

# **Chemistry A**

Advanced Subsidiary GCE

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

## **Mark Scheme for January 2011**

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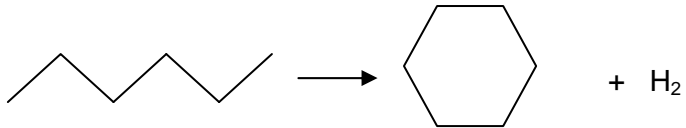
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Question		Answer	Mark	Guidance
1	(a)	<p>(The hydrocarbons have) different boiling points ✓</p> <p>The larger the molecules the stronger the van der Waals' forces ✓</p>	2	<p><b>PLEASE READ COMMENT ON PAGE 3</b></p> <p><b>ALLOW</b> longer chains have higher boiling points  <b>OR</b> separation based on boiling point  <b>OR</b> condense at different temperatures</p> <p><b>ALLOW</b> the larger molecular size more van der Waals' forces  <b>OR</b> longer chains have stronger van der Waals' force  <b>OR</b> the more electrons, the stronger the van der Waals' forces  <b>OR</b> the more surface contact the more van der Waals' forces  <b>IGNORE</b> surface area  <b>ALLOW ORA</b></p> <p>van der Waals must be seen at least once in correct context  <b>ALLOW</b> any 'recognisable' spelling of van der Waals', use of VDW is <b>not</b> sufficient</p> <p><b>DO NOT ALLOW</b> intermolecular force unless qualified as van der Waals' somewhere</p>
	(b)	(i) $C_nH_{2n}$ ✓	1	
		(ii) $C_6H_{14} \rightarrow C_6H_{12} + H_2$ ✓	1	<p><b>ALLOW</b> displayed, skeletal or structural formulae or combination in the equation</p> 

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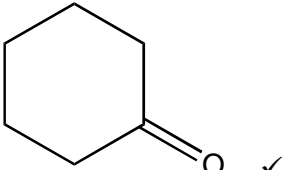
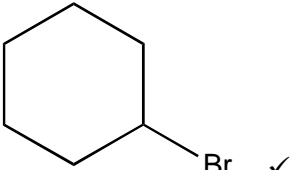
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1	(b)	(iii)	cyclohexane has more efficient combustion ✓	1	<p>Assume comments refer to cyclohexane unless specified otherwise</p> <p><b>ALLOW</b> cyclohexane allows smoother burning  <b>OR</b> cyclohexane increases octane number  <b>OR</b> cyclohexane reduces knocking  <b>OR</b> cyclohexane is less likely to produce pre-ignition  <b>OR</b> cyclohexane is a more efficient fuel  <b>OR</b> cyclohexane burns better <b>OR</b> easier to burn  <b>OR</b> cyclohexane combusts more easily  <b>OR</b> improves combustion  <b>DO NOT ALLOW</b> cyclohexane ignites more easily</p> <p><b>ALLOW</b> ORA for hexane</p> <p><b>IGNORE</b> cyclohexane increases volatility of fuel  <b>IGNORE</b> cyclohexane has a lower boiling point</p> <p>cyclohexane is a better fuel on its own is <b>NOT</b> sufficient  cyclohexane burns more cleanly on its own is <b>NOT</b> sufficient</p>
	(c)	(i)	<p><i>Unsaturated</i>: Contains (at least one) <b>carbon–carbon</b> double bond <b>OR</b> C=C <b>OR</b> multiple carbon–carbon bond ✓</p> <p><i>hydrocarbon</i>: Contains hydrogen and carbon <b>only</b> ✓</p>	2	<p><b>DO NOT ALLOW</b> just ‘contains a double bond’</p> <p><b>DO NOT ALLOW</b> ‘a mixture of carbon and hydrogen’  <b>OR</b> ‘contains carbon and hydrogen’  <b>OR</b> carbon and hydrogen molecules only</p>
		(ii)	<p>More than one hydrogen atom is substituted  <b>OR</b> ‘multisubstitution’ (by chlorine)  <b>OR</b> further substitution occurs ✓</p>	1	<p><b>ALLOW</b> can get dichloro-compounds (<b>IGNORE</b> numbering)  <b>ALLOW</b> reaction forms more than one <b>organic</b> product</p> <p><b>DO NOT ALLOW</b> ‘forms termination products’ on its own</p> <p>Reaction is not specific  <b>OR</b> reaction is difficult to control is <b>NOT</b> sufficient</p>

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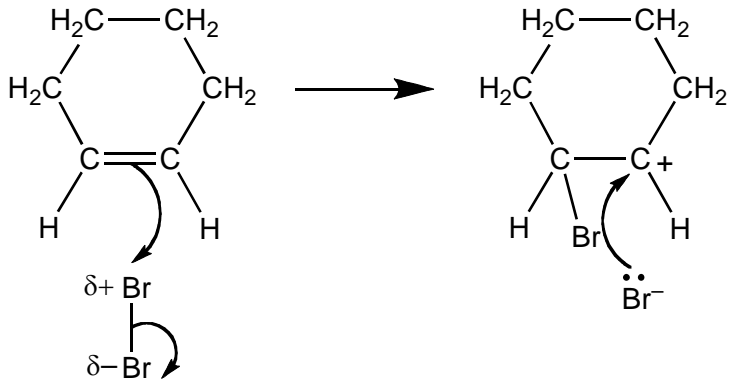
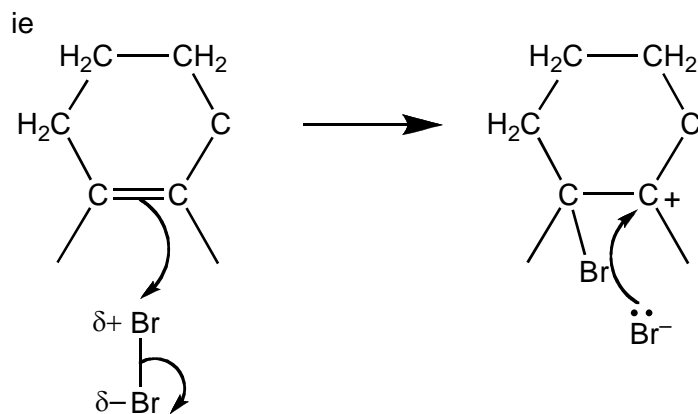
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Question			Answer	Mark	Guidance
1	(c)	(iii)	Contains a lone pair that can be donated ✓	1	<b>ALLOW</b> it can donate an electron pair 'lone pair' on its own is <b>NOT</b> sufficient
		(iv)	<p><b>A</b>  ✓</p> <p><b>B</b>  ✓</p>	2	<b>ALLOW</b> skeletal, displayed or structural formulae for <b>A</b> and <b>B</b> <b>ALLOW</b> combination of types of formulae as long as it is unambiguous  <b>DO NOT ALLOW</b> molecular formula For <b>A</b> , <b>ALLOW</b> carbonyl group on any carbon atom as it is still cyclohexanone  For <b>B</b> , <b>ALLOW</b> bromine atom on any carbon atom as it is still bromocyclohexane

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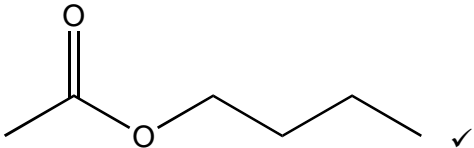
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Question	Answer	Mark	Guidance
1 (c) (v)	<p>Correct dipole on Br<sub>2</sub> / correct partial charges on Br<sub>2</sub> ✓</p> <p>Correct curly arrow from double bond to attack bromine atom and correct curly arrow to show heterolytic fission of Br–Br ✓</p> <p>Correct carbocation / carbonium ion drawn with the <b>full</b> positive charge shown: C<sup>+</sup> ✓</p> <p>Correct curly arrow from lone pair of Br<sup>–</sup> to correct carbon atom  <b>OR</b>  correct curly arrow from negative charge of Br<sup>–</sup> to correct carbon atom ✓</p> 	4	<p><b>ANNOTATE WITH TICKS AND CROSSES</b></p> <p>Curly arrow must come from covalent bonds and not atoms</p> <p><b>DO NOT ALLOW</b> C<sup>δ+</sup> for charge on carbonium ion</p> <p>Curly arrow from bromide ion can come from the negative charge or the lone pair  <b>DO NOT ALLOW</b> Br<sup>δ-</sup> instead of Br<sup>–</sup></p> <p>Lone pair does not need to be shown on Br<sup>–</sup> or used in mechanism</p> <p>Treat missing hydrogens on the CH<sub>2</sub> as a slip  Treat missing hydrogens on the double bond or carbonium ion as a slip providing a bond is shown</p> <p>ie</p>  <p><b>ALLOW</b> use of skeletal formulae in mechanism</p>
	<b>Total</b>	<b>15</b>	

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Question		Answer	Mark	Guidance
2	(a)		1	<b>IGNORE</b> any structural or displayed formula shown even if wrong (ie treat as rough working)
	(b)	<p>(<math>M_r</math> of all reactants <b>or</b> <math>M_r</math> of all products) is 134.0 <b>OR</b> 134  <b>OR</b>  (<math>M_r</math> of desired product) is 116.0 <b>OR</b> 116 ✓</p> <p>Atom economy = <math>100 \times \frac{116.0}{134.0}</math> ✓</p>	2	<p>Remember the marks are for the working out and not for the answer  <b>IGNORE</b> lack of decimal place in answer</p> <p><b>ALLOW</b> correct expressions to calculate the <math>M_r</math> or the atom economy eg</p> $\text{Atom economy} = 100 \times \frac{(6 \times 12) + (12 \times 1) + (2 \times 16)}{116 + 18}$ <p>Award 2 marks for this expression: <math>100 \times \frac{116.0}{134.0}</math> or similar expressions such as that above (subsumes 1st marking point)</p>
	(c) (i)	<p><b>acid</b> (catalyst) ✓</p> <p>heat <b>OR</b> reflux ✓</p>	2	<p><b>ALLOW</b> any acid, concentrated or dilute</p> <p><b>ALLOW</b> 'high temperature'  <b>OR</b> any temperature from 70 °C to 120 °C  Warm is <b>not</b> sufficient <b>but ALLOW</b> warm to 80 °C</p> <p><b>IGNORE</b> pressure</p>

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Question	Answer	Mark	Guidance
2 (c) (ii)	<p>maximum mass of ester than can be made is 9.7972973 (g) ✓</p> <p><math>\% \text{ yield} = \frac{6.57}{9.80} \times 100 \checkmark</math></p> <p><b>ALLOW</b> 2 or more sig figs up to calculated value but rounded up correctly, ie <b>ALLOW</b> <math>\frac{6.57}{9.797} \times 100</math> <b>OR</b> <math>\frac{6.57}{9.8} \times 100</math></p>	2	<p><b>ALLOW</b> moles of butan-1-ol = 0.08445946 <b>AND</b> moles of ester = 0.05663791</p> <p><b>OR</b> moles of butan-1-ol = <math>\frac{6.25}{74}</math> <b>AND</b> moles of ester = <math>\frac{6.57}{116}</math></p> <p>for one mark</p> <p><b>ALLOW</b> <math>\% \text{ yield} = \frac{0.05664}{0.08446} \times 100</math> for one mark</p> <p><b>ALLOW</b> 2 or more sig figs up to calculated value but rounded up correctly, ie <math>\frac{0.057}{0.084} \times 100</math> <b>OR</b> <math>\frac{0.0566}{0.0845} \times 100</math></p> <p>Remember the marks are for the working out</p>
(d)	<p>Link between yield <b>AND</b> explanation required:</p> <p>(high percentage) yield shows a high % conversion (of reactants into products) ✓</p> <p>Link between atom economy <b>AND</b> explanation required:</p> <p>(low) atom economy shows a <b>lot</b> of waste (product) <b>OR</b> (low) atom economy shows not much desired product ✓</p>	2	<p><b>ALLOW</b> percentage yield takes into account the practical difficulties of the process <b>OR</b> high % yield very little experimental loss of product <b>OR</b> high % yield because the process is not reversible <b>OR</b> most of reactants react to form products <b>DO NOT ALLOW</b> 'a lot of product made'</p> <p>There are waste products is <b>NOT</b> sufficient Reaction forms many products is <b>NOT</b> sufficient</p> <p><b>ALLOW</b> undesired product(s) as alternative for waste <b>IGNORE</b> a lot of by-products but ..... <b>ALLOW</b> a lot of <b>waste</b> by-products</p> <p><b>ALLOW</b> (low) atom economy shows a <b>lot</b> of HCl <b>OR</b> a lot of SO<sub>2</sub> is made <b>ALLOW</b> (low) atom economy shows not much ester / butyl ethanoate made</p>



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2 (e)	<p><b>NOTE: Comparison</b> essential throughout, ie <b>higher, less,</b> etc.</p> <p><b>ANY TWO FROM</b>            Less waste (products)  <b>OR</b> higher atom economy ✓</p> <p>Less toxic reactants  <b>OR</b> less toxic (waste) products  <b>OR</b> less corrosive reactants  <b>OR</b> less corrosive (waste) products  <b>OR</b> less harmful reactants  <b>OR</b> less harmful (waste) products  <b>OR</b> less hazardous reactants  <b>OR</b> less hazardous (waste) products ✓</p> <p>Cheaper starting materials  <b>OR</b> more readily available starting materials ✓</p> <p>Fewer steps  <b>OR</b> one step rather than two steps ✓</p>	2	<p><b>ALLOW</b> more sustainable</p> <p><b>ALLOW</b> poisonous for toxic</p> <p><b>IGNORE</b> 'dangerous'</p> <p>'Water is produced' is <b>not</b> sufficient</p> <p>Cheaper is <b>not</b> sufficient on its own</p> <p><b>IGNORE</b> less energy <b>OR</b> easier to carry out <b>OR</b> reversible</p>
	<b>Total</b>	<b>11</b>	

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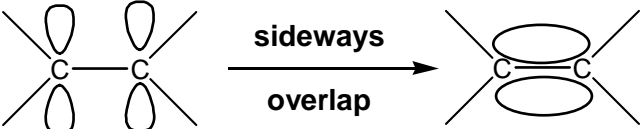
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Question		Answer	Mark	Guidance
3	(a)	(enthalpy change when) the number of moles of reactants ✓  as specified in the (balanced) equation react together ✓	2	<b>ALLOW</b> (enthalpy change when) the number of moles of products <b>ALLOW</b> molar quantities / amounts  Enthalpy change that occurs during a reaction is <b>not</b> sufficient
	(b) (i)	$Q = 50 \times 4.2 \times 11.0$ ✓  2.3 ✓	2	<b>ALLOW</b> 2310 J ✓ 2300j <b>ALLOW</b> use 4.18 for c which gives 2.299 J  <b>ALLOW</b> two marks for 2.31 / 2.310 with no working out <b>ALLOW ECF</b> ie Q divided by 1000 <b>IGNORE</b> any sign quoted
	(ii)	moles = 0.200 ✓	1	<b>ALLOW</b> 0.2 / 0.20
	(iii)	$\Delta H_r = 2 \times (2.3 \div 0.200)$ ✓  23 ✓  + sign ✓	3	<b>ALLOW ECF</b> from answer from $2 \times [(i) \div \text{answer to (ii)}]$  Answer from $2 \times [(i) \div \text{answer to (ii)}]$ must have only <b>2 sig figs</b>  + sign <b>must</b> be written for 'sign mark' + sign is <b>independent</b> of answer  <b>ALLOW</b> answers per mole of $\text{NH}_4\text{SCN}$ $\Delta H_r = 2.3 \div 0.200$ for one mark 12 for the second mark + sign for the third mark  <b>NOTE</b> If $c = 4.18$ has been used in <b>b(i)</b> , $\Delta H_r = +11$ by <b>ECF</b> for calculation per mole of $\text{NH}_4\text{SCN}$

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Question			Answer	Mark	Guidance
3	(c)	(i)	(Enthalpy change) when one mole of bonds ✓  of (gaseous covalent) bonds is broken ✓	2	<b>ALLOW</b> energy required rather than enthalpy change <b>DO NOT ALLOW</b> energy released  <b>DO NOT ALLOW</b> bonds formed
		(ii)	(Sideways) overlap of p orbitals ✓  Forming a $\pi$ /pi bond ✓	2	<b>IGNORE</b> reference to $\sigma$ bonds <b>IGNORE</b> incorrect diagram  This diagram would score one mark – the $\pi$ bond needs to be labelled for second mark   <b>2p orbitals</b>
		(iii)	$\pi$ bond is weaker (than the $\sigma$ bond) <b>OR</b> $\sigma$ bond is stronger (than the $\pi$ bond) ✓	1	There are two types of bonds is <b>not</b> sufficient <b>DO NOT ALLOW</b> $\pi$ bond is stronger than the $\sigma$ bond <b>ALLOW</b> the two bonds in double bond are not the same strength
		(iv)	bonds broken = (+)4010 <b>AND</b> bonds formed = (-)3931  Overall enthalpy change = +79 ✓	2	<b>ALLOW</b> Bonds broken = (+)690 <b>AND</b> bonds formed = (-)611 ✓  <b>ALLOW</b> 79 without a sign <b>ALLOW</b> -79 for one mark overall <b>ALLOW ECF</b> from incorrect enthalpy changes calculated for bonds broken and made

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Question			Answer	Mark	Guidance
3	(c)	(v)	Bond enthalpies may not be the same as the average bond enthalpy <b>OR</b> The idea that bonds have different strengths in different environments ✓	1	<b>DO NOT ALLOW</b> answers involving heat loss <b>OR</b> the use of non standard conditions  Average bond enthalpies are used is <b>NOT</b> sufficient
			<b>Total</b>	<b>16</b>	

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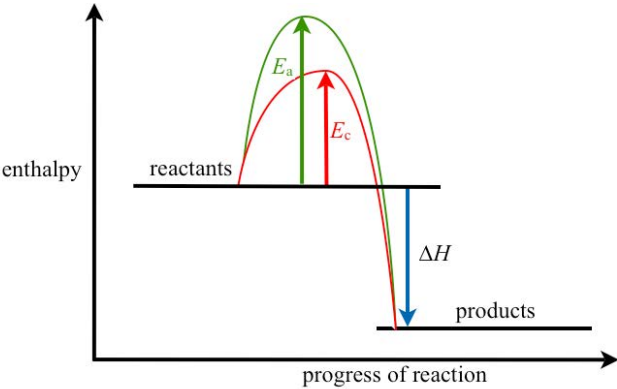
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4	(a)	(i)	$\text{Cl} + \text{O}_3 \rightarrow \text{ClO} + \text{O}_2 \checkmark$ $\text{ClO} + \text{O} \rightarrow \text{Cl} + \text{O}_2 \checkmark$	2	<b>ALLOW</b> any correct multiples <b>ALLOW</b> $\text{ClO} + \text{O}_3 \rightarrow 2\text{O}_2 + \text{Cl}$ <b>IGNORE</b> state symbols and dots
		(ii)	$\text{O}_3 + \text{O} \rightarrow 2\text{O}_2 \checkmark$	1	<b>ALLOW</b> any correct multiple <b>ALLOW</b> $2\text{O}_3 \rightarrow 3\text{O}_2$ <b>IGNORE</b> state symbols and dots
	(b)		Adsorption of reactants <b>OR</b> NO and CO attached to surface $\checkmark$  Bonds weaken in reactants $\checkmark$  Chemical reaction <b>OR</b> rearrangement of electrons $\checkmark$  Desorption $\checkmark$	4	<b>ANNOTATE WITH TICKS AND CROSSES</b>  <b>ALLOW</b> CO and NO (weakly) bonded to surface <b>OR</b> reactants bond to surface <b>OR</b> CO and NO form temporary bonds with the catalyst <b>DO NOT ALLOW</b> absorption  <b>ALLOW</b> bonds weaken in NO <b>OR</b> bonds weaken in CO <b>OR</b> activation energy is lowered  <b>ALLOW</b> bonds break and new bonds made in product <b>OR</b> $\text{N}_2$ and $\text{CO}_2$ made  <b>ALLOW</b> products leave the surface <b>OR</b> $\text{N}_2$ and $\text{CO}_2$ no longer bonded to surface <b>ALLOW</b> desorption <b>ALLOW</b> deabsorption if absorption given at start of answer

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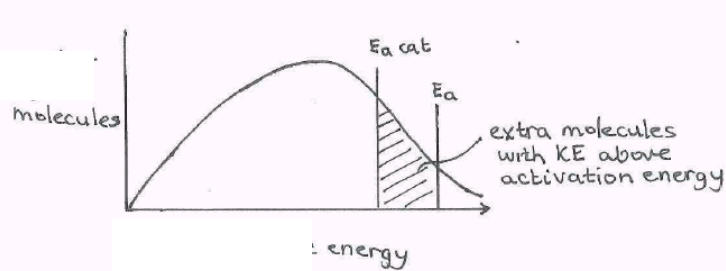
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4 (c)	<p>one activation energy labelled on enthalpy profile diagram ✓</p> <p>idea that activation energy is lowered ✓</p> <p>catalyst has a different reaction pathway  <b>OR</b> different reaction mechanism  <b>OR</b> two curves drawn on profile ✓</p> <p><b>QWC</b> – correct diagram of reaction profile for endothermic or exothermic reaction with products and reactants at different heights – y axis labelled as energy or enthalpy ✓</p>		<p><b>ANNOTATE WITH TICKS AND CROSSES</b></p> <p><b>ALLOW</b> double headed arrows on the activation energy label  <b>ALLOW</b> vertical line with no arrows  <b>DO NOT ALLOW</b> arrow just pointing downwards            Be generous with respect to the position of the line and the maximum of the curve</p> <p>marks can be awarded via, reaction profile, in words or from Boltzmann</p> <p><b>IGNORE</b> any enthalpy change label drawn</p>  <p><b>IGNORE</b> missing progress of reaction</p>

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4 (c)	<p>Drawing of Boltzmann distribution <b>AND</b> axes labelled (number of) molecules and energy ✓</p> <p>More molecules with energy above activation energy with a catalyst <b>OR</b>            More molecules that overcome the activation energy ✓</p> <p>More effective collisions <b>OR</b> more successful collisions ✓</p>	7	<p>Boltzmann distribution - must start at origin and must not end up at 0 on y-axis ie must not touch x-axis.  <b>DO NOT ALLOW</b> Boltzmann mark if two distributions are drawn one for non-catalysed and one for catalysed</p> <p><b>ALLOW</b> particles instead of molecules</p> <p><b>DO NOT ALLOW</b> atoms instead of particles</p>  <p><b>DO NOT ALLOW</b> more molecules have sufficient energy to react</p>

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4 (d)	<p><b>ANY FOUR FROM</b></p> <p>Enable reactions to occur with less waste  <b>OR</b> enable reactions to take place with higher atom economy  <b>OR</b> fewer undesired products ✓</p> <p>Enable reactions to happen with less toxic solvents/reactants  <b>OR</b> enable reactions to produce less toxic waste/side products ✓</p> <p>Reactions can happen at room temperature  <b>OR</b> reactions can happen at atmospheric pressure  <b>OR</b> reactions can happen at a lower pressure  <b>OR</b> reactions can happen at a lower temperature ✓</p> <p>Saves energy (costs) ✓</p> <p>Reduce carbon dioxide emissions  <b>OR</b> reduces amount of fuel burnt  <b>OR</b> reduces greenhouse gas emissions ✓</p> <p>Enable reactions to occur with more specificity  <b>OR</b> enable reactions to produce correct stereoisomer ✓</p>	4	<p><b>ANNOTATE WITH TICKS AND CROSSES</b></p> <p><b>ALLOW</b> make less hazardous waste  <b>ALLOW</b> corrosive, poisonous, harmful, hazardous as alternative to toxic  <b>DO NOT ALLOW</b> does not harm the environment  <b>IGNORE</b> dangerous</p> <p><b>IGNORE</b> less expensive  <b>IGNORE</b> reduces activation energy</p> <p><b>IGNORE</b> less pollution</p>
	<b>Total</b>	<b>18</b>	



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5 (a) (i)	$\text{CH}_3\text{CH}_2\text{I} + 2\text{NH}_3 \rightarrow \text{CH}_3\text{CH}_2\text{NH}_2 + \text{NH}_4\text{I}$ correct reactants ✓ correct products and balanced ✓	2	<b>ALLOW</b> $\text{CH}_3\text{CH}_2\text{I} + \text{NH}_3 \rightarrow \text{CH}_3\text{CH}_2\text{NH}_2 + \text{HI}$  <b>ALLOW</b> $\text{CH}_3\text{CH}_2\text{I} + \text{NH}_3 \rightarrow \text{CH}_3\text{CH}_2\text{NH}_3\text{I}$
	<div style="text-align: center;"> </div> <p>Correct curly arrow from the lone pair of ammonia to the carbon atom of C-Br ✓</p> <p>Correct dipole on C<sup>δ+</sup>-Br<sup>δ-</sup> bond <b>and</b> curly arrow showing the heterolytic fission of the C-Br bond ✓</p> <p>Correct missing product: Br<sup>-</sup> ✓</p>	3	<p>Curly arrow <b>must</b> start from the lone pair on nitrogen and go to the carbon atom  <b>DO NOT ALLOW</b> <math>\text{NH}_3^-</math> <b>OR</b> <math>^-\text{NH}_3</math>  <b>ALLOW</b> <math>\delta^-</math> on the N atom of <math>\text{NH}_3</math></p> <p>Curly arrow must start from the bond and go to the Br</p>

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5 (b)	<p><b>Effect of halogen in RX (3 marks)</b> Any correct comparison of rate <b>OR</b> reaction time between at least <b>TWO</b> of chloroalkane, bromoalkane and iodoalkane ✓</p> <p>Bond strength <b>OR</b> bond enthalpy/bond energy mentioned anywhere as a factor (even if reasoning is incorrect) ✓</p> <p>Any correct comparison of bond strength <b>OR</b> bond enthalpy/energy <b>OR</b> bond length <b>OR</b> ease of breaking of at least <b>TWO</b> of C–Cl, C–Br and C–I ✓</p>		<p><b>ANNOTATE WITH TICKS AND CROSSES</b> <b>Examples</b> chloroalkane reacts the slowest iodo compound reacts the fastest C–I bond is hydrolysed faster than C–Br C–Br has shorter reaction time than C–Cl</p> <p><b>DO NOT ALLOW</b> references to halogens as elements: <i>ie</i> chlorine is less reactive than bromine than iodine <b>DO NOT ALLOW</b> chloride, bromide and iodide</p> <p><b>ALLOW</b> this mark if mentioned within effect of halogen, branching <b>OR</b> temperature</p> <p><b>Examples</b> C–I bond is weaker than C–Br bond C–I bond is the weakest C–Cl bond is shorter than C–I bond C–Cl is strongest bond C–Br is broken more easily than C–Cl</p>

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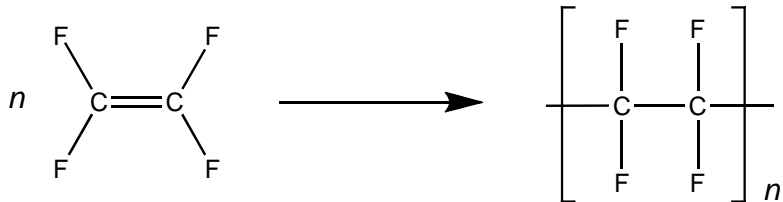
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Question	Answer	Mark	Guidance
5 (b)	<p><b>Effect of branching (2 marks)</b> Any correct comparison of rate or reaction time between at least <b>TWO</b> of the bromoalkanes ✓</p> <p>A sensible comparison of bond strength <b>OR</b> bond enthalpy/energy <b>OR</b> bond length <b>OR</b> ease of breaking of the C–Br bond in at least <b>TWO</b> of the bromoalkanes ✓</p> <p><b>Effect of temperature (2 marks)</b> <b>QWC</b> – Use of 50 °C and 60 °C using information in the table to show that rate increases with temperature ✓</p> <p>At higher temperature, particles have more energy <b>OR</b> At higher temperature, particles move faster ✓</p>	7	<p>Tertiary hydrolyses faster than secondary <b>OR</b> reaction time is less with tertiary than primary <b>OR</b> secondary hydrolyses faster than primary <b>OR</b> branched hydrolyses faster than straight chains <b>OR</b> primary hydrolyses the slowest <b>OR</b> tertiary hydrolyses the fastest <b>OR</b> when halogen on carbon 1 is hydrolysed slower than when halogen is on carbon 2 ✓</p> <p><b>DO NOT ALLOW</b> short chains hydrolyse faster than long chains</p> <p><b>Examples</b> C–Hal is weaker in tertiary halogenoalkane <b>OR</b> C–Br bond is stronger when it is bonded to carbon 1 rather than carbon 2</p> <p><b>ALLOW</b> an explanation based on relative stabilities of tertiary, secondary and/or primary carbocations</p> <p>Answer must <b>quote evidence</b> from the table to get this mark Rate increases with temperature is <b>NOT</b> sufficient</p> <p><b>ALLOW</b> more energy available to break the C–Hal bond <b>OR</b> more energy vibrates the C–Hal more so bond can break more easily <b>ALLOW</b> more successful collisions at higher temperature <b>ALLOW</b> more molecules exceed activation energy</p> <p><b>ALLOW ORA</b></p>

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Question	Answer	Mark	Guidance
5 (c) (i)	 <p data-bbox="331 454 593 486">Correct monomer ✓</p> <p data-bbox="331 622 571 654">Correct polymer ✓</p> <p data-bbox="331 821 1041 853">Balanced equation – correct use of <math>n</math> in the equation ✓</p>	3	<p data-bbox="1256 614 2045 750">Polymer must have <b>side</b> links (do not have to cut through bracket) <b>ALLOW a correct section of the polymer with side links</b> <b>ALLOW ECF</b> from wrong monomer, including use of FI for F</p> <p data-bbox="1256 821 2045 989"><math>n</math> on LHS can be at any height to the left of formula <b>AND</b> <math>n</math> on the RHS must be a subscript (essentially below the side link) On the LHS, <b>DO NOT ALLOW</b> <math>(C_2F_4)_n</math> (the <math>n</math> must be in front of the monomer)</p> <p data-bbox="1256 1021 2045 1053"><math>nC_2F_4 \rightarrow -(C_2F_4-)_n-</math> scores 1 mark for the correct use of <math>n</math></p>
	<p data-bbox="331 1098 784 1260">(PVC) produces hydrogen chloride <b>OR</b> produces acidic gases <b>OR</b> (PVC) produces phosgene <b>OR</b> produces toxic gases <b>OR</b> (PVC) produces dioxins ✓</p>	1	<p data-bbox="1256 1098 2045 1228"><b>ALLOW</b> produces poisonous gases <b>OR</b> produces gases that can kill <b>IGNORE</b> HF, Cl<sub>2</sub> and F<sub>2</sub> Makes a dangerous or harmful gas is <b>NOT</b> sufficient</p> <p data-bbox="1256 1260 2045 1364"><b>IGNORE</b> CO and CO<sub>2</sub> are greenhouse gases <b>IGNORE</b> chlorine radicals and ozone depletion <b>IGNORE</b> causes pollution</p>
<b>Total</b>		<b>16</b>	

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Question			Answer	Mark	Guidance
6	(a)	(i)	molecular ion is 58 <b>OR</b> $m/z$ is 58 ✓  (58 – (36 + 6) = 16) so $x = 1$ ✓	2	<b>ALLOW</b> peak on the right is 58 <b>OR</b> parent ion is 58 <b>ALLOW</b> 58 shown on the spectrum eg the peak is labelled with a number <b>OR</b> there is a ring around the peak  The $M_r$ <b>OR</b> molecular mass is 58 with no evidence is <b>not</b> sufficient  <b>ALLOW</b> $x = 1$ <b>ALLOW</b> Z is $C_3H_6O$
		(ii)	$CH_3CH_2CHO$ <b>OR</b> $CH_3COCH_3$ ✓	1	<b>ALLOW</b> displayed or skeletal formulae <b>ALLOW</b> combination of types of formulae as long as it is unambiguous  <b>ALLOW</b> other correct structures, eg enols, ethers and cyclic structures eg $CH_2=CHCH_2OH$ <b>OR</b> $CH_2=CHOCH_3$ <b>OR</b> structure of cyclopropanol  <b>DO NOT ALLOW</b> a structure showing H with 2 bonds, ie $OH-C$
		(iii)	$C_2H_5^+$ ✓	1	<b>ALLOW</b> $CH_3CH_2^+$ <b>OR</b> $COH^+$ <b>OR</b> $HCO^+$ The positive sign <b>must</b> be included
	(b)		$m/z$ values/peaks around 56 ✓	1	<b>ALLOW</b> peaks around 56 <b>OR</b> peak at 56 <b>OR</b> peaks around 55.8  <b>DO NOT ALLOW</b> peak at 55.8 <b>DO NOT ALLOW</b> peaks show the iron isotopes
	(c)	(i)	The <b>number</b> of $m/z$ values (around 32) ✓	1	<b>ALLOW</b> the <b>number</b> of peaks <b>IGNORE</b> any reference to molecular ion peak
		(ii)	Different isotopic abundance ✓	1	<b>ALLOW</b> different percentage of each isotope <b>OR</b> different isotopes present <b>ALLOW</b> sulfur atoms have different number of neutrons <b>OR</b> different mass numbers

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Question		Answer	Mark	Guidance
6	(d)	No absorption between 1640 and 1750 $\text{cm}^{-1}$ <b>AND</b> no (broad) absorption between 3200 and 3550 $\text{cm}^{-1}$ ✓	1	<b>ALLOW</b> the only significant absorption is at around 2850 to 3100 $\text{cm}^{-1}$ due to C–H bond <b>OR</b> There is an absorption around 2850 to 3100 $\text{cm}^{-1}$ due to C–H bond <b>AND no</b> absorptions by C=O and O–H bonds  <b>IGNORE</b> comments about C—O  <b>ALLOW</b> any values within the wavenumber range
	(e)	C=O because of absorption between 1640 and 1750 $\text{cm}^{-1}$  <b>AND</b> O–H (broad) absorption between 2500 to 3300 $\text{cm}^{-1}$ ✓  Carboxyl group <b>OR</b> carboxylic acid ✓	2	<b>ALLOW</b> any values within the wavenumber range <b>ALLOW</b> O–H (broad) absorption between 2500 to 3500 $\text{cm}^{-1}$ (from spectrum) <b>IGNORE</b> C–O  <b>ALLOW</b> carboxylic acid if linked with O–H absorption <b>IGNORE</b> alcohol, ester, aldehyde, ketone or amide
<b>Total</b>			<b>10</b>	

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Question		Answer	Mark	Guidance
7	(a)	<p><b>ANY THREE FROM</b></p> <p><math>C_6H_{12}O_6 \rightarrow 2CO_2 + 2C_2H_5OH</math> ✓</p> <p>Use of yeast/zymase at 25–45 °C <b>OR</b> warm with yeast/zymase ✓</p> <p>Anaerobic <b>OR</b> lack of oxygen ✓</p> <p>(Separate bioethanol) by (fractional) distillation ✓</p>	3	<p><b>IGNORE</b> state symbols</p> <p><b>ALLOW</b> correct multiples</p> <p><b>DO NOT ALLOW</b> yeast/zymase and heat <b>DO NOT ALLOW</b> yeast/zymase and reflux</p>
	(b) (i)	$C_{15}H_{30}O_2 + 21\frac{1}{2}O_2 \rightarrow 15CO_2 + 15H_2O$ ✓✓	2	<p><b>ALLOW</b> <math>\frac{43}{2}</math> for 21½</p> <p><b>DO NOT ALLOW</b> [O]</p> <p><b>ALLOW</b> one mark for correct products if equation is wrong</p>
	(ii)	(Energy needed) for processing biofuel makes carbon dioxide ✓	1	<b>ALLOW</b> (energy needed) for transport makes carbon dioxide
	(c)	<p><b>ANY THREE FROM</b></p> <p>Fossil fuels are finite resources <b>OR</b> biofuels are renewable ✓</p> <p>Allows fossil fuels to be used as a feedstock for organic compounds ✓</p> <p>Less food crops may be grown <b>OR</b> Land not used to grow food crops ✓</p> <p>(rain) forests have to be cut down to provide land <b>OR</b> deforestation ✓</p> <p>Shortage of fertile soils <b>OR</b> reduces fertility of soils ✓</p> <p>No risk of large scale pollution from exploitation of fossil fuels ✓</p>	3	<p><b>ANNOTATE WITH TICKS AND CROSSES</b></p> <p><b>ALLOW</b> fossil fuels are non-renewable <b>OR</b> plants are a renewable resource <b>OR</b> bio-fuels is (more) sustainable <b>OR</b> fossil fuels are not sustainable</p> <p><b>ALLOW</b> decrease the need for fossil fuels</p> <p>Destroys habitats is <b>NOT</b> sufficient</p> <p><b>IGNORE</b> comments about availability / fertilisers / pesticides</p>

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Question		Answer	Mark	Guidance
7	(d)	React with hydrogen <b>OR</b> hydrogenation ✓  Nickel catalyst ✓	2	<b>IGNORE</b> reference to pressure and temperature
	(e) (i)	Drawing of the <i>Z</i> isomer with the double bond shown in full ✓	1	Diagram must show a minimum of four carbon atoms and two hydrogen atoms and the correct orientation of the C=C double bond  <b>ALLOW</b> minor slips with rest of structure eg missing atoms, bonds and subscripts
	(ii)	Double bond does not rotate <b>OR</b> restricted rotation of the double bond ✓  Each carbon atom of double bond is bonded to (two) different groups ✓	2	<b>ALLOW</b> $\pi$ /pi bond does not rotate <b>IGNORE</b> 'bond does not move'  <b>ALLOW</b> each carbon atom of double bond is bonded to (two) different atoms <b>OR</b> each carbon atom of double bond is bonded to a hydrogen and a carbon/different group <b>OR</b> each end of the $\pi$ /pi-bond is bonded to different groups or atoms
<b>Total</b>			<b>12</b>	



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