy
Then 1 mole of C_4H_{10} releases $106/0.0629 = 1685$ kJ of energy
And 2 moles of C_4H_{10} releases $1685 \times 2 = 3370$ kJ of energy (3368) ($\Delta_r H = -3370$ kJ mol^{-1})

Must show sign

Tap

GAS
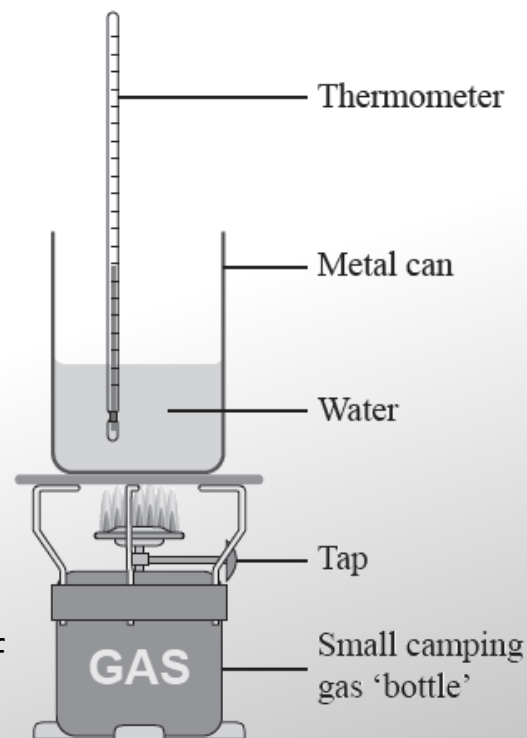
Small camping gas 'bottle'
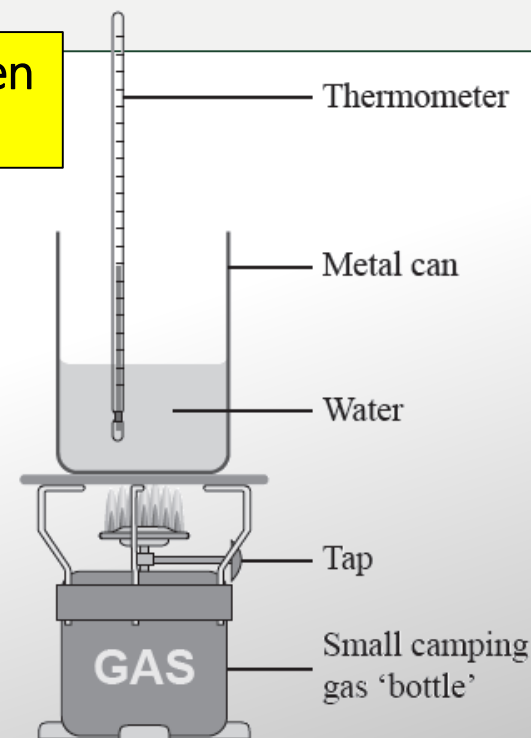
Question 2c: (ii) The accepted enthalpy change for the combustion reaction of butane gas, $C_4H_{10(g)}$, is $\Delta r\ H = -5754$ kJ mol^{-1}.
Explain why the result you calculated in part (c)(i) is different to the accepted value. In your answer, you should include at least TWO reasons.

TWO errors required for Excellence

Answer 2c: The results from this experiment are less than the accepted results, due to errors in the experimental design. The errors could include:

1. Some energy is used to heat the metal can and the air surrounding the experiment / the experiment was not conducted in a closed system
2. Incomplete combustion of butane.
3. Some butane may have escaped before being ignited.
4. The butane in the gas canister was impure.
5. Some water evaporated
6. Some energy was converted to light and sound
7. Not carried out under standard conditions

Therefore, not all of the energy released by the combustion of butane was transferred to heating the water.

- Thermometer
- Metal can
- Water
- Tap
- Small camping gas 'bottle'
- GAS

Question 1c(iii): Hexane, C_6H_{14}, like pentane, will combust (burn) in sufficient oxygen to produce carbon dioxide gas and water. Pentane combustion: $\Delta_rH\,° = -3509$ kJ mol^{-1}
Hexane combustion: $2C_6H_{14(l)} + 19O_{2(g)} \rightarrow 12CO_{2(g)} + 14H_2O_{(l)}$ $\Delta_rH\,° = -8316$ kJ mol^{-1}
Justify which alkane – pentane or hexane – will produce more heat energy when 125 g of each fuel is combusted in sufficient oxygen.
$M(C_5H_{12}) = 72.0$ g mol^{-1} $M(C_6H_{14}) = 86.0$ g mol^{-1}

An equation and n=m/M are required for this type of thermochemical calculation

Answer 1c(iii):

n (pentane) = 125 g / 72.0 g mol^{-1} = 1.74 mol
n (hexane) = 125 g / 86.0 g mol^{-1} = 1.45 mol
If 1 mole of pentane releases 3509 kJ energy,
then 1.74 mol of pentane:
1.74 × 3509 = 6106 kJ energy released.

Convert mass into mols

If 2 moles of hexane release 8316 kJ energy,
then 1 mole of hexane releases 4158 kJ energy.
So 1.45 mol of hexane 1.45 × 4158 = 6029 kJ energy releases.

So pentane releases more energy (77.0 kJ) than hexane,
per 125 g of fuel.

Make sure both fuels are compared for Excellence

Excellence Question

Question 1c: Thermite reactions occur when a metal oxide reacts with a metal powder. The equations for two thermite reactions are given below:

Reaction 1: $Fe_2O_{3(s)} + 2Al_{(s)} \rightarrow 2Fe_{(s)} + Al_2O_{3(s)}$ $\Delta_rH° = -852$ kJ mol^{-1}

Reaction 2: $3CuO_{(s)} + 2Al_{(s)} \rightarrow 3Cu_{(s)} + Al_2O_{3(s)})$ $\Delta_rH° = -1520$ kJ mol^{-1}
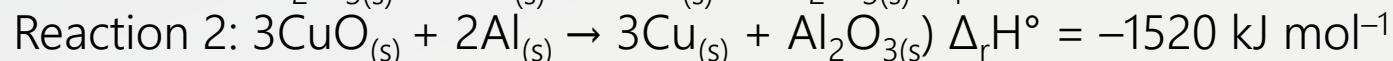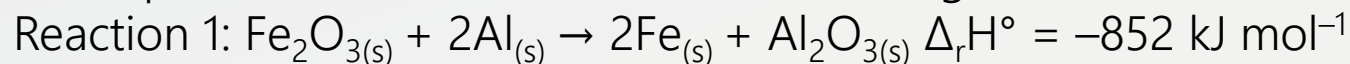
Use calculations to determine which metal oxide, iron(III) oxide, $Fe_2O_{3(s)}$, or copper(II) oxide, $CuO_{(s)}$, will produce more heat energy when 50.0 g of each metal oxide is reacted with aluminium powder, $Al_{(s)}$.

$M(Fe_2O_3) = 160$ g mol^{-1} $M(CuO) = 79.6$ g mol^{-1}

An equation and n=m/M are required for this type of thermochemical calculation
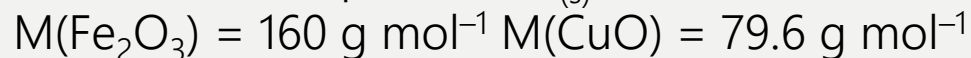
$n(Fe_2O_3) = \dfrac{50.0\ g}{160\ g\ mol^{-1}} = 0.313$mol

$n(CuO) = \dfrac{50.0\ g}{79.6\ g\ mol^{-1}} = 0.628$ mol

Reaction 1: If 1 mole of Fe_2O_3 releases 852 kJ energy
0.313 mol × 852 kJ mol^{-1} = 266 kJ energy released

Reaction 2: If 3 mole of CuO releases 1520 kJ energy
Then 1 mole of CuO releases 507 kJ energy
0.628 mol × 507 kJ mol–1 = 318 kJ energy released

So 50.0 g CuO releases more energy than 50.0 g Fe_2O_3
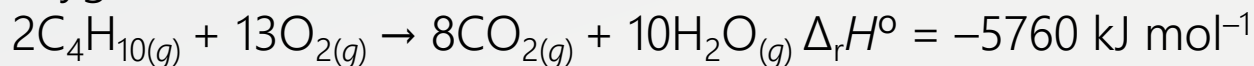OR
CuO releases more energy (52 kJ) than Fe_2O_3
(**Reaction 2** releases more energy.)

Convert mass into mols

Make sure both substances are compared for Excellence

Question 1c. (ii) The following is the equation for the combustion of butane gas in oxygen.

$2C_4H_{10(g)} + 13O_{2(g)} \rightarrow 8CO_{2(g)} + 10H_2O_{(g)}$ $\Delta_rH^o = -5760$ kJ mol^{-1}

The fuel cylinder for the stove contains 450 g of butane gas.

Calculate the energy released when this mass of butane gas is burned completely in oxygen.

Show your working and include appropriate units in your answer.

$M(C_4H_{10})$ = 58.0 g mol^{-1}
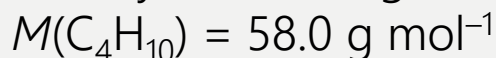
If 2 moles of butane produce 5760 kJ then 1 mole produces 2880 kJ.

n(butane) = 450 / 58.0

= 7.76 moles

q = 7.76 × 2880 = <u>**22 300 kJ released**</u>

alternative:

n(butane) = 450 / 58.0

= 7.76 moles

so n(reaction) = n(butane)/2 = 3.88 mol

So q = 5760 x 3.88 = <u>**2.23 x 10^4 kJ released**</u>

Question 2d. Methanol, $CH_3OH_{(\ell)}$, is made industrially by reacting carbon monoxide, $CO_{(g)}$, and hydrogen, $H_{2(g)}$.

$CO_{(g)} + 2H_{2(g)} \rightarrow CH_3OH_{(\ell)}$ Δ_rH° = –91.0 kJ mol^{-1}
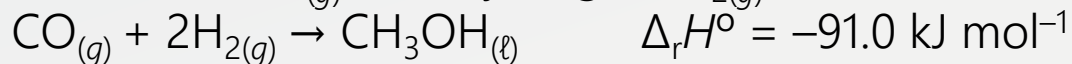
Calculate the volume of methanol made when 4428 kJ of energy is released.
The mass of 1.00 L of methanol is 0.790 kg.
$M(CH_3OH)$ = 32.0 g mol^{-1}

n(methanol) = 4428 / 91
= 48.7 moles

mass (methanol) = 48.66 × 32 = 1557 g

volume (methanol) = 1.56 / 0.790
= <u>1.97 L</u>

Question 3b. When magnesium, $Mg_{(s)}$, is burned, it produces a white powder according to the equation below.

$2Mg_{(s)} + O_{2(g)} \rightarrow 2MgO_{(s)}$ $\Delta_r H = -1203$ kJ mol^{-1}
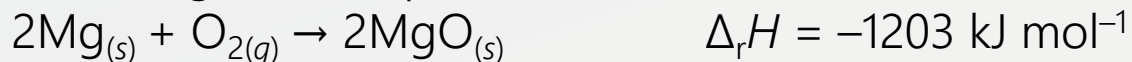
(i) Calculate the mass of oxygen required to produce 1804.5 kJ of energy.

$M(O) = 16.0$ g mol^{-1}

$$n(O_2(g)) = \frac{1804.5}{1203} = 1.5 \text{ moles}$$

$$m = n \times M = 1.5 \times 32 = 48.0 \text{ g}$$

(ii) Calculate the energy change when 100 g of $MgO_{(s)}$ is produced.

$M(MgO) = 40.3$ g mol^{-1}

$$n = \frac{m}{M} = \frac{100}{40.3} = 2.48$$

$$\text{Energy} = \Delta_r H \times n = \frac{-1203}{2} \times 2.48 = -1492.5 \text{ kJ} / -1493 \text{ kJ}$$

-1490 kJ (3 sf)

(either positive or negative values accepted)

Question 3c. A common industrial process is the extraction of metals from their ores. Aluminium is found naturally in aluminium oxide, and the oxygen is removed to produce the metal. Information is given below of the enthalpy change when aluminium, $Al_{(s)}$, is extracted.

$$2Al_2O_{3(s)} \rightarrow 4Al_{(s)} + 3O_{2(g)} \qquad \Delta_rH = 3350 \text{ kJ mol}^{-1}$$
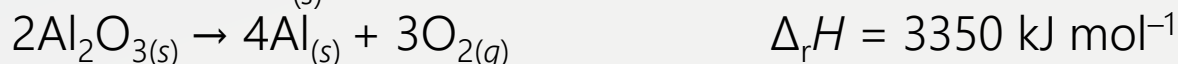
A production plant produces 65.0 kg (65 000 g) of aluminium per minute. Calculate how much energy is required per hour of production of aluminium. Round your answer to 3 significant figures. $M(Al) = 27.0 \text{ g mol}^{-1}$

$$65\,000 \times 60 = 3\,900\,000 \text{ g}$$

$$n(Al) = \frac{3\,900\,000}{27} = 144\,444 \text{ moles}$$

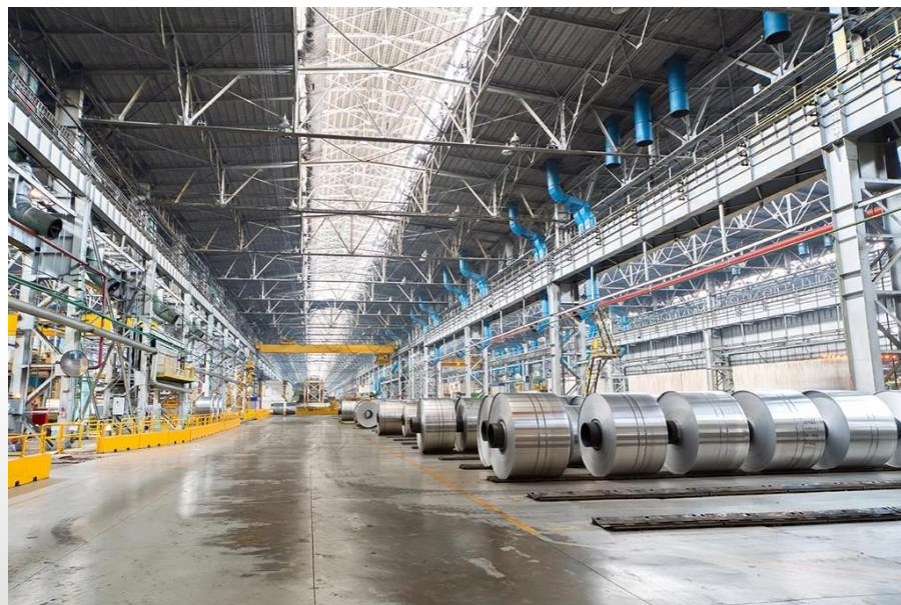$$\Delta_rH = \frac{144\,444}{4} \times 3350$$

$$= 121\,000\,000 \text{ kJ} = 1.21 \times 10^8 \text{ kJ (rounded to 3sf)}$$

OR $\dfrac{65000}{27} = 2407$ moles

$$\Delta_rH = \frac{2407}{4} \times 3350 = 2\,015\,862 \text{ kJ} \times 60$$

$$= 121\,000\,000 \text{ kJ} = 1.21 \times 10^8 \text{ kJ (rounded to 3sf)}$$

Bond enthalpy (also known as bond energy) is defined as the amount of energy required to break one mole of the stated bond.

The high values for bond enthalpy explains why some substances are very resistant to chemical attack and form very stable molecules

In a polyatomic (more than one atom) molecule, the bond strength between a given pair of atoms can vary slightly from one compound to another. The value given for bond enthalpy is the average of all these variations.

A multiple bond (double/triple) is always stronger than a single bond because more electrons bind the multiple bonded atoms together.
The table below shows some common average bond enthalpies.

Bond enthalpy /kJ mol^{-1}		Bond enthalpy /kJ mol^{-1}		thalpy /kJ mol^{-1}	
H - H	436	C - H	412	C = C	612
H - O	463	C - Cl	338	C o C	837
H - N	388	C - F	484	C = O	743
H - Cl	431	C - O	360	O = O	496
H - F	565	C - C	348	N ≡ N	944
F - F	158	O – O	146		
Cl - Cl	242				

Bond Enthalpy calculations

Bonds Broken – Endothermic   **Bonds formed – Exothermic**

$\Delta_r H° = \sum$ (energy of bonds broken) $- \sum$(energy of bonds formed)

Note: Bond energies calculated for gases. Convert using $\Delta_{vap}H°$ or $\Delta_{sub}H°$ if in solid or liquid state.

$$CO_{(g)} + H_2O_{(g)} \rightarrow H_{2(g)} + CO_{2(g)} \quad \Delta_r H° = -1.0 \text{ Kjmol}^{-1}$$

Bonds Broken		Bonds formed	
C≡O	995kJ	C=O x 2	2(743)kJ
H-O x 2	2(463)kJ	H-H	436kJ
	1921kJ		1922kJ

$\Delta_r H° = 1921\text{kJmol}^{-1} - 1922\text{kJmol}^{-1}$ The equation can also be arranged

$\Delta_r H° = -1.0 \text{ kJmol}^{-1}$ to calculate unknown bond energy

Bonds Broken – Endothermic   Bonds formed – Exothermic

$\Delta_rH° = \sum$ (energy of bonds broken) - $\sum$(energy of bonds formed)

Reactants:
Draw lewis diagrams to calculate the number and type of bond

Multiply the bond energy given by the number of bonds

Products:
Draw lewis diagrams to calculate the number and type of bond

$CO_{(g)} + H_2O_{(g)} \rightarrow H_{2(g)} + CO_{2(g)}$ $\Delta_rH° = -1.0 \ Kjmol^{-1}$

Bonds Broken		Bonds formed	
C≡O	995kJ	C=O x 2	2(743)kJ
H-O x 2	2(463)kJ	H-H	436kJ
	1921kJ		1922kJ

$\Delta_rH° = 1921kJmol^{-1} - 1922kJmol^{-1}$ The equation can also be arranged

$\Delta_rH° = -1.0 \ kJmol^{-1}$ to calculate unknown bond energy

Total the bond energy for reactant molecules

Total the bond energy for product molecules

bonds broken (reactants) minus bonds formed (product) = total enthalpy

Bond enthalpy is the change in enthalpy when the covalent bond, **in a gaseous molecule,** is broken. It is always a positive value because bond breaking always requires an input of energy.
Making bonds releases energy so generally speaking the more bonds a substance can form the more stable it will be.

The strength of a covalent bond depends on the **electrostatic attraction** between the positive nuclei and the shared electron pair. The larger the atomic radius of an atom (which increases down a group) the further the shared electron pair from the positive nucleus – which creates decreasing electrostatic attraction. Therefore the weaker the covalent bond and the lower the value of bond enthalpy.

The stronger a covalent bond, the higher the value of the bond enthalpy. The units are kJ mol^{-1}

Draw Lewis structures if not given

Excellence Question

Question 2c: Chlorine reacts with methane to form chloromethane and hydrogen chloride, as shown in the equation below.
$CH_{4(g)} + Cl_{2(g)} \rightarrow CH_3Cl_{(g)} + HCl_{(g)}$
Use the following bond enthalpies to calculate Δr $H°$ for this reaction.
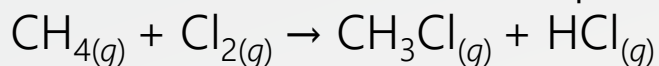
$\Delta_r H^o$ = ΣBond energies(bonds broken) – ΣBond energies(bonds formed)

Bond energies always given – use units on chart to remind you

Bonds broken | Bonds formed
C-H x 4   1656   C–Cl   324
Cl–Cl   242 C–H   3 × 1242
            1898 H-Cl   431
                          1997
= 1898 – 1997   = –99 kJ mol⁻¹

Alternative calculation that includes the breaking and reforming of only the bonds involved in the reaction
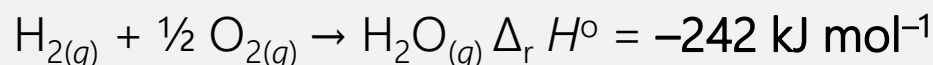(656 – 755 = 99kJ mol⁻¹)

Bond	Bond enthalpy /kJ mol⁻¹
H–Cl	431
C–H	414
C–Cl	324
Cl–Cl	242

Excellence Question

Question 1d: Hydrogen gas, $H_{2(g)}$, reacts with oxygen gas, $O_{2(g)}$, as shown by the following equation

$$H_{2(g)} + \tfrac{1}{2}\,O_{2(g)} \rightarrow H_2O_{(g)} \quad \Delta_r H^o = -242 \text{ kJ mol}^{-1}$$

Given the average bond enthalpies in the table below, calculate the average bond enthalpy of the $O - H$ bond in H_2O.

$\Delta_r H^o = \Sigma$Bond energies(bonds broken) $- \Sigma$Bond energies(bonds formed)

Bond energies always given – use units on chart to remind you

Bonds broken	Bonds formed
H–H = 436	$2 \times$ O–H
$\tfrac{1}{2} \times$ O=O = $\tfrac{1}{2} \times$ 498	
Total = 685 kJ mol^{-1}	Total = ?

-242 kJ mol^{-1} = 685kJ - ?

? = 685 − (− 242)

= 927 kJ mol^{-1}

Swap total and ?

Bond	Average bond enthalpy / kJ mol^{-1}
H–H	436
O=O	498

$2 \times$ O–H = 927 kJ mol^{-1}

but ONE   **O–H** = 464 (463.5) kJ mol^{-1}

Excellence Question

Question 1d: Ethene gas, $C_2H_{4\,(g)}$, reacts with bromine gas, $Br_{2(g)}$, as shown in the equation below.

Calculate the enthalpy change, $\Delta_r H°$, for the reaction between ethane and bromine gases, given the average bond enthalpies in the table below. Show your working and include appropriate units in your answers.

Draw Lewis structures if not given

$$\begin{array}{ccc} H & & H \\ & C=C & \\ H & & H \end{array} (g) \;+\; Br-Br\,(g) \;\rightarrow\; \begin{array}{cc} H & H \\ H-C-C-H & (g) \\ Br & Br \end{array}$$
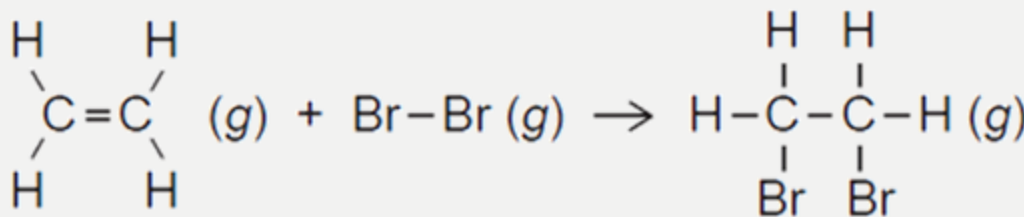
Bond energies always given – use units on chart to remind you

$\Delta_r H° = \Sigma$Bond energies(bonds broken) $- \Sigma$Bond energies(bonds formed)

Bonds broken	Bonds formed
C=C 614	C–C 346
Br–Br 193	C–Br 2×285
807	916

$= 807 - 916 \quad = -109$ kJ mol^{-1}

Alternative calculation that includes the breaking and reforming of four C-H bonds
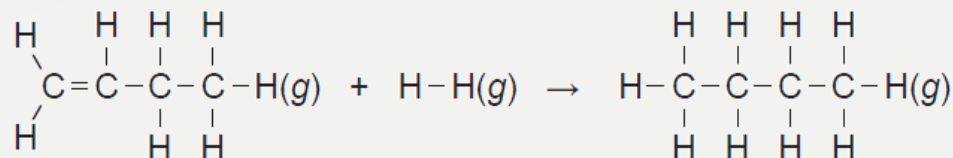(or $2463 - 2572$)

Bond	Average bond enthalpy/kJ mol^{-1}
Br–Br	193
C–C	346
C=C	614
C–Br	285
C–H	414

Draw Lewis structures if not given

Excellence Question

Question 3c: Calculate the enthalpy change, $\Delta_rH°$, for the reaction of but-1-ene gas, $C_4H_{8(g)}$, with hydrogen gas, $H_{2(g)}$, to form butane gas, $C_4H_{10(g)}$.
Use the average bond enthalpies given in the table below.

Bond energies always given – use units on chart to remind you

Bond breaking

C=C 614
C–C × 2   692
C–H × 8   3312
H–H 436
    5054 kJ mol^{-1}

Bond making

C–C × 3 1038
C–H × 10   4140
    5178 kJ mol^{-1}

$\Delta_rH°$ = Bond breaking – bond making
$\Delta_rH°$ = 5054 kJ mol^{-1} – 5178 kJ mol^{-1}
$\Delta_rH°$ = −124 kJ mol^{-1}

 Alternative calculation that includes the breaking and reforming of only the bonds involved in the reaction
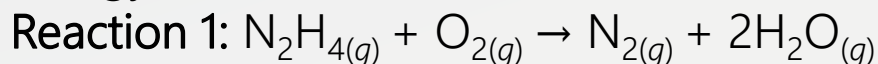$\Delta_rH°$ = 1050 – 1174 = −124 kJ mol^{-1}

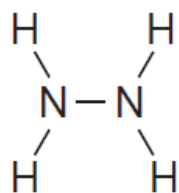| Bond	Average bond enthalpy / kJ mol^{-1}
C=C	614
C–C	346
C–H	414
H–H	436

Question 2c: Hydrazine, N_2H_4, is used as rocket fuel.

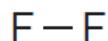Use calculations to determine which of **Reaction 1** or **Reaction 2** releases more energy.

Reaction 1: $N_2H_{4(g)} + O_{2(g)} \rightarrow N_{2(g)} + 2H_2O_{(g)}$

Reaction 2: $N_2H_{4(g)} + 2F_{2(g)} \rightarrow N_{2(g)} + 4HF_{(g)}$
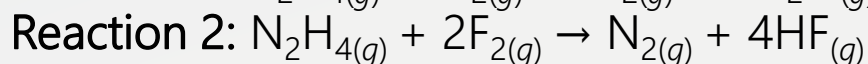
The structure of each chemical species is shown in the box below. Show your working and include appropriate units in your answer.
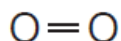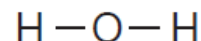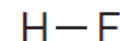

Use the average bond enthalpies given in the table below.

Bond	Average Bond enthalpy /kJ mol^{-1}	Bond	Average Bond enthalpy /kJ mol^{-1}
H–H	436	N–N	158
H–F	567	F–F	159
N–H	391	O=O	498
O–H	463	N≡N	945

Question 2c: Hydrazine, N_2H_4, is used as rocket fuel.
Use calculations to determine which of **Reaction 1** or **Reaction 2** releases more energy.

Reaction 1
Hydrazine and oxygen

Bond breaking		Bond making	
N–N	158	N≡N	945
N–H × 4	1564	O–H × 4	1852
O=O	498		2797
	2220		

Bond breaking – bond making
$2220 - 2797 = -577$ kJ mol^{-1}

Reaction 2
Hydrazine and Fluorine

N–N	158	N≡N	945
N–H × 4	1564	H–F × 4	2268
F–F × 2	318		3213
	2040		

Bond breaking – bond making
$2040 - 3213 = -1173$ kJ mol^{-1} (or -1170 kJ mol^{-1})
Reaction 2 releases more energy than **Reaction 1** (by 596 kJ mol^{-1}).

Question 1d. Nitrogen gas, $N_{2(g)}$, reacts with hydrogen gas, $H_{2(g)}$, to produce ammonia gas, $NH_{3(g)}$, as shown by the following equation:

$N_{2(g)} + 3H_{2(g)} \rightarrow 2NH_{3(g)}$ $\Delta_r H^° = -92.0$ kJ mol^{-1}

Calculate the average bond enthalpy of the **N–H** bond in 3, using the average bond enthalpies in the table below.

Bond	Average bond enthalpy kJ mol^{-1}
N≡N	945
H–H	436

$\Delta H = \Sigma E$(Bonds broken) $- \Sigma E$(Bonds made) $= -92.0$ kJ mol^{-1}

E(N≡N) $+ 3E$(H–H) $- 6E$(N–H) $= -92.0$ kJ mol^{-1}

$6E$(N–H) $= 2253 - (-92.0) = 2345$ kJ mol^{-1}

Therefore E(N–H) = 2345 / 6

= **391 kJ mol^{-1}**

Question 2c. When propane, $C_3H_{8(g)}$, is burned, it reacts with oxygen, $O_{2(g)}$, in the air to form water, $H_2O_{(g)}$, and carbon dioxide, $CO_{2(g)}$.

$C_3H_{8(g)} + 5O_{2(g)} \rightarrow 4H_2O_{(g)} + 3CO_{2(g)}$ $\Delta_rH = -2056$ kJ mol^{-1}

Calculate the average bond enthalpy of the C = O bond using the data below.
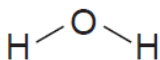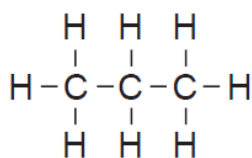
Bond	Average bond enthalpy/kJ mol^{-1}
C – C	348
C – H	413
O = O	495
O – H	463

Bond breaking
2 x C – C = 348 x 2 = 696
8 x C – H = 413 x 8 = 3304
5 x O = O = 495 x 5 = 2475
Total = **6475**

Bond making
8 x O – H = 463 x 8 = 3704
6 x C = O = 6x

Δ_rH = ΣBond energies (bonds broken) − ΣBond energies (bonds formed)
6475 − 3704 − 6x = −2056 kJ mol^{-1}
6x = + 2056 + 6475 − 3704 = 4827   x = 805 kJ mol^{-1}