

# **Chemistry A**

Advanced Subsidiary GCE

Unit **F322**: Chains, Energy and Resources

## **Mark Scheme for June 2012**

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All examiners are instructed that alternative correct answers and unexpected approaches in candidates' scripts must be given marks that fairly reflect the relevant knowledge and skills demonstrated.

Mark schemes should be read in conjunction with the published question papers and the report on the examination.

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
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## Annotations

Annotation	Meaning
	Benefit of doubt given
	Contradiction
	Incorrect response
	Error carried forward
	Ignore
	Not answered question
	Benefit of doubt not given
	Power of 10 error
	Omission mark
	Rounding error
	Error in number of significant figures
	Correct response

Annotation	Meaning
<b>DO NOT ALLOW</b>	Answers which are not worthy of credit
<b>IGNORE</b>	Statements which are irrelevant
<b>ALLOW</b>	Answers that can be accepted
( )	Words which are not essential to gain credit
—	Underlined words must be present in answer to score a mark
<b>ECF</b>	Error carried forward
<b>AW</b>	Alternative wording
<b>ORA</b>	Or reverse argument

### Subject-specific Marking Instructions

Annotations should be placed to clearly show where they apply within the body of the text (ie not in margins)

**Question 1** (c)(ii), (d)  
**Question 3** (c)(ii)  
**Question 4** (a)(i)  
**Question 5** (c)(i)  
**Question 6** (b)(iii)

All questions where an ECF has been applied.

### Checking additional pages

All the Additional Pages in the examination script must be checked to see if any candidates include any answers.

The only additional page is part of the last question, **6(b)(iii)**.

You must annotate page 20 with an omission mark ^ if the page is blank to show that you have checked this page.

## Generic comments

## ORGANIC STRUCTURES

For a 'structure' or 'structural formula',

- **ALLOW** correct structural **OR** displayed **OR** skeletal formula **OR** mixture of the above (as long as unambiguous)

For an alkyl group shown within a structure,

- **ALLOW** bond drawn to C or H,  
eg **ALLOW** CH<sub>3</sub>–, CH<sub>2</sub>–, C<sub>3</sub>H<sub>7</sub>–, etc
- **ALLOW** vertical 'bond' to any part of an alkyl group

For an OH group shown within a structure,

- **DO NOT ALLOW** formula with horizontal –HO **OR**
- **ALLOW** vertical 'bond' to any part of the OH group

For a CHO group shown within a structure,

- **DO NOT ALLOW** COH

For a 3-D structure,

• For bond in the plane of paper, a solid line is expected:	
• For bond out of plane of paper, a solid wedge is expected:	
• For bond into plane of paper, <b>ALLOW</b> :	
• <b>ALLOW</b> a hollow wedge for 'in bond' <b>OR</b> an 'out bond', provided it is different from the other in or out wedge eg:	

**NAMES**

Names including alkyl groups:

- **ALLOW** alkanyl, eg ethanyl (ie **IGNORE** 'an')
- **DO NOT ALLOW** alkol, eg ethol (ie 'an' is essential)

Names of esters:

- Two words are expected, eg ethyl ethanoate
- **ALLOW** one word, eg ethylethanoate

Names with multiple numbers and hyphens:

Use of 'e'

- **ALLOW** superfluous 'e', eg propane-1-ol ('e' is kept if followed by consonant)
- **ALLOW** absence of 'e', eg propan-1,2-diol ('e' is omitted if followed by vowel)

Hyphens separate name from numbers:

- **ALLOW** absence of hyphens, eg propane 1,2 diol

Multiple locant numbers must be clearly separated:

- **ALLOW** full stops: eg 1.2 **OR** spaces: 1 2
- **DO NOT ALLOW** eg 12

Locant numbers in formula must be correct

- **DO NOT ALLOW** propan-3-ol

Order of substituents should be alphabetical:

- **ALLOW** any order (as long as unambiguous), eg 2-chloro-3-bromobutane

**ABBREVIATIONS**

van der Waals' forces

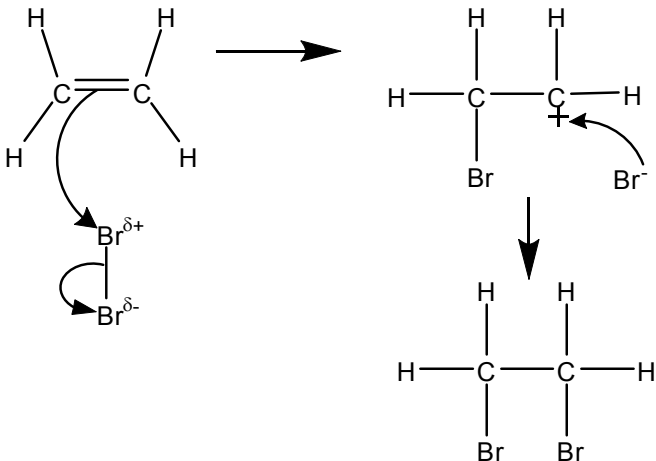
- **ALLOW** vdw forces **OR** VDW forces (and any combination of upper and lower cases)

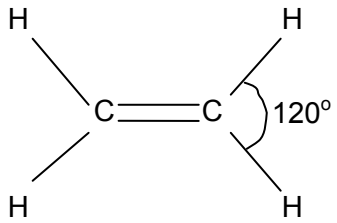
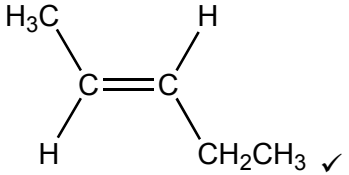
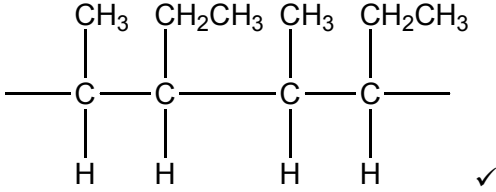
Question		Answer	Marks	Guidance
1	(a)	Because hydrocarbons have different boiling points ✓	1	<p><b>ALLOW</b> each fraction / component / substance / molecule / compound / fuel has a different boiling temperatures</p> <p><b>ALLOW</b> condense at different temperatures</p> <p><b>ALLOW</b> because van der Waals' forces differ with molecular size</p> <p><b>IGNORE</b> references to volatility</p> <p>different strength of intermolecular forces is <b>not</b> sufficient</p>
	(b)	<p><b>Any one from:</b></p> <p>Bio-fuels produce less carbon dioxide (overall) <b>OR</b> petrol or diesel produce more carbon dioxide (overall) ✓</p> <p>Bio-fuels are renewable <b>OR</b> petrol and diesel are non-renewable ✓</p> <p>Allows crude oil to be used to make other products <b>OR</b> petrochemicals (rather than petrol) <b>OR</b> Save crude oil <b>OR</b> no risk of large scale pollution from exploitation of crude oil ✓</p>	1	<p><b>ASSUME</b> 'they' or 'it' refers to biofuels</p> <p><b>ALLOW</b> bio-fuels are (more) carbon-neutral <b>OR</b> plants take up the carbon dioxide released during combustion</p> <p><b>ALLOW</b> lower carbon footprint</p> <p><b>ALLOW</b> plants are a renewable resource / crude oil non-renewable resource / bio-diesel is more sustainable / diesel is not sustainable / petrol and diesel are made from a finite resource / petrol and diesel will run out / bio-fuels will not run out</p> <p><b>ALLOW</b> decrease the need for fossil fuels</p> <p><b>IGNORE</b> can be used by diesel powered cars with or without any conversion</p>
	(c) (i)	Idea that <b>carbon-carbon</b> bonds can break anywhere ✓	1	<p>The answer must refer to carbon-carbon bonds or the carbon chain</p> <p><b>ALLOW</b> (carbon) <b>chain</b> can break anywhere</p> <p>Bonds can break anywhere is not sufficient</p>

Question			Answer	Marks	Guidance
1	(c)	(ii)	<p>Correct identification of <math>C_2H_3^+</math> for <math>m/z = 27</math> ✓</p> <p>Some indication to explain how the identity of propene was deduced  <b>OR</b> further analysis of the mass spectrum ✓</p> <p>Correct identification of the alkene as <math>C_3H_6</math> <b>OR</b> propene ✓</p> <p><math>C_{12}H_{26} \rightarrow C_3H_8 + 3C_3H_6</math> ✓</p>	4	<p><b>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</b></p> <p><b>ALLOW</b> <math>CHCH_2^+</math>  <b>DO NOT ALLOW</b> <math>C_2H_3</math> (the positive charge is essential) <b>OR</b> <math>CCH_3^+</math></p> <p><b>ALLOW</b> Molecular ion/<math>M^+</math>/<math>M</math> is <math>m/z = 42</math> <b>OR</b> <math>m/z = 15</math> is <math>CH_3</math>  <b>ALLOW</b> mass spectrum shows <math>M_r = 42</math>  <b>ALLOW</b> idea that alkane <math>C_{12}H_{26} - C_3H_8</math> can only give <math>3C_3H_6</math></p> <p><b>ALLOW</b> prop-1-ene  An incorrect formula for the alkene in the equation will not contradict this answer</p> <p><b>ALLOW</b> <math>C_3H_6</math> from its use in an equation even if the equation is wrong providing there has not been an attempt elsewhere to identify the alkene</p> <p><b>ALLOW</b> correct displayed <b>OR</b> structural <b>OR</b> skeletal <b>OR</b> molecular formulae in the equation</p>

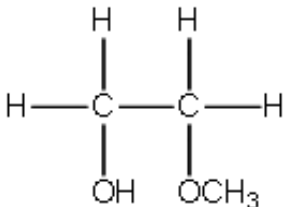
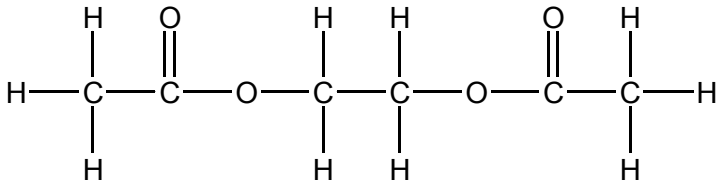


Question		Answer	Marks	Guidance
1	(d)	<p>React with bromine <b>OR</b> <math>C_2H_4 + Br_2 \rightarrow C_2H_4Br_2</math> ✓</p> <p>React with hydrogen bromide <b>OR</b> <math>C_2H_4 + HBr \rightarrow C_2H_5Br</math> ✓</p> <p>React with steam <b>OR</b> heat with water <b>OR</b> <math>C_2H_4 + H_2O(g) \rightarrow C_2H_5OH</math> ✓</p> <p>acid (catalyst) ✓</p>	9	<p><b>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</b></p> <p><b>ALLOW</b> reactants even from incorrect equations</p> <p><b>ALLOW</b> reactants or conditions over the arrow</p> <p><b>ALLOW</b> <math>Br_2</math> mark from the mechanism even if the mechanism is incorrect</p> <p><b>IGNORE</b> conditions unless they would lead to a different reaction with ethene</p> <p><b>IGNORE</b> conditions unless they would lead to a different reaction with ethene</p> <p><b>ALLOW</b> temperature range between 100–400 °C if quoted</p> <p><b>IGNORE</b> reference to pressure</p> <p><b>IGNORE</b> hydrolysis</p> <p>Hydration is not sufficient but <b>DO NOT ALLOW</b> hydrogenation</p> <p><b>ALLOW</b> <math>H_2SO_4</math> <b>OR</b> <math>H_3PO_4</math> <b>OR</b> <math>H^+</math></p> <p><b>DO NOT ALLOW</b> <math>HCl</math>, <math>HBr</math> etc.</p> <p><b>ALLOW</b> two stage process e.g. react with <math>HBr</math> one mark followed by <math>KOH(aq)</math> one mark</p>

Question	Answer	Marks	Guidance
	<p>Electrophilic addition ✓</p> <p>Curly arrow from double bond to attack <math>\text{Br}^{\delta+}</math> of <math>\text{Br}-\text{Br}</math> and breaking of <math>\text{Br}-\text{Br}</math> bond ✓</p> <p>Correct dipoles shown on <math>\text{Br}^{\delta+}-\text{Br}^{\delta-}</math> ✓</p> <p>Correct carbonium / carbocation ion drawn ✓</p> <p>Curly arrow from <math>\text{Br}^-</math> to the carbonium ion <b>and</b> correct product shown ✓</p> 		<p>Curly arrow must start from the double bond and not a carbon atom and go to the <math>\text{Br}^{\delta+}</math>; other curly arrow must start from <math>\text{Br}-\text{Br}</math> bond.</p> <p><b>ALLOW</b> attack of <math>\text{Br}-\text{Br}</math> if dipoles not shown  <b>DO NOT ALLOW</b> attack of <math>\text{Br}^{\delta-}</math></p> <p>Dipole must be partial charge and not full charge  <b>DO NOT ALLOW</b> any other partial charges eg on the double bond</p> <p>Carbocation needs a full charge and not a partial charge (charges do not need to be surrounded by a circle)  All atoms in the carbocation must be shown</p> <p><math>\text{Br}^-</math> curly arrow must come from one lone pair on <math>\text{Br}^-</math> ion <b>OR</b> from minus sign on <math>\text{Br}^-</math> ion  Lone pair does not need to be shown on <math>\text{Br}^-</math> ion</p> <p><b>ALLOW</b> mechanism which goes via a cyclic bromonium ion instead of the carbocation</p> <p><b>SEE EXTRA ADVICE ABOUT CURLY ARROWS ON PAGE 30</b></p>

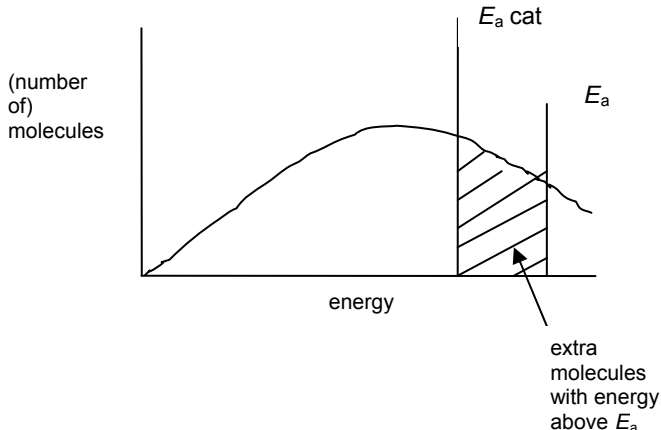
Question		Answer	Marks	Guidance
1	(e)	<p>Correct shape ✓</p>  <p>120° ✓</p> <p>Three areas of electron density repel each other ✓</p>	3	<p><b>IGNORE</b> any name of shape given</p> <p><b>ALLOW</b> 115–125°</p> <p><b>ALLOW</b> even if it is the C–C–H shown on a diagram.</p> <p><b>ALLOW</b> three or four electron pairs repel <b>OR</b> three or four bonds repel</p> <p><b>IGNORE</b> does not have any lone pairs</p> <p><b>DO NOT ALLOW</b> atoms repel / electrons repel</p> <p><b>DO NOT ALLOW</b> has lone pair which repels more</p>
	(f) (i)		1	<p><b>ALLOW</b> correct structural <b>OR</b> displayed <b>OR</b> skeletal formula</p> <p><b>OR</b> mixture of the above (as long as unambiguous)</p>
	(ii)		1	<p><b>ALLOW</b> correct structural <b>OR</b> displayed <b>OR</b> skeletal formula</p> <p><b>OR</b> mixture of the above (as long as unambiguous)</p> <p><b>ALLOW</b> CH<sub>3</sub> and C<sub>2</sub>H<sub>5</sub> groups above or below chain</p> <p><b>ALLOW</b> bond to ethyl and methyl group to any part of ethyl or methyl group</p> <p><b>IGNORE</b> any brackets drawn</p> <p><b>ALLOW</b> two or more repeat units but has to have a whole number of repeat units (<i>ie</i> does not have to be two)</p> <p>'End bonds' <b>MUST</b> be shown and can be dotted</p> <p><b>IGNORE</b> <i>n</i></p>
<b>Total</b>			<b>21</b>	

Question		Answer	Marks	Guidance
2	(a) (i)	$2\text{C}_2\text{H}_4 + \text{O}_2 \rightarrow 2\text{C}_2\text{H}_4\text{O}$ ✓	1	<b>ALLOW</b> molecular formulae <b>OR</b> correct structural <b>OR</b> displayed <b>OR</b> skeletal formula <b>OR</b> mixture of the above (as long as unambiguous) <b>ALLOW</b> correct multiples, including fractions, of this equation <b>IGNORE</b> state symbols <b>DO NOT ALLOW</b> [O]
	(ii)	$\text{C}_2\text{H}_4 + 3\text{O}_2 \rightarrow 2\text{CO}_2 + 2\text{H}_2\text{O}$ ✓	1	<b>ALLOW</b> molecular formulae <b>OR</b> correct structural <b>OR</b> displayed <b>OR</b> skeletal formula <b>OR</b> mixture of the above (as long as unambiguous) <b>ALLOW</b> correct multiples of this equation <b>IGNORE</b> state <b>symbols</b> <b>DO NOT ALLOW</b> [O]
	(b) (i)	✓ 	1	Only one carbon atom needs to have the correct partial charge <b>DO NOT ALLOW</b> partial charges on hydrogen atoms
	(ii)	Movement of an electron pair ✓	1	<b>ALLOW</b> movement of a lone pair <b>OR</b> movement of a bond <b>ALLOW</b> movement of two electrons
	(iii)	Heterolytic ✓  Both electrons (in the bond) go to the same atom <b>OR</b> (bond breaks) to make a cation and (a lone pair on the oxygen atom) <b>OR</b> bond pair becomes a lone pair on oxygen ✓	2	<b>MARK INDEPENDENTLY</b>  <b>ALLOW</b> one atom gets none of the bonded electrons <b>DO NOT ALLOW</b> both electrons go to a molecule <b>DO NOT ALLOW</b> makes a positive and a negative ion because in this example this is <b>not</b> true
	(iv)	It donates a <b>pair</b> of electrons ✓	1	<b>ALLOW</b> donates a lone pair <b>DO NOT ALLOW</b> it donates electrons
	(v)	idea that $\text{H}^+$ ion is used in <b>step 1 AND</b> made in <b>step 4</b> ✓	1	<b>ALLOW</b> $\text{H}^+$ ion is used at the <b>start AND</b> made at the <b>end</b> <b>IGNORE</b> overall $\text{H}^+$ is not used up in the mechanism

Question		Answer	Marks	Guidance
2	(b) (vi)		1	<p><b>ALLOW</b> correct structural <b>OR</b> displayed <b>OR</b> skeletal formula <b>OR</b> mixture of the above (as long as unambiguous) eg CH<sub>2</sub>OHCH<sub>2</sub>OCH<sub>3</sub></p> <p><b>ALLOW</b> vertical 'bond' to any part of the OH or OCH<sub>3</sub> group <b>DO NOT ALLOW</b> formula with horizontal –HO <b>OR</b> OH– <b>DO NOT ALLOW</b> formula with horizontal –CH<sub>3</sub>O <b>OR</b> OCH<sub>3</sub>–</p>
	(c)	<p>Ethane-1,2-diol has more OH groups (than ethanol) ✓</p> <p>Stronger hydrogen bonding (between ethane-1,2-diol molecules) ✓</p>	2	<p><b>ALLOW</b> has more hydroxyl groups <b>OR</b> has more hydroxy groups <b>OR</b> has more alcohol groups Ethane-1,2-diol has two OH groups is <b>NOT</b> sufficient but <b>ALLOW</b> ethane-1,2-diol has two OH groups and ethanol has one <b>DO NOT ALLOW</b> it has hydroxide (ions)</p> <p><b>ALLOW</b> more hydrogen bonds (between ethane-1,2-diol molecules) <b>IGNORE</b> hydrogen bonds with water</p>
	(d)	<p>One ester linkage drawn despite the rest of the structure ✓</p> <p>Correct structure for example CH<sub>3</sub>COOCH<sub>2</sub>CH<sub>2</sub>OOCCH<sub>3</sub> <b>OR</b></p> 	2	<p><b>ALLOW</b> correct structural <b>OR</b> displayed <b>OR</b> skeletal formula <b>OR</b> mixture of the above (as long as unambiguous)</p> <p><b>ALLOW</b> ester shown as all the atoms <b>OR</b> as –COOC– <b>OR</b> –CH<sub>2</sub>OOC– <b>OR</b> –CH<sub>2</sub>OCOC–</p> <p><b>IGNORE</b> molecular formula</p>

Question	Answer	Marks	Guidance
2 (e)	<p><i>Any two from:</i></p> $\begin{array}{c} \text{H} \quad \text{H} \\   \quad   \\ \text{O}=\text{C}-\text{C}=\text{O} \end{array} \quad \begin{array}{c} \text{O} \quad \text{O} \\    \quad    \\ \text{H}-\text{C}-\text{C}-\text{O}-\text{H} \end{array}$ $\begin{array}{c} \text{H} \quad \text{H} \\   \quad   \\ \text{H}-\text{O}-\text{C}-\text{C}=\text{O} \\   \\ \text{H} \end{array} \quad \begin{array}{c} \text{H} \quad \text{O} \\   \quad    \\ \text{H}-\text{O}-\text{C}-\text{C}-\text{O}-\text{H} \\   \\ \text{H} \end{array}$ $\text{H}-\text{O}-\overset{\text{O}}{\parallel}{\text{C}}-\overset{\text{O}}{\parallel}{\text{C}}-\text{O}-\text{H}$ <p>✓✓</p>	2	<p>Mark incorrect answers first</p> <ul style="list-style-type: none"> <li>• If one incorrect answer maximum of 1 mark</li> <li>• If two incorrect answers award 0 marks</li> </ul> <p><b>ALLOW</b> OH instead of <math>-\text{O}-\text{H}</math></p> <p><b>ALLOW</b> vertical 'bond' to any part of the OH  <b>DO NOT ALLOW</b> formula with horizontal <math>-\text{HO}</math> <b>OR</b> <math>\text{OH}-</math>  but <b>ALLOW ECF</b> if both displayed formulae are drawn this way</p> <p><b>ALLOW one mark if two correct structural OR skeletal formula OR mixture of the above (as long as unambiguous) are drawn</b></p>
	<b>Total</b>	<b>15</b>	

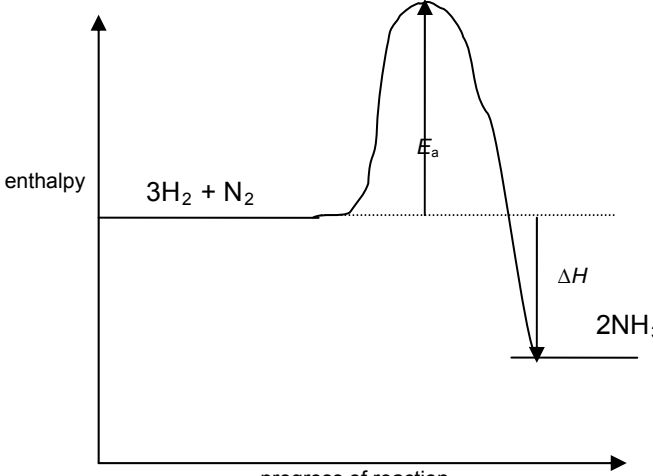
Question		Answer	Marks	Guidance
3	(a)	(equilibrium position shifts) to the left ✓  (because there are) fewer moles (of gas) on the reactant side <b>OR</b> (there are) more moles (of gas) on product side ✓ This explanation mark is dependent on the correct shift of the equilibrium	2	<b>Note: ALLOW</b> suitable alternatives for 'to left', eg: towards CH <sub>4</sub> or H <sub>2</sub> O / towards reactants <b>OR</b> in backward direction <b>OR</b> in reverse direction <b>OR</b> decreases yield of CO or H <sub>2</sub> /products <b>ALLOW</b> 'favours the left', as alternative for 'shifts equilibrium to left'  <b>ALLOW</b> fewer molecules on reactant side <b>OR</b> smaller volume on the left hand side <b>ALLOW ORA</b> if specified <b>IGNORE</b> responses in terms of rate
	(b)	(equilibrium position shifts) to the right ✓  (because forward) reaction is endothermic <b>OR</b> <b>reverse</b> reaction is exothermic ✓ This explanation mark is dependent on the correct shift of the equilibrium	2	<b>Note: ALLOW</b> suitable alternatives for 'to right', eg: towards CO or H <sub>2</sub> / towards products <b>OR</b> in forward direction <b>OR</b> increases yield of CO or H <sub>2</sub> /products <b>OR</b> decreases amount of CH <sub>4</sub> or H <sub>2</sub> O/reactants <b>ALLOW</b> 'favours the right', as alternative for 'shifts equilibrium to right'  <b>ALLOW</b> reaction takes in heat  <b>ALLOW reverse</b> reaction gives out heat  <b>ALLOW ORA</b> if specified <b>IGNORE</b> responses in terms of rate
	(c) (i)	Gives a high rate of reaction <b>OR</b> reaction is fast <b>OR</b> reasonable rate of reaction without shifting equilibrium too much to the left ✓	1	<b>ALLOW</b> if greater pressure used it increases safety risk <b>ALLOW</b> if greater pressure used it is more expensive <b>ALLOW</b> higher pressure will shift equilibrium position even more to the left  It is a compromise on its own is <b>not</b> sufficient but <b>ALLOW</b> compromise between rate and yield <b>OR</b> between rate and safety

Question	Answer	Marks	Guidance
3 (c) (ii)	<p>y-axis label is '(fraction of or number of) molecules' <b>AND</b> x-axis label is 'energy' <b>AND</b> correct curve ✓</p> <p>Lowers activation energy ✓</p> <p>More molecules with energy above activation energy with a catalyst <b>OR</b> more <b>effective</b> collisions <b>OR</b> more <b>successful</b> collisions ✓</p> 	3	<p><b>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</b></p> <p>Boltzmann distribution - must start at origin and must not end up at 0 on y-axis ie must not touch x-axis  <b>ALLOW</b> particles <b>OR</b> moles as y-axis label  <b>IGNORE</b> minor point of inflexion in the curve</p> <p><b>DO NOT ALLOW</b> two curves  <b>DO NOT ALLOW</b> atoms but credit atoms if used in a second marking point  <b>DO NOT ALLOW</b> enthalpy for x-axis label</p> <p><b>ALLOW</b> this mark from a labelled diagram</p> <p>more collisions per second is not sufficient</p>



Question		Answer	Marks	Guidance
3	(d)	<p><b>FIRST, CHECK THE ANSWER ON ANSWER LINE</b>  <b>IF answer = 91.2 (%) award 3 marks</b></p> <p>theoretical amount of hydrogen = <math>3.75 \times 10^7</math> (mol) ✓</p> <p>actual amount of hydrogen made = <math>3.42 \times 10^7</math> (mol) ✓</p> <p>% = 91.2 ✓</p>	3	<p><b>IF</b> there is an alternative answer, check to see if there is any <b>ECF</b> credit possible using working below</p> <p>IF ECF, <b>ANNOTATE WITH TICKS AND CROSSES, etc</b></p> <p>Answer must have <b>three</b> significant figures  <b>ALLOW</b> ECF from incorrect theoretical and actual amounts of hydrogen</p> <p><b>ALLOW</b> answer that uses grams rather than tonnes where theoretical amount of hydrogen = 37.5 (mol) and actual amount of hydrogen = 34.2 (mol)</p> <p><b>ALLOW</b> alternative approach based on the mass of hydrogen rather than the amount of hydrogen  Theoretical amount of hydrogen = <math>3.75 \times 10^7</math> (mol) ✓  Theoretical mass of hydrogen made = 75 (tonnes) ✓  Percentage = 91.2 ✓</p>
	(e) (i)	$\text{CO} + 2\text{H}_2 \rightarrow \text{CH}_3\text{OH}$ ✓	1	<p><b>ALLOW</b> correct multiples  <b>ALLOW</b> <math>\text{CH}_4\text{O}</math>  <b>IGNORE</b> state symbols</p>
	(ii)	<p><b>Any two from:</b></p> <p>Carbon monoxide is toxic <b>OR</b> poisonous ✓</p> <p>Increases atom economy of the process <b>OR</b> gives 100% atom economy ✓</p> <p>Methanol is a fuel ✓</p>	2	<p><b>IGNORE</b> harmful or dangerous</p> <p><b>ALLOW</b> uses a waste product <b>OR</b> CO is then a desired product <b>OR</b> CO is no longer a waste product <b>OR</b> reduces amount of waste product</p> <p><b>ALLOW</b> other uses of methanol eg petrol additive, solvent or organic feedstock</p>

Question		Answer	Marks	Guidance
3	(f)	Unsaturated (vegetable) oils OR oils containing C=C bonds ✓  (reacted with hydrogen) in the presence of a nickel catalyst ✓	2	<b>ALLOW</b> unsaturated fats OR unsaturated lipids OR unsaturated ester <b>ALLOW</b> oils become more saturated <b>IGNORE</b> unsaturated compound <b>DO NOT ALLOW</b> unsaturated hydrocarbon  <b>ALLOW</b> Pt OR Pd
<b>Total</b>			<b>16</b>	

Question	Answer	Marks	Guidance
4 (a) (i)	<p>2NH<sub>3</sub> added as product ✓</p> <p><math>\Delta H</math> labelled with product below reactant <b>AND</b> arrow downwards ✓</p> <p><math>E_a</math> labelled correctly <b>AND</b> above reactants ✓</p> 	3	<p><b>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</b></p> <p><b>IGNORE</b> state symbol <b>ALLOW</b> product mark even if product line above the reactant line</p> <p><b>ALLOW</b> -92 as a label for <math>\Delta H</math> <b>ALLOW</b> this line even if it has a small gap at the top and bottom ie does not quite reach reactant or product line</p> <p>The curve must be drawn for this marking point</p> <p><b>IGNORE</b> arrows at both ends of activation energy line but <b>DO NOT ALLOW</b> arrow pointing down The <math>E_a</math> line must go to maximum (or near to the maximum) on the curve <b>ALLOW</b> if the line clearly shows an activation energy and is not an enthalpy change <b>ALLOW</b> this line even if it has a small gap at the top and bottom ie does not quite reach the maximum or reactant line</p>

Question			Answer	Marks	Guidance
4	(a)	(ii)	$-46 \text{ (kJ mol}^{-1}\text{)} \checkmark$	1	<b>DO NOT ALLOW</b> 46 with no sign
		(iii)	Any value between +1 to +249 (kJ mol <sup>-1</sup> ) ✓	1	+ sign is not needed
		(iv)	$+342 \text{ (kJ mol}^{-1}\text{)} \checkmark$	1	+ sign is not needed
	(b)	(i)	$2\text{CO} + 2\text{NO} \rightarrow 2\text{CO}_2 + \text{N}_2 \checkmark$	1	<b>ALLOW</b> correct multiples

Question			Answer	Marks	Guidance
4	(b)	(ii)	<p>CO and NO are adsorbed (onto surface) <b>OR</b> reactants are adsorbed (onto surface) ✓</p> <p>weakening of bonds <b>OR</b> chemical reaction <b>OR</b> new bonds are made <b>OR</b> carbon dioxide and nitrogen are made ✓</p> <p>CO<sub>2</sub> and N<sub>2</sub> desorbs (from the surface) <b>OR</b> products desorbs (from the surface) ✓</p>	3	<p><b>ALLOW</b> CO and NO stick onto surface <b>OR</b> CO and NO form weak attractions to the surface <b>OR</b> gases are adsorbed onto surface <b>OR</b> gases bond to surface</p> <p><b>NOT</b> absorb but <b>allow</b> ecf for deabsorb later on</p> <p><b>ALLOW</b> lowers activation energy</p> <p><b>IGNORE</b> alternative pathway</p> <p>Requires less energy is not sufficient</p> <p><b>ALLOW</b> products leave (the surface) <b>OR</b> products diffuse away (from surface) <b>OR</b> weak attraction to surface is broken</p> <p><b>ALLOW</b> deadsorb</p>

Question			Answer	Marks	Guidance
4	(c)	(i)	<p><b>Any two from:</b></p> <p>IR (spectroscopy) ✓</p> <p>Mass spectrometry ✓</p> <p>UV (spectroscopy) ✓</p> <p>NMR ✓</p> <p>GC ✓</p>	2	<p><b>ALLOW</b> mass spec / MS / mass spectroscopy</p> <p><b>ALLOW</b> atomic absorption / AAS</p> <p><b>IGNORE</b> satellite imaging or thermal imaging</p>
		(ii)	<p><b>Any one from:</b></p> <p>Idea that pollution travels (across country) borders</p> <p><b>OR</b> idea that all countries contribute towards pollution</p> <p><b>OR</b> Cooperation means that scientists can share ideas</p> <p><b>OR</b> scientists can warn governments of risk</p> <p><b>OR</b> world-wide legislation can be introduced</p> <p><b>OR</b> allows monitoring of pollution in different countries</p> <p><b>OR</b> richer countries can help poorer countries introduce pollution controls</p> <p><b>OR</b> One country cannot control pollution unless all countries do ✓</p>	1	<p><b>ALLOW</b> some countries produce more pollution than others</p> <p><b>ALLOW</b> so protocols can be developed</p>
	(d)		<p><b>Step 1</b> <math>\text{NO} + \text{O}_3 \rightarrow \text{NO}_2 + \text{O}_2</math> ✓</p> <p><b>Step 2</b> <math>\text{NO}_2 + \text{O} \rightarrow \text{NO} + \text{O}_2</math> ✓</p> <p><b>overall</b> <math>\text{O}_3 + \text{O} \rightarrow 2\text{O}_2</math> ✓</p>	3	

Question			Answer	Marks	Guidance
4	(e)	(i)	Reaction gives NO <b>OR</b> reaction gives NO <sub>2</sub> <b>OR</b> reaction gives a mixture of oxides <b>OR</b> activation energy too high <b>OR</b> rate of reaction is too slow ✓	1	<b>ALLOW</b> makes a mixture of oxides/products <b>ALLOW</b> reaction cannot be carried out experimentally <b>ALLOW</b> reaction does not take place nitrogen and oxygen do not react together is <b>not</b> sufficient <b>IGNORE</b> heat loss to surroundings <b>IGNORE</b> reference to bond enthalpy being a mean value
		(ii)	<b>FIRST, CHECK THE ANSWER ON ANSWER LINE</b> <b>IF answer = +82 (kJ mol<sup>-1</sup>) award 2 marks</b> <b>IF answer = -82 (kJ mol<sup>-1</sup>) award 1 mark</b>  $\Delta H = 193 - 111$ ✓  $= +82$ ✓	2	<b>ALLOW</b> 82  <b>ALLOW</b> one mark for -82 <b>ALLOW</b> one mark for +304 / -304
			<b>Total</b>	<b>19</b>	

Question		Answer	Marks	Guidance
5	(a)	<p><b>FIRST, CHECK THE ANSWER ON ANSWER LINE</b>  <b>IF answer = <math>-162 \text{ (kJ mol}^{-1}\text{)}</math> award 3 marks</b></p> <p>Energy associated with bond breaking = 3354  <b>OR</b> <math>(2 \times 805) + (4 \times 436) \checkmark</math></p> <p>Energy associated with bond making = 3516  <b>OR</b> <math>(4 \times 415) + (4 \times 464) \checkmark</math></p> <p>Enthalpy change = <math>-162 \checkmark</math></p>	3	<p><b>IF</b> there is an alternative answer, check to see if there is any <b>ECF</b> credit possible using working below.</p> <p>IF ECF, <b>ANNOTATE WITH TICKS AND CROSSES, etc</b></p> <p><b>IGNORE</b> sign</p> <p><b>IGNORE</b> sign</p> <p><b>ALLOW</b> ECF from wrong additions of energy associated with bond breaking and/or from bond making</p> <p><b>ALLOW</b> two marks for <math>(+162, (+)6870, -6870</math> or <math>(+)766</math></p> <p><b>ALLOW</b> one mark for <math>-766</math></p>
	(b) (i)	<p>Absorbs IR radiation <math>\checkmark</math></p> <p>Bonds vibrate <math>\checkmark</math></p>	2	<p><b>IGNORE</b> absorbs heat</p> <p><b>ALLOW</b> IR re-radiated</p> <p><b>DO NOT ALLOW</b> absorbs UV radiation</p> <p><b>DO NOT ALLOW</b> blocks IR radiation</p> <p><b>ALLOW</b> bonds stretch <b>OR</b> bonds bend</p> <p><b>IGNORE</b> molecule vibrates/rotates</p> <p><b>DO NOT ALLOW</b> bonds break</p>



Question			Answer	Marks	Guidance
5	(b)	(ii)	<p><b>Any two from:</b></p> <p>(liquid) injected deep into the oceans ✓</p> <p>Stored in (old) geological formations  <b>OR</b> stored underground in rocks  <b>OR</b> stored in (old) mines  <b>OR</b> stored in (old) oil wells ✓</p> <p>Stored by reaction with metal <u>oxides</u>  <b>OR</b> reaction to form (solid) <u>carbonates</u>  <b>OR</b> stored as a <u>carbonate</u>  <b>OR</b> equation to show formation of metal carbonate ✓</p>	2	<p><b>DO NOT ALLOW</b> reference to carbon being stored – the answer must either refer to carbon dioxide or not mention the name of the stored substance</p> <p><b>ALLOW store deep</b> in the oceans <b>OR</b> on the <b>sea-bed</b> ✓  <b>ALLOW</b> stored deep under the sea  <b>DO NOT ALLOW</b> dissolve CO<sub>2</sub> in the sea <b>OR</b> stored in ocean</p> <p><b>ALLOW</b> stored under the sea bed  <b>ALLOW</b> pumped into oil wells to force last bit of oil out</p> <p><b>IGNORE</b> mineral storage</p>

Question			Answer	Marks	Guidance
5	(c)	(i)	<p>Homolytic ✓</p> <p><math>\text{Br}_2 \longrightarrow 2\text{Br}</math> ✓</p> <p><math>\text{Br} + \text{C}_2\text{H}_6 \longrightarrow \text{HBr} + \text{C}_2\text{H}_5</math> ✓</p> <p><math>\text{C}_2\text{H}_5 + \text{Br}_2 \longrightarrow \text{C}_2\text{H}_5\text{Br} + \text{Br}</math> ✓</p> <p><math>\text{Br} + \text{C}_2\text{H}_5 \longrightarrow \text{C}_2\text{H}_5\text{Br}</math></p> <p><b>OR</b> <math>\text{Br} + \text{Br} \longrightarrow \text{Br}_2</math></p> <p><b>OR</b> <math>\text{C}_2\text{H}_5 + \text{C}_2\text{H}_5 \longrightarrow \text{C}_4\text{H}_{10}</math> ✓</p> <p><b>Two</b> names of steps linked to appropriate equations ✓</p> <p><b>OR</b></p> <p><b>three</b> names of steps linked to appropriate equations ✓✓</p>	7	<p><b>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</b></p> <p><b>IGNORE</b> dot for radical and any state symbols for all equations</p> <p>If more than one termination step is written they must all be correct to be awarded the mark</p> <p><b>DO NOT ALLOW</b> termination steps with H</p> <p>initiation step linked to correct equation</p> <p>propagation step linked to one equation in which there is a radical on the left and a radical on the right</p> <p>termination step linked to equation involving two radicals:</p> <p>If no equations are given to link the names of the step then award one mark for mention of all three steps</p> <p>If halogen other than bromine do not give equation mark for initiation and only give one mark for all three terms linked to appropriate equations</p> <p>If hydrocarbons other than ethane are used <b>DO NOT ALLOW</b> any marks for the equations in the propagation steps</p>

Question			Answer	Marks	Guidance
5	(c)	(ii)	<p><b>Any two from:</b></p> <p>More than one C–H bond can be substituted <b>OR</b> multi-substitution can occur <b>OR</b> more than one substitution can happen ✓</p> <p>Lots of termination steps ✓</p> <p>termination steps can give products that will also react with (bromine) radicals ✓</p>	2	<p><b>ALLOW</b> equations or examples of multi substitution</p> <p><b>ALLOW</b> an equation to illustrate formation of other products eg butane</p> <p><b>ALLOW</b> examples of other products that can be formed in termination steps eg bromobutane</p> <p><b>ALLOW</b> examples of products eg butane reacting with bromine radicals to give bromobutane</p>
			<b>Total</b>	<b>16</b>	

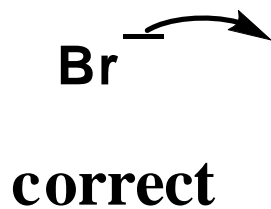
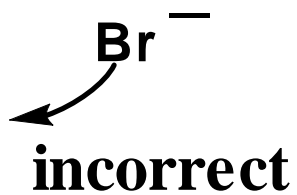
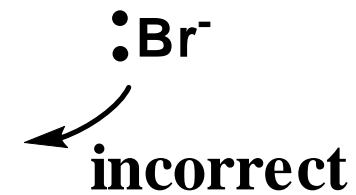
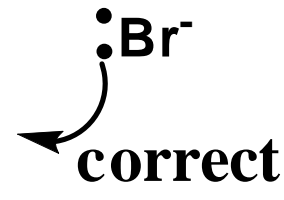
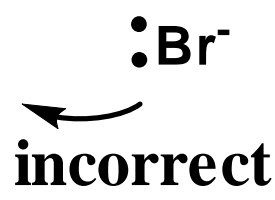
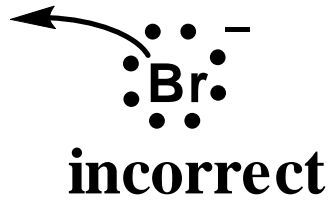
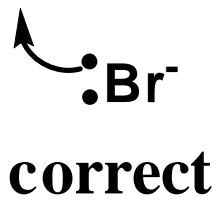
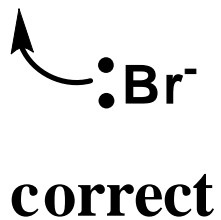
Question	Answer	Marks	Guidance
6 (a)	1-bromopentane reacts faster <b>OR</b> 1-chloropentane reacts slower ✓  C–C/ stronger bond (than C–Br bond) <b>OR</b> C–C/ shorter bond (than C–Br bond) <b>OR</b> C–C/ bond is harder to break <b>OR</b> needs more energy to break C–C/ bond <b>OR</b> bond enthalpy of C–C/ greater (than C–Br bond) ✓	2	<b>ALLOW</b> takes more time to react <b>ALLOW</b> chloro compound reacts slower than bromine compound <b>DO NOT ALLOW</b> bromine reacts faster than chlorine  <b>ALLOW</b> ORA  Answer must refer to the C–C/ bond or C–Br bonds
(b) (i)	$\text{CH}_3\text{—CH}_2\text{—CH}_2\text{—CH}_2\text{—I} \quad \checkmark$ $\begin{array}{c} \text{CH}_3\text{—CH}_2\text{—CH—CH}_3 \\   \\ \text{I} \end{array} \quad \checkmark$ $\begin{array}{c} \text{CH}_3 \\   \\ \text{CH}_3\text{—C—I} \\   \\ \text{CH}_3 \end{array} \quad \checkmark$ $\begin{array}{c} \text{CH}_3 \\   \\ \text{CH}_3\text{—C—CH}_2\text{—I} \\   \\ \text{H} \end{array} \quad \checkmark$	4	<b>ALLOW</b> correct structural <b>OR</b> displayed <b>OR</b> skeletal formula <b>OR</b> mixture of the above (as long as unambiguous) n.b. C <sub>2</sub> H <sub>5</sub> is unambiguous but C <sub>3</sub> H <sub>7</sub> is ambiguous  <b>IGNORE</b> incorrect name  Mark incorrect answers first of all. <ul style="list-style-type: none"> <li>• One incorrect answers maximum 3 marks</li> <li>• Two incorrect answers maximum 2 marks</li> <li>• Three incorrect answers maximum 1 mark</li> <li>• Four incorrect answers scores 0 mark</li> </ul> <b>ALLOW</b> as a slip one stick with no H on in a displayed formula

Question			Answer	Marks	Guidance
6	(b)	(ii)	$C_4H_{10}O$ ✓	1	<b>IGNORE</b> any structures drawn <b>DO NOT ALLOW</b> $C_4H_9OH$

Question			Answer	Marks	Guidance
6	(b)	(iii)	<p>infrared</p> <p>1700–1730 <math>\text{cm}^{-1}</math> indicates carbonyl group ✓</p> <p>broad 2900 <math>\text{cm}^{-1}</math> indicates O–H bond <b>AND</b> it is a <b>carboxylic acid</b> ✓</p> <p><b>explanation mark</b>  <b>B</b> has a branched structure because of relationship to methylpropene  <b>OR</b>  <b>C</b> has a branched structure because of relationship to methylpropene  <b>OR</b>  <b>C</b> must be a primary alcohol because it is oxidised to a carboxylic acid <b>OR</b> a primary alcohol because it reacts with acidified dichromate to make a carboxylic acid  <b>OR</b>  <b>C</b> cannot be a tertiary alcohol because it is oxidised <b>OR</b> cannot be a tertiary alcohol because it does react with acidified dichromate  ✓</p>	6	<p><b>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</b></p> <p><b>LOOK ON THE SPECTRUM</b> for labeled absorbances which can be given credit</p> <p><b>ALLOW</b> has a C=O bond because it has absorbance within range 1640–1750 <math>\text{cm}^{-1}</math></p> <p><b>ALLOW</b> 2900 <math>\text{cm}^{-1}</math> indicates O–H in carboxylic acid  <b>ALLOW</b> has O–H bond in carboxylic acid because it has absorbance within range 2500–3300 <math>\text{cm}^{-1}</math>  The presence of carboxylic acid can be anywhere in the text including the structure for <b>D</b></p> <p>If two marking points from the explanation mark are given both must be correct</p>

Question	Answer	Marks	Guidance
	<p><b>B</b> is <math display="block">\begin{array}{c} \text{CH}_3 \\   \\ \text{CH}_3 - \text{C} - \text{CH}_2 - \text{I} \\   \\ \text{H} \end{array}</math> ✓</p> <p><b>C</b> is <math display="block">\begin{array}{c} \text{CH}_3 \\   \\ \text{CH}_3 - \text{C} - \text{CH}_2 - \text{OH} \\   \\ \text{H} \end{array}</math> ✓</p> <p><b>D</b> is <math display="block">\begin{array}{c} \text{CH}_3 \\   \\ \text{CH}_3 - \text{C} - \text{COOH} \\   \\ \text{H} \end{array}</math> ✓</p>		<p><b>ALLOW</b> correct structural <b>OR</b> displayed <b>OR</b> skeletal formula <b>OR</b> mixture of the above (as long as unambiguous)</p> <p><b>IGNORE</b> incorrect names for <b>B</b>, <b>C</b> and <b>D</b></p> <p>Mark correct branched structures first of all.</p> <p>If there are no correct branched structures and <b>C</b> is <math>\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{OH}</math> then <b>ALLOW</b> one mark for <math>\text{CH}_3\text{CH}_2\text{CH}_2\text{COOH}</math> and one mark for <math>\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{I}</math></p>
	<b>Total</b>	<b>13</b>	

## APPENDIX 1





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