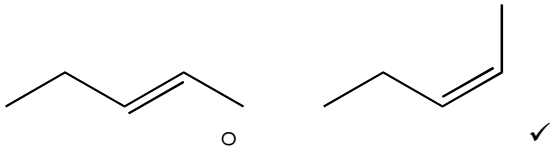
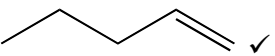
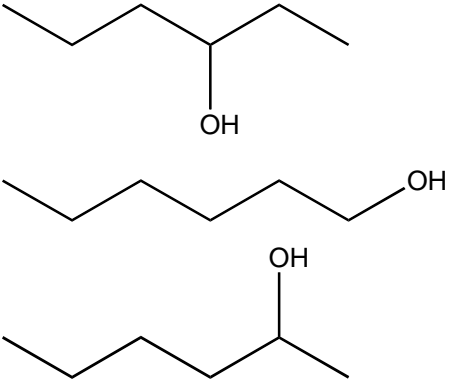


| Question | Expected Answers | Marks | Additional Guidance |
|----------|---|-------|--|
| 1 (a) | <p>Structural isomer compounds with the same molecular formula ✓ but with different structural formulae ✓</p> <p>Stereoisomer compounds with the same structural formula ✓ but with different arrangements in space ✓</p> <p>Evidence of using M_r of 70 to calculate molecular formula of C_5H_{10} ✓</p> <p>F and G are</p>  <p>Correct identification of the <i>E</i> and <i>Z</i> isomers ✓</p> <p>H is</p>  <p>E/Z happens because double bonds restricts rotation ✓</p> <p>different groups on each carbon of the double bond ✓</p> | 11 | <p>ALLOW same molecular formula ✓ but different structures ✓ Second marking point is DEPENDENT on first mark</p> <p>ALLOW compounds with the same structure 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 $CH_3CH_2CH_2CHCH_2$</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 i.e. <i>E</i>-pent-2-ene, <i>Z</i>-pent-2-ene and pent-1-ene</p> <p>ALLOW ecf on structures if wrong molecular formula used or consistent error or slip such as having just sticks</p> |

| Question | Expected Answers | Marks | Additional Guidance |
|----------|--|-----------|--|
| (b) | <p>from IR absorption, J contains O–H OR from IR J is an alcohol ✓</p> $C : H : O = \frac{70.59}{12.0} : \frac{13.72}{1.0} : \frac{15.69}{16.0}$ <p>OR 5.8825 : 13.72 : 0.9806 ✓</p> <p>empirical formula = C₆H₁₄O ✓</p> <p>(from mass spectrum), <i>M_r</i> = 102 ✓</p> <p>evidence that it has been shown that the empirical formula is the molecular formulae e.g. <i>M_r</i> of C₆H₁₄O = 102 so empirical formula is molecular formula ✓</p>  <p>One mark for each correct structure ✓ ✓ ✓</p> | 8 | <p>This is a QWC mark</p> <p>ALLOW two marks for correct empirical formula with no working out</p> <p>This is a QWC mark</p> <p>ALLOW structural or displayed formulae IGNORE incorrect names</p> <p>ALLOW one minor slip in drawing structures e.g. one missing hydrogen but ALLOW ecf for bigger slips such as showing just sticks and no hydrogen atoms ALLOW bond to H in OH</p> <p>ALLOW one mark for three isomers of C₆H₁₃OH whether branched or unbranched as a catch mark if no other mark has been awarded for the structures</p> <p>If more than three isomers of C₆H₁₃OH drawn</p> <ul style="list-style-type: none"> • 1 branched and 3 unbranched award two marks • any other combination award one mark <p>ALLOW one mark for hexan-1-ol, hexan-2-ol and hexan-3-ol if structures not drawn</p> |
| | Total | 19 | |

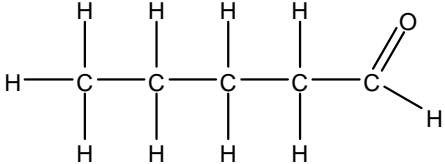
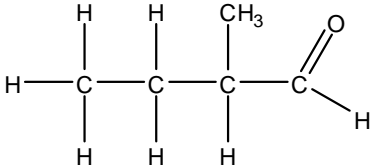
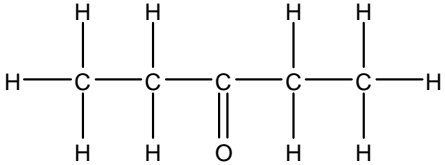
| Question | | er | Marks | Guidance |
|----------|-----|---|-------|---|
| 2 | (a) | (a compound) with no double bond (or triple bond) ✓ containing hydrogen and carbon only ✓ | 2 | ALLOW contains single bonds only ALLOW it contains just carbon and hydrogen DO NOT ALLOW a mixture of carbon and hydrogen OR only carbon and hydrogen molecules |
| | (b) | CH ₂ ✓ | 1 | ALLOW H ₂ C |
| | (c) | D and I OR F and G OR F and H ✓ | 1 | DO NOT ALLOW G and H |
| | (d) | (i) Cyclic hydrocarbons have more efficient combustion ✓ | 1 | The answer must relate to combustion or burning Assume 'they' refers to the cyclic hydrocarbons ALLOW cyclic hydrocarbons allow smoother burning OR cyclic hydrocarbons increase octane number OR cyclic hydrocarbons reduce knocking OR cyclic hydrocarbons are less likely to produce pre-ignition OR cyclic hydrocarbons are more efficient fuels OR cyclic hydrocarbons burn better OR easier to burn OR cyclic hydrocarbon combust more easily OR improves combustion DO NOT ALLOW cyclic hydrocarbons ignite more easily ALLOW ora for straight chain hydrocarbons IGNORE cyclic hydrocarbons increase volatility of fuel IGNORE cyclic hydrocarbons have a lower boiling point cyclic hydrocarbons are a better fuel on their own is NOT sufficient cyclic hydrocarbons burn more cleanly on their own is NOT sufficient |

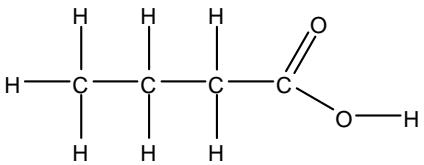
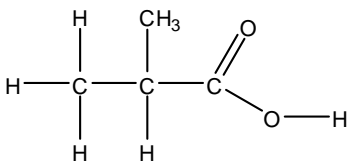
| Question | | Answer | Marks | Guidance |
|----------|----------|---|-------|--|
| | (d) (ii) | $C_7H_{16} \rightarrow C_7H_{14} + H_2$ ✓ | 1 | <p>ALLOW molecular formulae OR correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous)</p> <p>DO NOT ALLOW cycloheptane structure in equation</p> |
| | (e) | <p>D has more surface (area of) contact OR D is a bigger molecule ✓</p> <p>D has more van der Waals' forces OR C have fewer van der Waals' forces ✓</p> | 2 | <p>Both answers need to be comparisons Assume 'it' refers to D</p> <p>ALLOW has more electrons OR longer (carbon) chain OR higher molecular mass IGNORE surface area ALLOW ORA</p> <p>ALLOW D has stronger van der Waals' forces / larger VDW / greater VDW OR C has weaker van der Waals' forces OR C has smaller VDW</p> <p>ALLOW more VDW forces More intermolecular forces is not sufficient</p> <p>DO NOT ALLOW reference to bonds breaking or more bonds present unless it is clear that that the bonds are VDW</p> |
| | (f) | <p>Same structural formula ✓</p> <p>Different arrangement of groups around a double bond OR different arrangement (of atoms) in space ✓</p> | 2 | <p>ALLOW have the same structure / displayed formula / skeletal formula</p> <p>Stereoisomers have the same formula or molecular formula is not sufficient</p> <p>ALLOW different three dimensional arrangement</p> |

| Question | | er | Marks | Guidance |
|--------------|-------|--|-----------|--|
| | (g) | $C_7H_{16} + 11O_2 \rightarrow 7CO_2 + 8H_2O$ Correct reactants and products ✓ Balancing ✓ | 2 | ALLOW molecular formulae OR correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) ALLOW any correct multiple IGNORE state symbols Balancing is dependent on correct formulae |
| | (h) | $C_{16}H_{34} \rightarrow C_8H_{18} + 2C_4H_8$ ✓ | 1 | ALLOW molecular formulae OR correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) ALLOW any correct multiple ALLOW structural OR displayed OR skeletal formulae in equation ALLOW but-1-ene IGNORE state symbols |
| | (i) | (| 1 | ALLOW the 'part' (of the molecule or compound) that reacts ALLOW the group of atoms that gives the chemical properties ALLOW group of atoms which indicates the homologous series |
| | (ii) | 8 ✓ | 1 | |
| | (iii) | has an unpaired electron ✓ | 1 | ALLOW plural i.e. unpaired electrons has a lone OR single OR free electron is not sufficient |
| Total | | | 16 | |

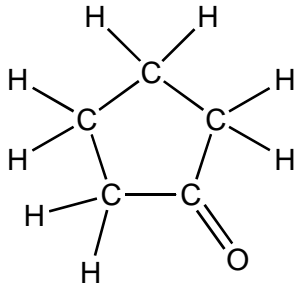
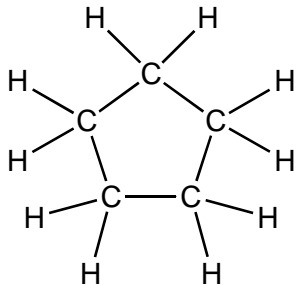
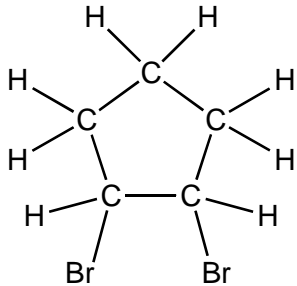
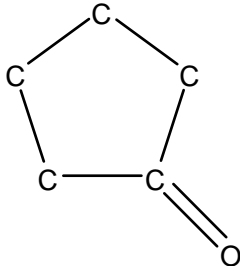
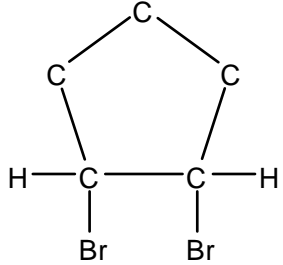
| Question | | Answer | Marks | Guidance |
|----------|-----|---|-------|--|
| 3 | (a) | <p>Any three from:</p> <p>Process 1 has a high atom economy OR has 100% atom economy OR a greater atom economy OR makes only the desired product ✓</p> <p>Process 1 has a very efficient conversion of reactants to products OR not much waste of starting material ✓</p> <p>Process 1 uses a lower pressure ✓</p> <p>Process 1 uses up toxic carbon monoxide ✓</p> <p>Process 1 uses methanol which can be produced from biomass ✓</p> | 3 | <p>Assume it refers to Process 1</p> <p>ALLOW process 1 has no waste OR process 1 has no co-products OR process 1 needs less separation OR process 1 has fewer other products OR gives only one product ALLOW ORA if process 2 is specified</p> <p>ALLOW ORA if process 2 is specified high percentage yield is not sufficient DO NOT ALLOW if percentage yield is explicitly linked to more waste (products) e.g. process 1 has a high percentage yield so makes little waste (product) scores 0 marks but process 1 makes no waste (product) and it has a high percentage yield scores 1 mark</p> <p>ALLOW ORA if process 2 is specified</p> <p>IGNORE process 2 comes from crude oil a non-renewable source ALLOW process 1 starts from a renewable source if the source is specified e.g. wood, municipal waste or sewage</p> <p>IGNORE reference to catalyst and rate of reaction</p> |

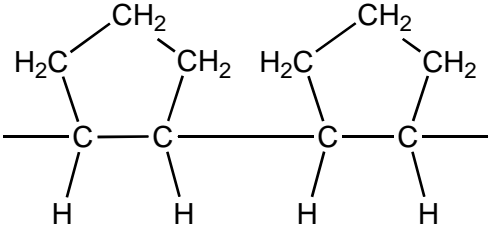
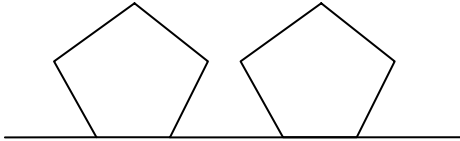
| Question | er | Marks | Guidance |
|----------|--|-------|--|
| (b) (| <p>Contains C=O bond because of absorption between 1700 and 1740 cm^{-1} (from the spectrum) ✓</p> <p>does not contain an O–H bond ✓</p> <p>(So was a) ketone OR aldehyde ✓</p> <p>$M_r = 86$ ✓</p> <p>Correct structure ✓</p> | 5 | <p>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</p> <p>ALLOW contains a carbonyl group because of absorption within range 1640–1750 cm^{-1} OR contains an aldehyde, ketone or carboxylic acid because of absorption within range 1640–1750 cm^{-1} ✓</p> <p>Mention of only an aldehyde or a ketone is not sufficient it needs reference to the wavenumber</p> <p>LOOK FOR THIS MARK ON THE SPECTRUM</p> <p>ALLOW not a carboxylic acid ✓</p> <p>ALLOW does not have any other characteristic absorbance due to other functional groups</p> <p>ALLOW (so was a) carbonyl compound</p> <p>ALLOW this mark if a structure of an aldehyde or a ketone is given even if the structure has an incorrect number of carbon atoms</p> <p>ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous)</p> <p>LOOK FOR AN ALDEHYDE or KETONE with FIVE carbon atoms OR a DIALDEHYDE, DIONE OR an OXOALDEHYDE with FOUR carbon atoms – a comprehensive list of correct structures is shown on page 34</p> <p>IGNORE incorrect name</p> <p>DO NOT ALLOW COH for an aldehyde</p> |

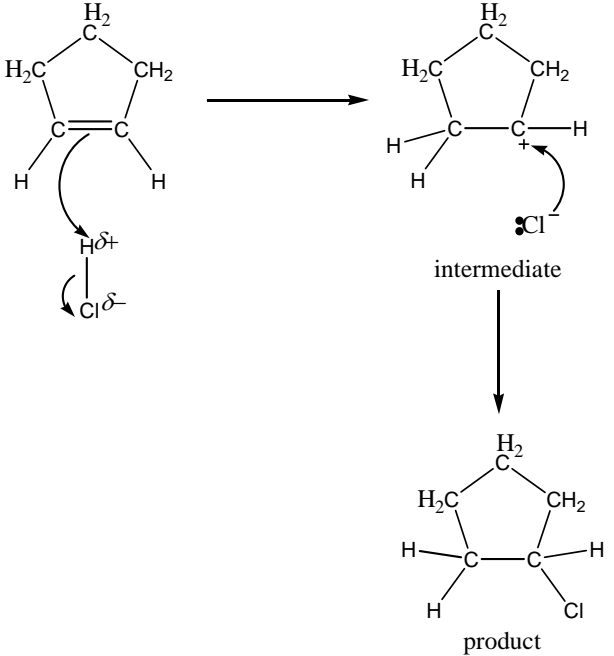
| Question | er | Marks | Guidance |
|----------|---|-------|---|
| | <p style="text-align: center;">  pentanal </p> <p>OR</p> <p style="text-align: center;">  2-methylbutanal </p> <p>OR</p> <p style="text-align: center;">  pentan-3-one </p> | | <p>ALLOW as a slip one stick with no H on in a displayed formula</p> |

| Question | er | Marks | Guidance |
|----------|--|-------|--|
| (b) (i) | <p>Correct structure ✓</p> <p>Name of the structure drawn ✓</p>  <p>butanoic acid</p> <p>OR</p>  <p>2-methylpropanoic acid</p> | 2 | <p>ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous)</p> <p>All bonds and all hydrogen atoms must be shown in a displayed formula within this question</p> <p>Name must correspond to the correct structure for two marks ALLOW butanoic acid or 2-methylpropanoic acid if the structure drawn is incorrect There is no ECF in this question</p> <p>ALLOW CH₃CH₂CH₂COOH</p> <p>ALLOW (CH₃)₂CHCOOH</p> <p>ALLOW methylpropanoic acid</p> |

| Question | | er | Marks | Guidance |
|----------|-----|---|-----------|--|
| | (c) | <p>Use of propan-1-ol ✓</p> $\text{CH}_3\text{COOH} + \text{C}_3\text{H}_7\text{OH} \rightarrow \text{CH}_3\text{COOCH}_2\text{CH}_2\text{CH}_3 + \text{H}_2\text{O}$ <p>Correct formulae for the ester ✓ Correctly balanced equation ✓</p> <p>Add H_2SO_4 OR acid catalyst OR H^+ ✓</p> | 4 | <p>ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous)</p> <p>ALLOW from the equation propanol OR $\text{C}_3\text{H}_7\text{OH}$ is not sufficient</p> <p>ALLOW molecular formula OR correct structural OR displayed OR skeletal formula OR mixture of the above</p> <p>ALLOW propan-2-ol in the equation</p> <p>ALLOW conditions mark over the arrow in the equation</p> |
| | | Total | 14 | |

| Question | er | Marks | Guidance |
|----------|---|-------|---|
| 4 (a) |  <p>compound P ✓</p>  <p>compound Q ✓</p>  <p>compound R ✓</p> | 3 | <p>ALLOW structures with missing hydrogen atoms on the carbon atoms that do not take part in the reaction. i.e. all hydrogen atoms must be shown in Q but not in P and R</p> <p>For example for the structures of P and R</p>   |

| Question | er | Marks | Guidance |
|----------|---|-------|---|
| (b) | Orange OR brown to colourless ✓ | 1 | ALLOW shades of orange OR yellow OR brown DO NOT ALLOW red alone DO NOT ALLOW any response that includes precipitate OR solid, irrespective of colour |
| (c) | Two or more repeat units ✓  | 1 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) Must have at least two repeat units and the free bonds at the end ALLOW free bonds with dotted lines All carbon–carbon bonds in the polymer chain must be shown IGNORE any brackets drawn IGNORE any missing hydrogen atoms on the CH ₂ groups ALLOW skeletal formula  |

| Question | er | Marks | Guidance |
|----------|---|-------|--|
| (d) | <p>Curly arrow from double bond to attack hydrogen of H-Cl/ and breaking of H-Cl/ bond ✓</p> <p>Correct dipole shown on H-Cl ✓</p> <p>Correct carbonium ion drawn ✓</p> <p>Curly arrow from Cl⁻ to the carbonium ion ✓</p>  <p>Correct product ✓</p> | 5 | <p>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</p> <p>Curly arrow must start from the double bond and not a carbon atom; other curly arrow must start from H-Cl/ bond</p> <p>DO NOT ALLOW dipoles on double bond</p> <p>Dipole must be partial charge and not full charge</p> <p>Carbocation needs a full charge and not a partial charge (charges do not need to be in a circle)</p> <p>Cl⁻ curly arrow must come from one lone pair on Cl⁻ ion OR from minus sign on Cl⁻ ion</p> <p>Lone pair does not need to be shown on Cl⁻ ion</p> <p>ALLOW structures with missing hydrogen atoms on the CH₂ groups</p> |

| Question | er | Marks | Guidance |
|----------|---|-------|--|
| (e) | <p>Nucleophilic substitution ✓</p> <p>Heterolytic (fission) spelt correctly ✓</p> <p>dipole shown on C—Cl bond, C^{δ+} and Cl^{δ-} ✓</p> <p>curly arrow from HO⁻ to carbon atom of C—Cl bond ✓</p> <p>curly arrow from C—Cl bond to the chlorine atom and formation of Cl ✓</p> | 5 | <p>ANNOTATE ANSWER WITH TICKS AND CROSSES ETC</p> <p>Dipole must be partial charge and not full charge</p> <p>HO⁻ curly arrow must come from one lone pair on O of HO⁻ ion OR from minus sign on HO⁻ ion</p> <p>curly arrow must start from C—Cl bond and not from C atom</p> <p>ALLOW structures with missing hydrogen atoms on the CH₂ groups</p> <p>ALLOW S_N1 mechanism dipole shown on C—Cl bond, C^{δ+} and Cl^{δ-} ✓ curly arrow from C—Cl bond to the Cl atom and Cl shown ✓ curly arrow from HO⁻ to correct carbonium ion ✓</p> |
| | Total | | |