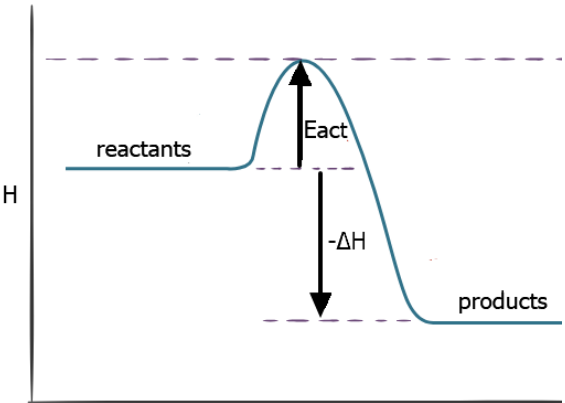
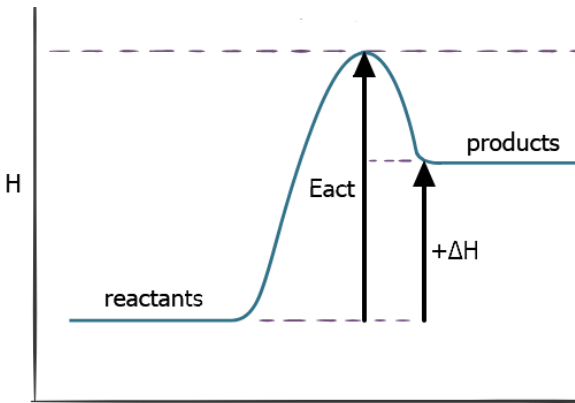



<p>Types of solids</p> <p>Ionic – between metal & non-metal (exception AlCl_3 covalent molecular); particles = IONS, attraction = ionic / electrostatic bond</p> <p>Covalent – between non-metal & non-metal.</p> <ul style="list-style-type: none"> • COVALENT NETWORK; particles = atoms; attraction = covalent bond; E.g. diamond, graphite and silicon dioxide SiO_2. • COVALENT MOLECULAR; particles = molecules; attraction = weak intermolecular forces; E.g. H_2O, I_2, CO_2. <p>Metallic – bonding between metal atoms; particles involved = atoms; attraction = metallic bond; E.g. Ag</p>	<p>Defining bond types</p> <p>Metallic: attraction between loosely held valence electrons & positively charged nuclei of neighbouring atoms. (Or “metal cations in sea of delocalised electrons”).</p> <p>Ionic: electrostatic attraction between oppositely charged ions.</p> <p>Covalent: bond in which one or more pairs of electrons are shared by two atoms. It’s intramolecular (between atoms <u>within</u> the molecule).</p> <p>Intermolecular: weak attraction between molecules (inter = between)</p>	<p>Explaining properties</p> <p>Hardness – indicates strong metallic, covalent or ionic bond – much energy needed to overcome attraction.</p> <p>High m.pt, b.pt – indicates strong attraction between atoms or ions (metallic, covalent network or ionic bond).</p> <p>Brittle – indicates ionic bond (when like force causes like charged ions to line up, they repel)</p> <p>Low m.pt, b.pt, soft - indicates weak attraction / weak intermolecular forces between molecules (molecular covalent)</p> <p>Malleable, ductile – indicates the non-directional attraction due to metallic bond.</p> <p>Conducts electricity – has “mobile charge carriers” – delocalised electrons in metals (and graphite) OR ions in molten or aqueous ionic substances. (NOTE: In solid ionic substances the ions are NOT free to move and so these solids are insulators).</p> <p>Solubility - non-polar substances e.g. I_2 dissolve in non-polar solvents e.g. hexane, because similar weak intermolecular forces exist between I_2 and hexane molecules as did between I_2 and I_2 molecules and between cyclohexane and cyclohexane molecules.</p> <p>Polar covalent molecules <u>and</u> many ionic solids dissolve in polar solvents (e.g. H_2O) due to attraction between charged particles. (DO NOT call ionic solids “polar” – they are not!)</p>
<p>Lewis structures</p> <p>Show only valence electrons. The number of valence electrons an atom has = its group # or group # – 10. E.g. Al – group 13. $13 - 10 = 3$ valence electrons. (Or work it out from Aluminium’s atomic number of 13 which gives an electron arrangement of 2, 8, 3).</p> <p>Pairs of electrons are drawn as ●● or x x</p> <p>Simple molecules have no more than four electron pairs about any atom (including multiple-bonded species). Be e.g. in BeCl_2 and B e.g. in BCl_3 are electron deficient (don’t have octet around Be or B, the central atom).</p> <div style="text-align: center;"> </div> <p>Usually, the central atom is already known. Otherwise the atom with the lowest electronegativity is the central atom. NOTE: H can never be a central atom.</p>	<p>Other useful Lewis diagrams</p> <div style="text-align: center;"> </div>	<p>Polar or non polar molecules</p> <p>To predict if a <u>bond</u> is polar, consider electronegativity, the ability of atoms <i>in a bond</i> to attract electrons to themselves. Inc \rightarrow periodic table, inc \uparrow a group. Remember (most)...F O N/Cl S/C H (less). Show EN with $\delta+$ and $\delta-$ above atoms. Predicting if <u>molecule</u> is polar or not. <i>Polar molecules:</i> contain polar bonds AND their lack of molecule symmetry means dipoles <u>do not</u> cancel out, e.g. H_2O. <i>Non-polar molecules:</i> usually contain polar bonds BUT the molecule symmetry means dipoles <u>do</u> cancel out e.g. $\text{O}=\text{C}=\text{O}$; OR do not contain polar bonds e.g. where atoms of same electronegativity are bonded, e.g. Cl-Cl.</p> <div style="text-align: center;"> </div>
<p>Bond angles and shapes of molecules</p> <p>Based on repulsion of regions of electron density. 2 regions – linear 180°, 3 regions – trigonal planar 120°, 4 regions tetrahedral – 109°.</p> <p>Shapes of molecules are based around repulsion of regions of electron density, bonding and non-bonding BUT ultimately depend on the positions of the atoms. E.g. NH_3</p> <ul style="list-style-type: none"> • 4 regions of electron density around the central N atom which repel each other to get as far away from each other as possible, taking up the <i>arrangement</i> of a regular tetrahedron. • 3 regions are bonding, one non-bonding – and so the shape of the molecule is trigonal pyramid / trigonal pyramidal, with bond angle of approximately 109°. 		

<p>Exothermic and endothermic reactions</p> <p>In an exothermic reaction, the reactants are at a higher energy level as compared to the products. The products are more stable than the reactants. Overall Δ_rH for the reaction is negative; energy is released in the form of heat.</p> <p>In the case of an endothermic reaction, the reactants are at a lower energy level compared to the products. The products are less stable than the reactants. The overall Δ_rH for the reaction is positive, i.e., energy is absorbed from the surroundings.</p>	<p style="text-align: center;">Exothermic</p>  <p style="text-align: center;">Reaction progress</p>	<p style="text-align: center;">Endothermic</p>  <p style="text-align: center;">Reaction Progress</p>	
<p>Enthalpy changes associated with the making and breaking of chemical bonds</p> <p>Average bond enthalpies have the units of kJ mol^{-1}. They show the energy required to break 1 mol of a particular bond.</p> <p>Average bond enthalpies are always listed as positive (+) numbers.</p>	<p style="text-align: center;">$\Delta_rH^\ominus = \sum(\text{bonds broken}) + \sum(\text{bonds formed})$ where \sum means “the sum of”</p> <p>You will need to decide if the value is + or - depending upon whether bonds are being broken (+) or formed (-).</p> <p>To break bonds is endothermic; to make the same bonds is exothermic</p> <p>e.g if $\text{O}=\text{O} \rightarrow \text{O} + \text{O} + 498 \text{ kJ mol}^{-1}$, then $\text{O} + \text{O} \rightarrow \text{O}=\text{O} - 498 \text{ kJ mol}^{-1}$</p> <p>Note: To turn $\text{C}=\text{C}$ into $\text{C}-\text{C}$, you must break $\text{C}=\text{C}$ and then make $\text{C}-\text{C}$.</p>		
<p>Exothermic and endothermic reactions including energy (enthalpy) changes associated with differing amounts of substances</p> <p>Calculations of energy changes using Δ_rH and reaction stoichiometry $n = cV$</p> <p>e.g. $\text{CH}_4(\text{g}) + 2\text{O}_2(\text{g}) \rightarrow \text{CO}_2(\text{g}) + 2\text{H}_2\text{O}(\text{l})$ $\Delta_rH = -890 \text{ kJ mol}^{-1}$. When 1 mol of $\text{CH}_4(\text{g})$ is completely burnt $\Delta_rH = -890 \text{ kJ mol}^{-1}$ OR 890 kJ of energy are released. $M(\text{CH}_4) = 16.0 \text{ g mol}^{-1}$.</p> <p>When 16.0 g of $\text{CH}_4(\text{g})$ is completely burnt $\Delta_rH = -890 \text{ kJ mol}^{-1}$ OR 890 kJ of energy are released. (Don't say -890 kJ of energy are released).</p> <p>What is the enthalpy change when 46.2 g of methane is burnt? Either do</p> <ul style="list-style-type: none"> by ratio: 16.0 g of $\text{CH}_4(\text{g})$ $\Delta_rH = -890 \text{ kJ mol}^{-1}$, so 46.2 g = $46.2/16.0 \times -890 = -2569$ $\Delta_rH = -2570 \text{ kJ}$ (3 s.f.) by mol: $n = m/M$ $n = 46.2/16.0 = 2.89 \text{ mol}$. For 1 mol, $\Delta_rH = -890 \text{ kJ mol}^{-1}$. For 2.89, $\Delta_rH = 2.89 \times -890 = -2570 \text{ kJ}$ (3 s.f.) 			
<p>Changes of state</p> <p>Solid \rightarrow Liquid \rightarrow Gas</p> <p>Endothermic; bond breaking.</p>		<p>Gas \rightarrow Liquid \rightarrow Solid</p> <p>Exothermic; bond making</p>	