



# **Chemistry A**

Advanced Subsidiary GCE

Unit F322: Chains, Energy and Resources

## Mark Scheme for January 2011

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Mark schemes should be read in conjunction with the published question papers and the Report on the Examination.

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Any enquiries about publications should be addressed to:

OCR Publications PO Box 5050 Annesley NOTTINGHAM NG15 0DL

Telephone:0870 770 6622Facsimile:01223 552610E-mail:publications@ocr.org.uk

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

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

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

Question		Answer	Mark	Guidance
	(v)	Correct dipole on $Br_2$ / correct partial charges on $Br_2 \checkmark$ Correct curly arrow from double bond to attack bromine	4	ANNOTATE WITH TICKS AND CROSSES Curly arrow must come from covalent bonds and not
		atom and correct curly arrow to show heterolytic fission of Br–Br $\checkmark$		atoms
		Correct carbocation / carbonium ion drawn with the full positive charge shown: $C^* \checkmark$		<b>DO NOT ALLOW</b> $C^{\delta+}$ for charge on carbonium ion
		Correct curly arrow from lone pair of Br <sup>-</sup> to correct carbon atom <b>OR</b>		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>
		correct curly arrow from negative charge of $Br^-$ to correct carbon atom $\checkmark$		Lone pair does not need to be shown on Br <sup>-</sup> or used in mechanism
		$H_2C - CH_2 $ $H_2C - CH_2$ $H_2C - CH_2 $ $H_2C - CH_2$ $H_2C - CH_2$		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 ie
		H = H = H = H = H		$\begin{array}{c c} H_2C \longrightarrow CH_2 & H_2C \longrightarrow CH_2 \\ H_2C & C & \longrightarrow H_2C & C \end{array}$
		$\delta + Br$ $Br$ $Br$		
				$\delta + Br$ $Br$ $Br$
				ALLOW use of skeletal formulae in mechanism
		Total	15	

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

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

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

Q	Question		Answer	Mark	Guidance
3	(a)		(enthalpy change when) the number of moles of reactants ✓		ALLOW (enthalpy change when) the number of moles of products ALLOW molar quantities / amounts
			as specified in the (balanced) equation react together $\checkmark$	2	Enthalpy change that occurs during a reaction is <b>not</b> sufficient
	(b)	(i)	Q = 50 × 4.2 × 11.0 ✓		<b>ALLOW</b> 2310 J ✓ 2300j <b>ALLOW</b> use 4.18 for <i>c</i> which gives 2.299 J
			2.3 ✓	2	ALLOW two marks for 2.31 / 2.310 with no working out ALLOW ECF ie Q divided by 1000 IGNORE any sign quoted
		(ii)	moles = 0.200 ✓	1	ALLOW 0.2 / 0.20
		(iii)	$\Delta H_{\rm r} = 2 \times (2.3 \div 0.200) \checkmark$		ALLOW ECF from answer from 2 × [(i) ÷ answer to (ii)]
			23 ✓		Answer from 2 × [(i) ÷ answer to (ii)] must have only 2 sig figs
			+ sign ✓	3	<ul> <li>+ sign must be written for 'sign mark'</li> <li>+ sign is independent of answer</li> </ul>
					<b>ALLOW</b> answers per mole of NH <sub>4</sub> 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 NH <sub>4</sub> SCN

G	Question		Answer	Mark	Guidance
3	(c)	(i)	(Enthalpy change) when one mole of bonds ✓ of (gaseous covalent) bonds is broken ✓	2	<ul> <li>ALLOW energy required rather than enthalpy change</li> <li>DO NOT ALLOW energy released</li> <li>DO NOT ALLOW bonds formed</li> </ul>
		(ii)	(Sideways) overlap of p orbitals ✓ Forming a π/pi bond ✓	2	IGNORE reference to $\sigma$ bonds IGNORE incorrect diagram This diagram would score one mark – the $\pi$ bond needs to be labelled for second mark $\overbrace{\c}$ $\overbrace{\c}$
		(iii)	π bond is weaker (than the σ bond) <b>OR</b> σ bond is stronger (than the π 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	<ul> <li>ALLOW Bonds broken = (+)690</li> <li>AND bonds formed = (−)611 ✓</li> <li>ALLOW 79 without a sign</li> <li>ALLOW –79 for one mark overall</li> <li>ALLOW ECF from incorrect enthalpy changes calculated for bonds broken and made</li> </ul>

Q	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
			Total	16	

Qı	Question		Answer	Mark	Guidance
4	(a)	(i)	$CI + O_3 \rightarrow CIO + O_2 \checkmark$		ALLOW any correct multiples
			$CIO + O \rightarrow CI + O_2 \checkmark$	2	ALLOW CIO + $O_3 \rightarrow 2O_2$ + CI
					IGNORE state symbols and dots
		(ii)	$O_3 + O \rightarrow 2O_2 \checkmark$	1	ALLOW any correct multiple
					ALLOW $2O_3 \rightarrow 3O_2$
					IGNORE state symbols and dots
	(b)				ANNOTATE WITH TICKS AND CROSSES
			Adsorption of reactants OR NO and CO attached to surface ✓ Bonds weaken in reactants ✓ Chemical reaction OR rearrangement of electrons ✓ Desorption ✓	4	<ul> <li>ALLOW CO and NO (weakly) bonded to surface</li> <li>OR reactants bond to surface</li> <li>OR CO and NO form temporary bonds with the catalyst</li> <li>DO NOT ALLOW absorption</li> <li>ALLOW bonds weaken in NO</li> <li>OR activation energy is lowered</li> <li>ALLOW bonds break and new bonds made in product</li> <li>OR N<sub>2</sub> and CO<sub>2</sub> made</li> <li>ALLOW products leave the surface</li> <li>OR N<sub>2</sub> and CO<sub>2</sub> no longer bonded to surface</li> <li>ALLOW deadsorption</li> <li>ALLOW deabsorption if absorption given at start of answer</li> </ul>

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

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

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

Q	uest	ion	Answer	Mark	Guidance
5	(a)	(i)	CH <sub>3</sub> CH <sub>2</sub> I + 2NH <sub>3</sub> → CH <sub>3</sub> CH <sub>2</sub> NH <sub>2</sub> + NH <sub>4</sub> I correct reactants $\checkmark$ correct products and balanced $\checkmark$	2	ALLOW $CH_3CH_2I + NH_3$ $\rightarrow CH_3CH_2NH_2 + HI$ ALLOW $CH_3CH_2I + NH_3 \rightarrow CH_3CH_2NH_3I$
		(ii)	$CH_{3}CH_{2} \xrightarrow{H}_{G} \xrightarrow{h}_{H} \xrightarrow{h}_{Br} \xrightarrow{h}_{CH_{3}} CH_{3}CH_{2} \xrightarrow{H}_{H} \xrightarrow{h}_{H} \xrightarrow{h}_{H} = Br^{-}$ $H^{-}_{H} \xrightarrow{h}_{H} $		Curly arrow <b>must</b> start from the lone pair on nitrogen and go to the carbon atom <b>DO NOT ALLOW</b> NH <sub>3</sub> <sup>-</sup> <b>OR</b> <sup>-</sup> NH <sub>3</sub> <b>ALLOW</b> δ– on the N atom of NH <sub>3</sub>
			Correct dipole on $C^{\delta_+}$ –Br <sup><math>\delta</math></sup> bond <b>and</b> curly arrow showing the heterolytic fission of the C–Br bond $\checkmark$		Curly arrow must start from the bond and go to the Br
			Correct missing product: Br <sup>-</sup> ✓	3	

Qı	uestio	on	Answer	Mark	Guidance
5	(b)		<i>Effect of halogen in RX (3 marks)</i> Any correct comparison of rate <b>OR</b> reaction time between at least <b>TWO</b> of chloroalkane, bromoalkane and iodoalkane ✓		ANNOTATE WITH TICKS AND CROSSES <i>Examples</i> 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–CI <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
			Bond strength <b>OR</b> bond enthalpy/bond energy mentioned anywhere as a factor (even if reasoning is incorrect) $\checkmark$		<b>ALLOW</b> this mark if mentioned within effect of halogen, branching <b>OR</b> temperature
			Any correct comparison of bond strength OR bond enthalpy/energy OR bond length OR ease of breaking of at least <b>TWO</b> of C–CI, C–Br and C–I ✓		Examples C-I bond is weaker than C-Br bond C-I bond is the weakest C-CI bond is shorter than C-I bond C-CI is strongest bond C-Br is broken more easily than C-CI

F322

## Mark Scheme

## January 2011

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

Q	uesti	ion	Answer	Mark	Guidance
Q 5	(c)	on (i)	Answer $ \begin{array}{c}                                     $	<u>Mark</u>	Guidance         Polymer must have side links         (do not have to cut through bracket)         ALLOW a correct section of the polymer with side links         ALLOW ECF from wrong monomer, including use of FI for F $n$ on LHS can be at any height to the left of formula         AND $n$ on the RHS must be a subscript         (essentially below the side link)         On the LHS, DO NOT ALLOW (C <sub>2</sub> F <sub>4</sub> ) <sub>n</sub> (the $n$ must be in front of the monomer)
		(ii)	<ul> <li>(PVC) produces hydrogen chloride</li> <li>OR produces acidic gases</li> <li>OR (PVC) produces phosgene</li> <li>OR produces toxic gases</li> <li>OR (PVC) produces dioxins ✓</li> </ul>	1	$nC_2F_4 \rightarrow -(-C_2F_4-)_n$ - scores 1 mark for the correct use of $n$ <b>ALLOW</b> produces poisonous gases <b>OR</b> produces gases that can kill <b>IGNORE</b> HF, $Cl_2$ and $F_2$ Makes a dangerous or harmful gas is <b>NOT</b> sufficient <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
			Total	16	

Q	uesti	ion	Answer	Mark	Guidance
6	(a)	(i)	molecular ion is 58 <b>OR</b> <i>m</i> / <i>z</i> is 58 ✓		<ul> <li>ALLOW peak on the right is 58 OR parent ion is 58</li> <li>ALLOW 58 shown on the spectrum eg the peak is labelled with a number</li> <li>OR there is a ring around the peak</li> <li>The <i>M</i><sub>r</sub> OR molecular mass is 58 with no evidence is <b>not</b> sufficient</li> </ul>
			$(58 - (36 + 6) = 16)$ so $x = 1 \checkmark$	2	ALLOW $x = 1$ ALLOW Z is C <sub>3</sub> H <sub>6</sub> O
		(ii)	CH <sub>3</sub> CH <sub>2</sub> CHO <b>OR</b> CH <sub>3</sub> COCH <sub>3</sub> ✓	1	<ul> <li>ALLOW displayed or skeletal formulae</li> <li>ALLOW combination of types of formulae as long as it is unambiguous</li> <li>ALLOW other correct structures, eg enols, ethers and cyclic structures eg CH<sub>2</sub>=CHCH<sub>2</sub>OH OR CH<sub>2</sub>=CHOCH<sub>3</sub> OR structure of cyclopropanol</li> <li>DO NOT ALLOW a structure showing H with 2 bonds, ie OH—C</li> </ul>
		(iii)	C <sub>2</sub> H <sub>5</sub> <sup>+</sup> ✓	1	ALLOW CH <sub>3</sub> CH <sub>2</sub> <sup>+</sup> OR COH <sup>+</sup> OR HCO <sup>+</sup> The positive sign <b>must</b> be included
	(b)		m/z values/peaks around 56 ✓	1	ALLOW peaks around 56 OR peak at 56 OR peaks around 55.8 DO NOT ALLOW peak at 55.8 DO NOT ALLOW peaks show the iron isotopes
	(c)	(i)	The <b>number</b> of $m/z$ values (around 32) $\checkmark$	1	ALLOW the number of peaks IGNORE any reference to molecular ion peak
		(ii)	Different isotopic abundance ✓	1	ALLOW different percentage of each isotope OR different isotopes present ALLOW sulfur atoms have different number of neutrons OR different mass numbers

### Mark Scheme

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

Qı	uesti	on	Answer	Mark	Guidance	
7	(a)		ANY THREE FROM		IGNORE state symbols	
			$C_6H_{12}O_6 \rightarrow 2CO_2 + 2C_2H_5OH \checkmark$		ALLOW correct multiples	
			Use of yeast/zymase at 25–45 ⁰C OR warm with yeast/zymase ✓		DO NOT ALLOW yeast/zymase and heat DO NOT ALLOW yeast/zymase and reflux	
			Anaerobic <b>OR</b> lack of oxygen ✓	3		
			(Separate bioethanol) by (fractional) distillation ✓			
	(b)	(i)	$C_{15}H_{30}O_2 + 21\frac{1}{2}O_2 \rightarrow 15CO_2 + 15H_2O \checkmark \checkmark$	2	<b>ALLOW</b> $\frac{43}{2}$ for 21 <sup>1</sup> / <sub>2</sub>	
					<b>DO NOT ALLOW</b> [O] <b>ALLOW</b> one mark for correct products if equation is wrong	
		(ii)	(Energy needed) for processing biofuel makes carbon dioxide ✓	1	ALLOW (energy needed) for transport makes carbon dioxide	
	(c)		ANY THREE FROM Fossil fuels are finite resources OR biofuels are renewable ✓		ANNOTATE WITH TICKS AND CROSSES ALLOW fossil fuels are non-renewable OR plants are a renewable resource OR bio-fuels is (more) sustainable OR fossil fuels are not sustainable	
			Allows fossil fuels to be used as a feedstock for organic compounds $\checkmark$		ALLOW decrease the need for fossil fuels	
			Less food crops may be grown OR Land not used to grow food crops ✓			
			(rain) forests have to be cut down to provide land <b>OR</b> deforestation ✓		Destroys habitats is <b>NOT</b> sufficient	
			Shortage of fertile soils OR reduces fertility of soils ✓		IGNORE comments about availability / fertilisers / pesticides	
			No risk of large scale pollution from exploitation of fossil fuels $\checkmark$	3		

Q	uesti	on	Answer	Mark	Guidance
7	(d)		React with hydrogen <b>OR</b> hydrogenation ✓		
			Nickel catalyst ✓	2	IGNORE reference to pressure and temperature
	(e)	(i)	Drawing of the Z 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 OR restricted rotation of the double bond ✓ Each carbon atom of double bond is bonded to (two) different groups ✓	2	<ul> <li>ALLOW π/pi bond does not rotate IGNORE 'bond does not move'</li> <li>ALLOW each carbon atom of double bond is bonded to (two) different atoms</li> <li>OR each carbon atom of double bond is bonded to a hydrogen and a carbon/different group</li> <li>OR each end of the π/pi-bond is bonded to different groups or atoms</li> </ul>
			Total	12	

OCR (Oxford Cambridge and RSA Examinations) 1 Hills Road Cambridge CB1 2EU

**OCR Customer Contact Centre** 

#### 14 – 19 Qualifications (General)

Telephone: 01223 553998 Facsimile: 01223 552627 Email: general.qualifications@ocr.org.uk

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