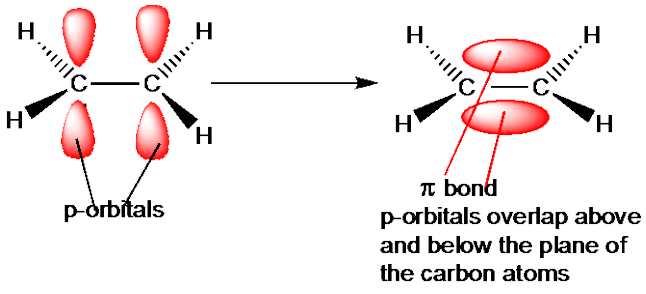
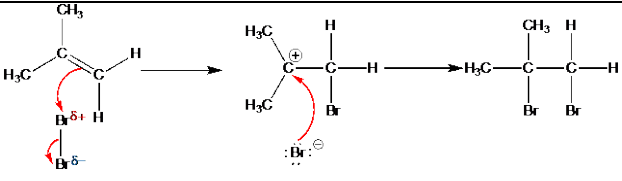
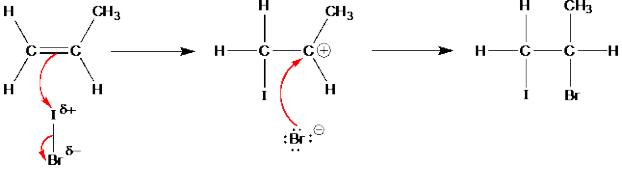
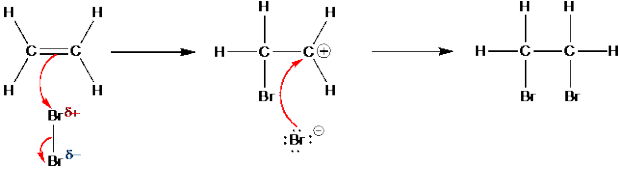
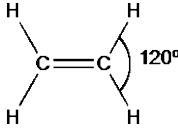
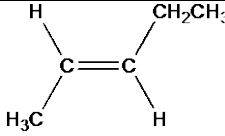


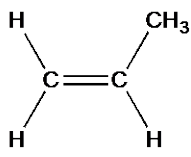
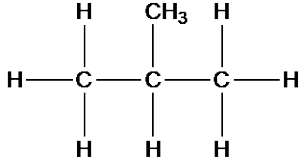
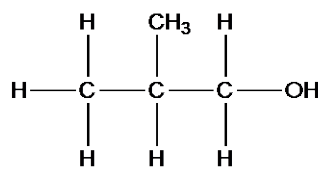
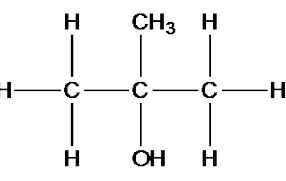
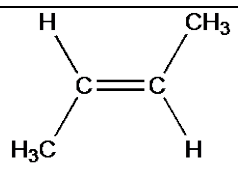
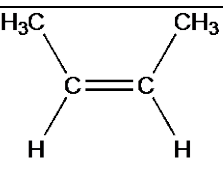
Question number	Answer	Marks	Guidance
1 (a)	An unsaturated compound contains double or triple covalent bonds. A hydrocarbon is a compound of carbon and hydrogen only	B1 B1	
1 (b)	a: 109.5° b: 120°	B1 B1	
1 (c)	 <p>1 mark for sideways overlap of p orbitals 1 mark for diagram showing π bond</p>	B1 x 2	
2 (a)	An electrophile is an electron pair acceptor	B1	
2 (b)	 <p>1 mark for curly arrow from C=C to $\text{Br}^{\delta+}$ of Br_2 1 mark for curly arrow from Br-Br and correct dipole 1 mark for correct carbocation AND curly arrow from Br^- to C^+ 1 mark for product</p>	B1 x 4	
3 (a) (i)	 <p>1 mark for curly arrow from C=C to $\text{I}^{\delta+}$ of IBr 1 mark for curly arrow from I-Br and correct dipole 1 mark for correct carbocation AND curly arrow from Br^- to C^+</p>	B1 x 3	<p>Curly arrow must start from bond and go to correct atom DO NOT ALLOW partial charges on carbon-carbon double bond</p> <p>DO NOT ALLOW $\delta+$ on carbon atom The positive charge must be associated with the carbon atom and not with a bond Make certain the carbonium ion includes the iodine atom</p>

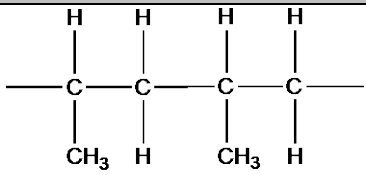
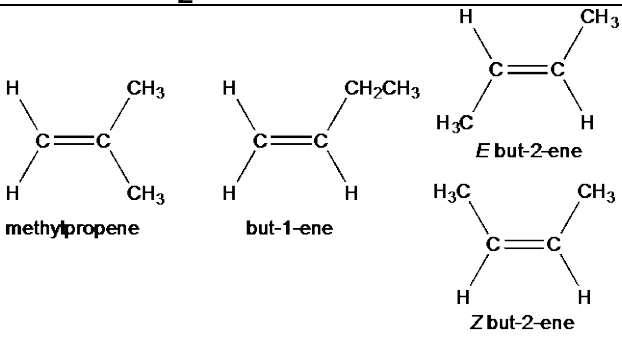
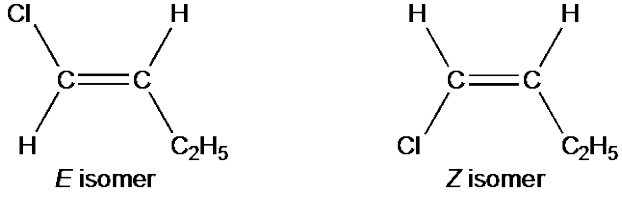
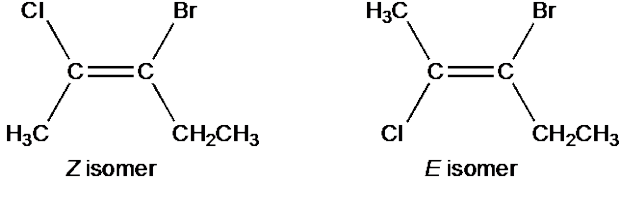
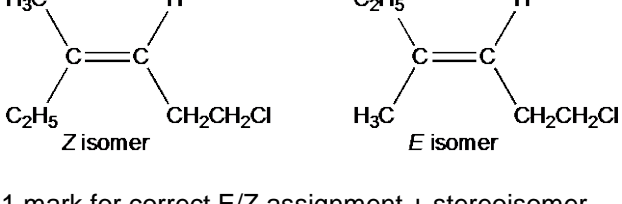
Question number	Answer	Marks	Guidance
			<p>Curly arrow must come from any lone pair or the negative sign of the bromide ion</p> <p>The lone pair on the bromide ion does not need to be shown</p>
3 (a) (ii)	Electrophilic addition	B1	
3 (a) (iii)	$ \begin{array}{c} \text{H} \quad \text{CH}_3 \\ \quad \\ \text{H}-\text{C}-\text{C}-\text{H} \\ \quad \\ \text{Br} \quad \text{I} \end{array} $	B1	<p>ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) eg $\text{CH}_2\text{BrCHICH}_3$</p> <p>IGNORE any name given</p>
4 (a)	<p>1,2-dibromoethane: Use bromine</p> <p>$\text{C}_2\text{H}_4 + \text{Br}_2 \rightarrow \text{C}_2\text{H}_4\text{Br}_2$ bromoethane: Hydrogen bromide</p> <p>$\text{C}_2\text{H}_4 + \text{Br}_2 \rightarrow \text{C}_2\text{H}_5\text{OH}$ Ethanol: steam</p> <p>with an acid catalyst (e.g. conc H_2SO_4) $\text{C}_2\text{H}_4 + \text{H}_2\text{O} \rightarrow \text{C}_2\text{H}_5\text{OH}$</p> <p>Name of mechanism: electrophilic addition</p>  <p>1 mark for curly arrow from $\text{C}=\text{C}$ to $\text{Br}^{\delta+}$ of Br_2</p> <p>1 mark for curly arrow from $\text{Br}-\text{Br}$ and correct dipole</p> <p>1 mark for correct carbocation</p> <p>1 mark for curly arrow from Br^- to C^+</p>	<p>B1</p> <p>B1</p> <p>B1</p> <p>B1</p> <p>B1</p> <p>B1</p> <p>B1 x 4</p>	<p>ALLOW reactants even from incorrect equations</p> <p>ALLOW reactants or conditions over the arrow</p> <p>ALLOW Br_2 mark from the mechanism even if the mechanism is incorrect</p> <p>IGNORE conditions unless they would lead to a different reaction with ethane</p> <p>IGNORE conditions unless they would lead to a different reaction with ethane</p> <p>ALLOW temperature range between $100\text{--}400^\circ\text{C}$ if quoted</p> <p>IGNORE reference to pressure</p> <p>IGNORE hydrolysis</p> <p>Hydration is not sufficient but DO NOT ALLOW Hydrogenation</p> <p>ALLOW H_2SO_4 OR H_3PO_4 OR H^+</p> <p>DO NOT ALLOW HCl, HBr etc.</p> <p>ALLOW two stage process e.g. react with HBr one mark followed by $\text{KOH}(\text{aq})$ one mark</p>

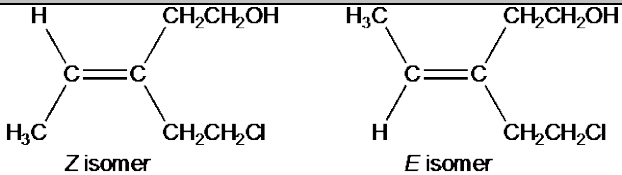
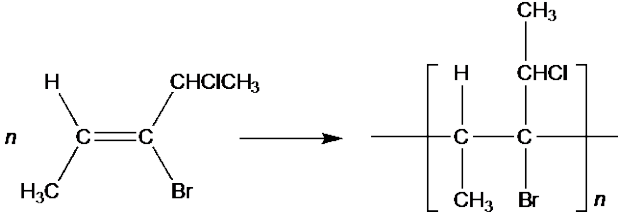
Question number	Answer	Marks	Guidance
			<p>Curly arrow must start from the double bond and not a carbon atom and go to the $\text{Br}^{\delta+}$; other curly arrow must start from $\text{Br}-\text{Br}$ bond.</p> <p>ALLOW attack of $\text{Br}-\text{Br}$ if dipoles not shown</p> <p>DO NOT ALLOW attack of $\text{Br}^{\delta-}$</p> <p>Dipole must be partial charge and not full charge</p> <p>DO NOT ALLOW 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>Br^- curly arrow must come from one lone pair on Br^- ion OR from minus sign on Br^- ion Lone pair does not need to be shown on Br^- ion</p> <p>ALLOW mechanism which goes via a cyclic bromonium ion instead of the carbocation</p>
4 (b)	 <p>3 areas of electron density surround each carbon atom and repel one another as far apart as possible</p> <p>Bond angle: 120°</p>	B1 B1 B1	<p>IGNORE any name of shape given</p> <p>ALLOW $115-125^\circ$</p> <p>ALLOW even if it is the $\text{C}-\text{C}-\text{H}$ shown on a diagram.</p> <p>ALLOW three or four electron pairs repel OR three or four bonds repel</p> <p>IGNORE does not have any lone pairs</p> <p>DO NOT ALLOW atoms repel / electrons repel</p> <p>DO NOT ALLOW has lone pair which repels more</p>
4 (c) (i)		B1	<p>ALLOW correct structural OR displayed OR skeletal formula</p> <p>OR mixture of the above (as long as unambiguous)</p>

Question number	Answer	Marks	Guidance
4 (c) (ii)		B1	<p>ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous)</p> <p>ALLOW CH₃ and C₂H₅ groups above or below chain</p> <p>ALLOW bond to ethyl and methyl group to any part of ethyl or methyl group</p> <p>IGNORE any brackets drawn</p> <p>ALLOW two or more repeat units but has to have a whole number of repeat units (ie does not have to be two)</p> <p>'End bonds' MUST be shown and can be dotted</p> <p>IGNORE <i>n</i></p>
5 (a)	<p style="text-align: center;">mixture of isomers</p> <p>1 mark for each product</p>	B1 x 4	<p>ALLOW skeletal formula OR displayed formulae</p> <p>IGNORE molecular formulae</p> <p>IF two answers given e.g. name and structure then both must be correct to be given a mark</p> <p>ALLOW methylpropane OR (CH₃)₃CH</p> <p>ALLOW 1,2-dibromo-methylpropane OR CH₂BrCBr(CH₃)₂</p> <p>ALLOW 1-bromo-methylpropane OR CH₂BrCH(CH₃)₂</p> <p>ALLOW 2-bromo-methylpropane OR CH₃CBr(CH₃)₂</p> <p>ALLOW ecf if wrong carbon skeleton is used in all of the structures mark first structure wrong and then apply ecf for the rest</p>

Question number	Answer	Marks	Guidance
5 (b)	<p>1 mark for curly arrow from C=C to Cl^{δ+} of Cl₂</p> <p>1 mark for curly arrow from Cl—Cl and correct dipole</p> <p>1 mark for correct carbocation AND curly arrow from Cl[−] to C⁺</p>	B1 x 3	<p>Curly arrow must start from the double bond and not a carbon atom, other curly arrow must start from Cl—Cl bond</p> <p>ALLOW curly arrow from any part of chloride ion The chloride ion does not need to show a lone pair</p> <p>Dipole must be partial charge and not full charge Carbocation needs a full charge and not a partial charge (charges do not need to be surrounded by a circle)</p> <p>ALLOW carbocation on carbon 1 where electrophile attacks carbon 2 i.e. ⁺CH₂CCl(CH₃)₂</p>
6 (a)	<p>Structural isomers are compounds with the same molecular formula</p> <p>but different structural formulae</p> <p>Stereoisomers have same structural formulae</p> <p>but different arrangements in space</p> <p>The double C=C bond which does not rotate</p> <p>The alkene needs to have two different groups attached to each carbon atom of C=C bond</p> <p>Evidence for molecular formula: F, G and H have molecular formula of C₅H₁₀ as <i>M_r</i> is 5 × 12 + 10 × 1 = 70</p> <p>1 mark for each structure of F, G and H (3 marks total)</p> <p>E and Z isomers identified</p>	<p>B1</p> <p>B1</p> <p>B1</p> <p>B1</p> <p>B1</p> <p>B1</p> <p>B1</p> <p>B1 x 4</p>	<p>ALLOW same molecular formula but different structures</p> <p>Second marking point is DEPENDENT on first mark</p> <p>ALLOW compounds with the same structure</p> <p>Second marking point is DEPENDENT on first mark</p> <p>This is the QWC mark</p> <p>IGNORE wrong names of F, G and H</p> <p>ALLOW structural or displayed formulae for F, G and H e.g. H is CH₃CH₂CH₂CHCH₂</p> <p>ALLOW identification using <i>trans</i> and <i>cis</i> and ALLOW this marking point as identification of another example of identifying <i>E/Z</i> or <i>cis</i> and <i>trans</i> if not done for F and G</p> <p>ALLOW one mark if no structures drawn but correct names given for F, G and H</p>

Question number	Answer	Marks	Guidance
			i.e <i>E</i> -pent-2-ene, <i>Z</i> -pent-2-ene and pent-1-ene ALLOW ecf on structures if wrong molecular formula used or consistent error or slip such as having just sticks
7	<p>$C : H = 82.8/12.0 : 17.2/1 = 6.9 : 17.2$</p> <p>Empirical formula = C_2H_5</p> <p>Molecular formula = $C_2H_5 \times 58/29$ $= C_2H_5 \times 2 = C_4H_{10}$</p> <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">  <p>A</p> </div> <div style="text-align: center;">  <p>B</p> </div> </div> <div style="display: flex; justify-content: space-around; align-items: center; margin-top: 20px;"> <div style="text-align: center;">  <p>C</p> </div> <div style="text-align: center;">  <p>D</p> </div> </div> <p>1 mark for each structure</p> <p>B: $C_4H_8 + H_2 \rightarrow C_4H_{10}$</p> <p>C and D: $C_4H_8 + H_2O \rightarrow C_4H_9OH$</p>	<p>B1</p> <p>B1</p> <p>B1</p> <p>B1 x 4</p> <p>B1</p> <p>B1</p>	
8	<div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">  <p><i>E</i> isomer</p> </div> <div style="text-align: center;">  <p><i>Z</i> isomer</p> </div> </div> <p>Q and R:</p> <p>2 marks for Q and R identified 1 mark for correct labelling of <i>E</i> and <i>Z</i> isomers</p>	B1 x 3	
9 (a)	Addition polymerisation	B1	
9 (b)	Propene	B1	

Question number	Answer	Marks	Guidance
9 (c)	 <p style="text-align: center;">E</p>	B1	
10	 <p>1 mark for each structure</p> <p>Methylpropene, but-1-ene and but-2-ene are structural isomers</p> <p>But-2-ene has <i>E</i> and <i>Z</i> stereoisomers</p>	B1 x 4 B1 B1	
11 (a)	 <p>1 mark for correct <i>E/Z</i> assignment + stereoisomer</p>	B1	
11 (b)	 <p>1 mark for correct <i>E/Z</i> assignment + stereoisomer</p>	B1	
11 (c)	 <p>1 mark for correct <i>E/Z</i> assignment + stereoisomer</p>	B1	

Question number	Answer	Marks	Guidance
11 (d)	 <p style="text-align: center;">Z isomer E isomer</p> <p>1 mark for correct E/Z assignment + stereoisomer</p>	B1	
12	 <p>1 mark for repeat unit being correct</p> <p>1 mark for monomer and balancing</p>	B1 x 2	