

GCE

Chemistry A

Advanced Subsidiary GCE

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

Mark Scheme for January 2013

| Q | uestion | Answer | Marks | Guidance |
|---|---------|--|-------|--|
| 1 | (a) | C₃H ₇ ✓ | 1 | ALLOW H ₇ C ₃ |
| | (b) | Saturated Only has (carbon to carbon) single bonds ✓ | 2 | ALLOW does not contain any (carbon to carbon) double bonds ALLOW all of the carbon atoms are bonded to four other atoms |
| | | Hydrocarbon Contains (the elements) hydrogen and carbon only ✓ | | DO NOT ALLOW contains hydrogen and carbon DO NOT ALLOW a mixture of carbon and hydrogen only DO NOT ALLOW hydrogen and carbon molecules only |
| | (c) | | 1 | |
| | (d) | as branching increases the boiling point decreases OR the more branched the isomers of hexane are the lower the boiling point ✓ | 3 | ALLOW ORA throughout First marking point must compare boiling point and branching for all three isomers |
| | | branched isomers have less surface (area) of contact OR branched fewer points of contact (than unbranched) ✓ | | Reference to just surface area / closeness of molecules is not sufficient |
| | | (the more branched the) fewer van der Waals' forces OR (the more branched) has weaker van der Waals' forces OR Less energy required to break van der Waal's forces ✓ | | ALLOW vdw forces OR VDW forces (and any combination of upper and lower cases) DO NOT ALLOW VDW mark if answer states that these are between atoms or answer implies that these are bonds |
| | (e) | $C_{10}H_{22} \rightarrow C_6H_{14} + C_4H_8$ OR $C_{10}H_{22} \rightarrow C_6H_{14} + 2C_2H_4 \checkmark$ | 1 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) IGNORE state symbols |

| Q | Question | | Answer | | Marks | Guidance |
|---|----------|------|---|----------------------|-------|--|
| 1 | (f) | (i) | $C_4H_{10} + 2Cl_2 \rightarrow C_4H_8Cl_2 + 2H_8$ | Cl✓ | 1 | IGNORE state symbols |
| | | (ii) | Isomer 1 | Isomer 2 | 2 | |
| | | | | 1,3-dichlorobutane ✓ | | Must be a displayed formula |
| | | | Correct displayed formula eg: | | | ALLOW absence of hyphens 1 and 3 must be clearly separated: ALLOW full stops: 1.3 OR spaces: 1 3 DO NOT ALLOW 13 |
| | (g) | (i) | covalent bond breaking ✓ | | 2 | ALLOW covalent bond is split |
| | | | one electron (from the bond pair) g OR makes (two) radicals ✓ | oes to each atom | | IGNORE particle for atom DO NOT ALLOW molecule or compound for atom DO NOT ALLOW to each molecule or to each reactant ALLOW one electron goes to each product / species IGNORE homolytic fission equations |
| | | (ii) | $Cl + C_4H_9Cl \rightarrow C_4H_8Cl$ $C_4H_8Cl + Cl_2 \rightarrow C_4H_8Cl_2 + C$ | | 2 | IGNORE dots even if incorrect |
| | (h) | | $C_4H_{10} + 4\frac{1}{2}O_2 \rightarrow 4CO + 5H_2O$ OR $C_4H_{10} + 2\frac{1}{2}O_2 \rightarrow 4C + 5H_2O \checkmark$ | | 1 | ALLOW any correct multiples for these equations eg $2C_4H_{10} + 9O_2 \rightarrow 8CO + 10 H_2O$ IGNORE state symbols ALLOW equations for incomplete combustion that give CO_2 with CO and/or C eg $C_4H_{10} + 4O_2 \rightarrow 3CO + C + 5H_2O$ |
| | | | | Total | 16 | |

| Q | uesti | on | Answer | Marks | Guidance |
|---|-------|----|---|-------|---|
| 2 | (a) | | (enthalpy change for the) formation of one mole (of P_4O_{10}) \checkmark from (constituent) elements OR from P_4 /phosphorus and O_2 /oxygen \checkmark | 2 | ALLOW energy required OR energy released ALLOW makes one mole of product/substance/molecule/ compound ALLOW made from P and O ₂ OR made from two elements IGNORE comments related to standard conditions |
| | (b) | | FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer = -368 (kJ mol ⁻¹) award 3 marks (+)2984 +(+)6 × 286 OR (+)2984 +(+)1716 OR (+)4700 ✓ | 3 | IF there is an alternative answer, check to see if there is any ECF credit possible using working below. See list below for marking of answers from common errors. IGNORE sign |
| | | | | | |
| | | | (−)1267 × 4 OR (−)5068 ✓ | | IGNORE sign |
| | | | -368 ✓ | | ALLOW ECF for enthalpy change of products – enthalpy change of reactants |
| | | | | | ALLOW for 2 marks: +368 cycle wrong way around OR -1798 no × 6 OR (+)3433 no x 4 OR -3352 missing 2984 OR (+) 9768 product the wrong sign around OR (-) 9768 reactants the wrong sign |
| | | | | | ALLOW for 1 mark: (+)1798 no x 6 and cycle wrong way around OR -3433 cycle wrong way around and not × 4 OR (+)3352 missing 2984 and cycle wrong way around OR (+)2003 no x 6 or x 4 OR (+)449 missing 2984 and x 4 OR -4782 missing 2984 and x 6 Note: There may be other possibilities. |

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| | | - | |
|------|---|----|----|
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| Q | uestion | Answer | Marks | Guidance |
|---|---------|---|-------|--|
| | (c) | P ₄ + 5O ₂ + 6H ₂ O → 4H ₃ PO ₄ ✓ Only the desired product is made ✓ Second marking point can only be awarded if the equation is correct. | 2 | ALLOW there are no waste products OR there are no by-products OR there is only one product. DO NOT ALLOW it is an addition reaction |
| | | Total | 7 | |

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|---|----------|-------|---|---------------|---|
| C | Question | | Answer | Marks | Guidance |
| 3 | (a) | (i) | acid ✓ | 1 | ALLOW named mineral acid or correct formula eg phosphoric acid, H ₃ PO ₄ , sulfuric acid, H ₂ SO ₄ or H ⁺ DO NOT ALLOW any carboxylic acids |
| | | (ii) | $C_5H_{12}O \rightarrow C_5H_{10} + H_2O \checkmark$ | 1 | DO NOT ALLOW use of C₅H₁₁OH |
| | | (iii) | structural isomerism have the same molecular formula ✓ but different structural formulae ✓ | 4 | Same formula is not sufficient ALLOW different structure OR different displayed formula OR different skeletal formula Different formula or different arrangement of atoms is not sufficient ALLOW different structural arrangement (of atoms) |
| | | | stereoisomerism have the same structural formula ✓ but different arrangement (of atoms) in space ✓ | | ALLOW have the same structure Stereoisomers have the same formula or molecular formula is not sufficient ALLOW different spatial arrangements (of atoms) |

| Quest | ion | Answer | Marks | Guidance |
|-------|------|--|-------|---|
| 3 (a) | (iv) | CH ₃ H CH ₂ CH ₃ H CH ₂ CH ₃ B H CH ₂ CH ₂ CH ₃ C Correct structure for A Correct structure for C Correct structure for C | 3 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above A and B must clearly show cis and trans configuration eg A B Answers to A and B are interchangeable C: CH ₂ CHCH ₂ CH ₂ CH ₃ ALLOW -C ₂ H ₅ group in A or B or -CH ₂ C ₂ H ₅ in C DO NOT ALLOW -C ₃ H ₇ group in C |

| Question | Answer | Marks | Guidance |
|-----------|--|-------|--|
| 3 (a) (v) | carbon–carbon double bond ✓ Each carbon atom in the double bond is attached to (two) different groups/atoms ✓ | 2 | IGNORE comments about rotation ALLOW carbon double bond ALLOW Each carbon atom of the double bond is attached to a H and an alkyl group DO NOT ALLOW functional groups for groups DO NOT ALLOW the carbon atoms are attached to different groups "Each carbon atom in the double bond" implies a carbon—carbon double bond for the first marking point |
| (b) | + [O] + H ₂ O Correct skeletal structure of product ✓ Balanced equation ✓ | 2 | Balancing mark can only be awarded if the equation has a correct skeletal formula for the product |
| (c) | distil with H ₂ SO ₄ / heat under reflux with H ₂ SO ₄ / K ₂ Cr ₂ O ₇ CH ₃ CH ₂ CH ₂ CHO CH ₃ CH ₂ CH ₂ CHO CH ₃ CH ₂ CH ₂ CHO | 2 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) eg H H C C C C C C C C C C C C C C C C C |
| | Total | 15 | |

| C | uestion | Answer | Marks | Guidance |
|---|---------|---|-------|--|
| 4 | (a) | FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer = 681 (kJ) award 3 marks | 3 | IF there is an alternative answer, check to see if there is any ECF credit possible using working below. |
| | | Evidence of dividing 1000 by 24 | | ALLOW 41.7 up to calculator value 41.6666667 correctly rounded. |
| | | | | ALLOW $\frac{1000}{24}$ for first marking point if not calculated |
| | | Evidence of dividing by 3 and multiplying by 49 in the calculation | | ALLOW energy released per mole = 16.3 ✓ |
| | | energy released = 681 (kJ) ✓ | | IGNORE (–) sign in the answer |
| | | (MUST BE TO 3 SIG FIGS) | | Common Incorrect answers 0.392 scores 2 marks 392000 scores 2 marks |

| Q | uesti | on | Answer | Marks | Guidance |
|---|-------|----|--|--------|--|
| 4 | (b) | | enthalpy $ \begin{array}{c} E_a \\ 3H_2(g) + CO_2(g) \\ & CH_3OH(g) + H_2O(g) \end{array} $ progress of reaction | 3 | |
| | | | CH ₃ OH and H ₂ O added as product AND shown below the reactants ✓ | | IGNORE state symbols for the products |
| | | | ΔH labelled with arrow pointing towards products product line if no products stated \checkmark | or the | IF there is no ΔH labelled then ALLOW –49 only as an alternative label for ΔH IF ΔH is labelled then IGNORE any numerical value DO NOT ALLOW $-\Delta H$ ALLOW this arrow even if it has a small gap at the top and bottom i.e. does not quite reach reactant or product line |
| | | | E _a labelled correctly AND above reactants ✓ | | ALLOW (+) 225 only as an alternative label for E_a ALLOW arrows at both ends of activation energy line The E_a line must point to maximum (or near to the maximum) on the curve ALLOW this line even if it has a small gap at the top and bottom ie does not quite reach the maximum or reactant line ALLOW A_E or E_a for activation energy |

| Answer | Marks | Guidance |
|---|---|---|
| (+)49 ✓ | 1 | DO NOT ALLOW –49 |
| (+)274 ✓ | 1 | DO NOT ALLOW –274 ALLOW answer to (c) + 225 as ECF |
| (equilibrium position shifts) to the left ✓ (Forward) reaction is exothermic OR reaction gives out heat | 2 | ALLOW 'favours the left', as alternative for 'shifts equilibrium to left' Note: ALLOW suitable alternatives for 'to left', eg: towards CO ₂ / H ₂ OR towards reactants OR in backward |
| OR reverse reaction is endothermic OR reverse reaction takes in heat ✓ | | direction OR in reverse direction OR decreases yield of CH ₃ OH /products |
| The explanation mark is dependent on the correct shift of the equilibrium | | IGNORE responses in terms of rate |
| (equilibrium position) shifts to the left ✓ | 2 | ALLOW 'favours the left', as alternative for 'shifts equilibrium to left' Note: ALLOW suitable alternatives for 'to left', eg: towards CO ₂ / H ₂ OR towards reactants OR in backward direction OR in reverse direction OR decreases yield of CH ₃ OH /products |
| | | IGNORE responses in terms of rate |
| Right-hand side has fewer (gaseous) moles/molecules ✓ ORA | | ALLOW four moles on the left and two moles on the right ALLOW more moles of reactants or fewer moles of products |
| The explanation mark is dependent on the correct shift of the equilibrium | | ASSUME "goes the side with more gas molecules" implies from equation that more molecules on the left OR "goes to side with fewer gas molecules" implies from equation that fewer molecules are on the right |
| | (+)49 ✓ (+)274 ✓ (equilibrium position shifts) to the left ✓ (Forward) reaction is exothermic OR reaction gives out heat OR reverse reaction is endothermic OR reverse reaction takes in heat ✓ The explanation mark is dependent on the correct shift of the equilibrium (equilibrium position) shifts to the left ✓ Right-hand side has fewer (gaseous) moles/molecules ✓ ORA The explanation mark is dependent on the correct shift | (+)49 ✓ 1 (+)274 ✓ 1 (equilibrium position shifts) to the left ✓ 2 (Forward) reaction is exothermic OR reaction gives out heat OR reverse reaction is endothermic OR reverse reaction takes in heat ✓ The explanation mark is dependent on the correct shift of the equilibrium (equilibrium position) shifts to the left ✓ 2 Right-hand side has fewer (gaseous) moles/molecules ✓ ORA The explanation mark is dependent on the correct shift |

| Question | Answer | Marks | Guidance |
|----------|--|-------|---|
| (g) | Adsorption of reactants OR adsorption of gases OR H₂ and CO₂ attached to surface ✓ | 3 | ALLOW CO ₂ and H ₂ (weakly) bonded to surface OR reactants bond to surface OR CO ₂ and H ₂ form temporary bonds with the catalyst DO NOT ALLOW absorption |
| | Bonds weaken in reactants OR chemical reaction OR activation energy decreases ✓ | | ALLOW bonds weaken in H ₂ OR bonds weaken in CO ₂ OR C=O bonds weaken OR bonds break and new bonds made in product OR H ₂ O and CH ₃ OH made |
| | Desorption of products OR desorption of H₂O and CH₃OH ✓ | | ALLOW products leave the surface/catalyst OR H ₂ O and CH ₃ OH no longer bonded to surface/catalyst ALLOW deadsorption OR adsorb from for desorption ALLOW diffuse away for desorption |
| | Total | 15 | |

| Q | uestio | | Marks | Guidance |
|---|--------|---|-------|---|
| 5 | (a) | Answer FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer = 90% award 3 marks amount of dichloroethane = $\frac{19800000}{99.0}$ OR 200000 (mol) OR 2 × 10 ⁵ (mol) \checkmark amount of chloroethene = $\frac{11250000}{62.5}$ OR 180000 (mol) OR 1.8 × 10 ⁵ (mol) \checkmark Calculates percentage yield = $\frac{180000}{200000}$ × 100 = 90 % \checkmark | 1 ' ' | IF there is an alternative answer, check to see if there is any ECF credit possible using working below. ALLOW approach based on mass for 2nd and 3rd marks Theoretical mass of chloroethene = 200000×62.5 OR 12500000 (g) OR $1.25 \times 10^7 \text{ (g)} \checkmark$ Calculates percentage yield = $\frac{11250000}{12500000} \times 100 = 90 \% \checkmark$ ALLOW approach based on grams rather than tonnes: $n(\text{dichloroethane}) = \frac{19.80}{99.0}$ OR $0.2 \text{ (mol)} \checkmark$ $n(\text{chloroethane}) = \frac{11.25}{62.5}$ OR 0.18 (mol) OR theoretical mass chloroethane = 0.2×62.5 OR $12.5 \text{ g} \checkmark$ % yield = $\frac{0.18}{0.20} \times 100 = 90 \%$ OR $\frac{11.25}{12.5} \times 100 = 90 \% \checkmark$ ALLOW ECF throughout from wrong M_r value(s) with final % yield to 2 or more significant figures DO NOT ALLOW final mark for an answer above 100% Note: |
| | | | | If this is the only working seen award no marks ie $\frac{11.25 \times 10^6}{19.80 \times 10^6} \times 100 = 56.81\%$ |

| Q | uestion | Answer | Marks | Guidance |
|---|---------|---|-------|--|
| 5 | (b) | FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer = (+)62 award 3 marks | 3 | IF there is an alternative answer, check to see if there is any ECF credit possible. |
| | | ΔH for bonds broken = 2691 (kJ mol ⁻¹) \checkmark | | IGNORE sign ALLOW 1106 (C–Cl, C–C and C–H bonds) |
| | | ΔH for bond formed = 2629 (kJ mol ⁻¹) \checkmark | | IGNORE sign ALLOW 1044 (H–Cl and C=C bonds) |
| | | $\Delta H = (+)62 \text{ (kJ mol}^{-1}) \checkmark$ | | ECF based on bonds broken – bonds formed |
| | | | | ALLOW 2 marks for –62 |
| | (c) | Displayed formulae of monomer and polymer required for the marks. | 3 | Polymer must have side links (do not have to cut through bracket and can be dotted lines) |
| | | $\begin{array}{cccccccccccccccccccccccccccccccccccc$ | | ALLOW a correct section of the polymer with side links as below would score two marks as the equation is not balanced CI |
| | | Only the correct polymer on right hand side ✓ | | DO NOT ALLOW ECF from wrong monomer |
| | | | | 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) |
| | | A correctly balanced equation using displayed formulae for any monomer and matching polymer including the correct use of $n \checkmark$ | | The equation below would be worth 1 mark for balancing n H C H H H N H H H N H H H N H H |

| Quest | ion | Answer | Marks | Guidance |
|-------|-------|--|-------|--|
| (d) | (i) | React with an alkali OR react with a base/carbonate OR Bubble through water (to make HCl(aq)) OR dissolve in water ✓ | 1 | ALLOW react with a named alkali or base eg calcium carbonate, calcium hydroxide, magnesium oxide, ammonia ALLOW an appropriate chemical formula IGNORE use of gas scrubbers |
| | (ii) | Sort and recycle ✓ Organic feedstock OR cracked ✓ | 2 | ALLOW separate and recycle or sorting and remoulding ALLOW use for the production organic compounds OR synthesis gas ALLOW the production of plastics or monomers or new polymers |
| | (iii) | (Bio) degradable (polymers) OR compostable (polymers) OR soluble (polymers) OR photodegradable (polymers) ✓ | 1 | IGNORE a named polymer if degradable DO NOT ALLOW any addition polymer eg PTFE |
| | | Total | 13 | |

| C | Questi | ion | Answer | Marks | Guidance |
|---|--------|-----|---|-------|--|
| 6 | (a) | | Bond breaking absorbs energy AND bond forming releases energy ✓ More energy released than absorbed ✓ The second marking point is dependent on the correct identification of the energy changes during bond breaking and bond making | 2 | ALLOW bond breaking is endothermic AND bond forming is exothermic DO NOT ALLOW bond forming requires energy ALLOW more energy is released when the bond in the products are formed than is required to break the bonds in the reactants ALLOW exothermic change transfers more energy than endothermic change OR bond forming transfers more energy than bond breaking OR '(the sum of the) bond enthalpies in the products is greater than the (sum of the) bond enthalpies in the reactants' OR '(the sum of the) bond enthalpies of the bonds made is greater than (the sum of) the bond enthalpies of the bonds broken' OR more energy associated with bond making than with bond breaking IGNORE reference to strong and weak bonds IGNORE reference to number of bonds broken or made IGNORE enthalpy of products is less than enthalpy of reactants |
| | (b) | (i) | (C=O) bond vibrates (more) OR bond bends (more) OR bond stretches (more) ✓ | 1 | IGNORE molecule vibrates/rotates "It" refers to the molecule and is insufficient DO NOT ALLOW any reference to bond breaking. DO NOT ALLOW a stated bond if not present in CO ₂ eg C-O, C-H |

| Question | Answer | Marks | Guidance |
|----------|---|-------|--|
| (ii) | Any two from: (injected) deep into the oceans / sea ✓ | 2 | DO NOT ALLOW reference to carbon being stored – the answer must either refer to carbon dioxide or not mention the name of the stored substance. Assume "it" refers to CO ₂ |
| | | | DO NOT ALLOW dumping waste at the bottom of the sea |
| | | | ALLOW on the sea-bed |
| | | | DO NOT ALLOW dissolve CO ₂ in the sea OR (stored) in ocean |
| | (Stored) in geological formations OR (stored) deep in rocks OR (stored) in old mines OR (stored) in old oil wells OR old gas fields ✓ | | DO NOT ALLOW geographical formations ALLOW stored under the sea (bed) ALLOW pumped into oil wells to force last bit of oil out DO NOT ALLOW buried underground |
| | (Stored) by reaction with metal oxides OR reaction to form (solid) carbonates OR (stored) as a carbonate | | DO NOT ALLOW react with metals to form carbonates |
| | OR equation to show formation of metal carbonate ✓ | | IGNORE mineral storage |

| Question | Answer | Marks | Guidance |
|----------|---|-------|---|
| (c) | Any two from: | 2 | |
| | Energy demand Low(er) temperature (can be used) OR reduces CO₂ emissions (from burning fossil fuels) ✓ | | ALLOW 'allows use of room temperature' OR 'allows use of a lower pressure' OR uses less fuel |
| | | | IGNORE lower energy demand OR lower activation energy IGNORE cheaper IGNORE less greenhouse gases OR reduces global warming |
| | Specificity enzymes have a great deal of specificity ✓ | | ALLOW making specific isomers / enantiomers ALLOW for making pure products ALLOW generating specified products |
| | Atom economy greater atom economy OR less waste ✓ | | ALLOW increases atom economy |
| | Toxicity can reduce use of toxic solvents OR reduces use of toxic catalysts OR reduces the use of toxic reactants ✓ | | ALLOW reduce use of hazardous/toxic/harmful/poisonous chemicals ALLOW enzymes are non toxic IGNORE can be reused |

| Question | Answer | Marks | Guidance |
|----------|--|-------|---|
| (d) | Catalyst lowers the activation energy (because of a different reaction pathway) ✓ Diagram of Boltzmann distribution ✓ | 5 | Can be scored from the diagram by correctly labelling $E_{a cat}$ closer to the origin than E_a Boltzmann distribution must start at origin AND must not touch x -axis at high energy |
| | axes labelled (number of) molecules and energy ✓ E _{a cat} | | DO NOT ALLOW Boltzmann distribution mark if two curves drawn |
| | (number of) | | DO NOT ALLOW Boltzmann distribution curve bending upwards at higher energy |
| | molecule | | ALLOW particles instead of molecules DO NOT ALLOW the first use of atoms but credit atoms if used in a subsequent marking point |
| | energy extra molecules with energy above E _a | | DO NOT ALLOW enthalpy on x-axis instead of energy |
| | Greater proportion of molecules with energy above activation energy with catalyst ✓ | | ALLOW more molecules with energy above activation energy (with a catalyst) OR more molecules overcome the activation energy (with a |
| | | | catalyst) OR more molecules have enough energy to react (with a catalyst) OR more molecules are able to react at lower energies |
| | more effective collisions OR more successful collisions OR increased frequency of successful collisions ✓ | | More collisions OR more frequent collisions are not sufficient |
| | Total | 12 | |

| Question | Answer | Marks | Guidance |
|----------|--|-------|--|
| 7 | Nucleophilic substitution reaction | 5 | ANNOTATE ANSWER WITH TICKS AND CROSSES ETC |
| | correct equation for the reaction \checkmark CH ₂ CHCH ₂ CH ₂ Cl + KOH \rightarrow CH ₂ CHCH ₂ CH ₂ OH + KCl | | ALLOW CH ₂ CHCH ₂ CH ₂ Cl + OH ⁻ → CH ₂ CHCH ₂ CH ₂ OH + Ct |
| | OR $C_4H_7Cl + KOH \rightarrow C_4H_7OH + KCl$ | | ALLOW $C_4H_7Cl + OH^- \rightarrow C_4H_7OH + Cl^-$ |
| | | | ALLOW correct molecular OR structural OR displayed OR skeletal formula OR mixture of the above |
| | correct product of the reaction ✓ H H H H C C C C C OH H H H | | For structure of the product ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) if seen ONCE in equation, mechanism or drawn out eg CH ₂ CHCH ₂ CH ₂ OH |
| | Mechanism | | |
| | $\begin{array}{cccccccccccccccccccccccccccccccccccc$ | | |
| | dipole shown on C—C l bond: C^{δ^+} and Cl^{δ^-} in the correct chloroalkene \checkmark | | curly arrow must start from one lone pair on O atom of ⁻ OH ion OR from negative charge on the O atom of the ⁻ OH ion |
| | curly arrow from HO [−] to carbon atom of C–C <i>l</i> bond AND curly arrow from C–C <i>l</i> bond to chlorine atom ✓ | | Lone pair does not need to be shown on ⁻ OH ion |
| | formation of C1 - ✓ | | |

| Question | Answer | , , , | , · |
|----------|---|-------|--|
| Question | Answer Nucleophilic substitution continued (S_N1) Step 1: H H Step 2: H H H H H H H H H H H H H | Marks | Guidance ANNOTATE ANSWER WITH TICKS AND CROSSES ETC ALLOW S _N 1 mechanism dipole shown on C–CI bond, C ^{δ+} and CI ^{δ−} in correct chloroalkene ✓ curly arrow from C–CI bond to halogen atom and Cl ⁻ ✓ curly arrow from OH to correct carbocation ✓ curly arrow must start from one lone pair on O atom of OH ion OR from negative charge on the O atom of the OH ion Lone pair does not need to be shown on OH ion |

| Question | Answer | Marks | Guidance |
|----------|--|-------|---|
| 7 | Electrophilic addition | 6 | ANNOTATE ANSWER WITH TICKS AND CROSSES ETC |
| | correct equation for the reaction ✓ | | |
| | CH ₂ CHCH ₂ CH ₂ CI + HBr → CH ₃ CHBrCH ₂ CH ₂ CI OR CH ₂ CHCH ₂ CH ₂ CI + HBr → CH ₂ BrCH ₂ CH ₂ CH ₂ CI | | ALLOW correct molecular OR structural OR displayed OR skeletal formula OR mixture of the above. |
| | Indication that there are two possible addition products ✓ | | eg C_4H_7Cl + HBr $\rightarrow C_4H_8BrCl$ |
| | Correct product ✓ | | For the structure of the product ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) if seen ONCE in equation, mechanism or drawn out eg CH ₂ BrCH ₂ CH ₂ CH ₂ Cl or CH ₃ CHBrCH ₂ CH ₂ Cl |
| | Mechanism Curly arrow from C=C of correct chloroalkene to attack the H atom in HBr ✓ | | curly arrow must start from covalent bonds and not atoms Lone pair does not need to be shown on ion or used in mechanism |
| | Correct dipole on H–Br: $H^{\delta+}$ and $Br^{\delta-}$ AND curly arrow from H–Br bond to Br \checkmark | | DO NOT ALLOW any other partial charges eg shown on double bond |
| | Correct carbocation / carbonium ion with the full positive charge shown: C ⁺ | | DO NOT ALLOW C^{δ^+} for charge on carbonium ion. |
| | correct curly arrow from lone pair of Br⁻ to correct carbon atom OR correct curly arrow from negative charge of Br⁻ to correct carbon atom ✓ | | Curly arrow from Br ⁻ can start from the negative charge or the lone pair DO NOT ALLOW delta negative, i.e. Br ⁵⁻ |

| Question | Answer | Marks | Guidance |
|----------|--|-------|----------|
| 7 | Electrophilic addition continued | | |
| | $\begin{array}{c ccccccccccccccccccccccccccccccccccc$ | | |
| | H H H H H H H H H H H H H H H H H H H | | |
| | $\begin{array}{c ccccccccccccccccccccccccccccccccccc$ | | |
| | H H H H H H H C C C C C C C C C C C C C | | |
| | heterolytic fission for both mechanisms and not contradicted ✓ | 1 | |

| Ougation | A nower | · , | T . |
|----------|--|-------|----------|
| Question | Answer | Marks | Guidance |
| | ALTERNATIVE APPROACH The Candidate who reacts with KOH followed by HBr | | |
| | Award all marks for the nucleophilic substitution mechanism as per the marking scheme You can award all marks for the electrophilic addition mechanism; however the product will be one of the following: | | |
| | $ \begin{bmatrix} & & & & & & & & & & & & & & & & & & $ | | |
| | The mechanism will be the same except the -Cl will now be replaced by -OH at every stage | | |
| | Total | 12 | |

| Question | Answer | Marks | Guidance |
|----------|---|-------|---|
| 8 | IR spectrum (absorbance between) 3200–3550 cm ⁻¹ indicates –OH AND X is an alcohol ✓ | 1 | LOOK ON THE SPECTRUM for labelled absorbance which can be given credit ALLOW an absorbance within the range 3100 to 3700cm ⁻¹ from the spectrum. Answer must give –OH and alcohol for the mark. IGNORE phenol DO NOT ALLOW carboxylic acid (there is no carbonyl group present in the spectrum) |
| | Formula mole ratio C: H: O $\frac{0.600}{12}:\frac{0.133}{1.0}:\frac{0.267}{16} \text{ OR } 0.0500:0.133:0.0167 \checkmark$ $\frac{0.05}{0.0167}:\frac{0.133}{0.0167}:\frac{0.0167}{0.0167} \text{ OR } 3:8:1 \text{ OR } C_3H_8O \checkmark$ Candidate links C_3H_8O to 60 such as C_3H_8O has M_r 60 OR C_3H_8O has $m/z=60$ \checkmark | 3 | Must be a clear link between the formula and the M_r OR m/z ALLOW evidence of M_r , eg $(12 \times 3) + (8 \times 1) + 16$; $36 + 8 + 16 = 60$ ALLOW alternative approach for empirical formula and evidence that 60 is equal to C_3H_8O $M_r = 60$ Carbon Hydrogen $60 \times \frac{60}{100} = 36 60 \times \frac{13.3}{100} = 8$ $36/12 = 3 C 8/1 = 8H$ $36 + 8 = 44 60 - 44 = 16 \text{ so } 1 O C_3H_8O$ |

| Question | Answer | Marks | Guidance |
|----------|---|-------|--|
| 8 | Identification and equation | 6 | |
| | X is CH ₃ CH ₂ CH ₂ OH OR CH ₃ CHOHCH ₃ OR either CH ₃ CH ₂ CH ₂ OH or CH ₃ CHOHCH ₃ ✓ | | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) |
| | | | IGNORE names |
| | QWC Stated in words that Y must be an ester because it is made from the reaction of a carboxylic acid AND X (propan-1-ol OR propan-2-ol OR an alcohol) ✓ | | ALLOW a carboxylic acid reacts with an alcohol to give an ester. IGNORE ethanoic acid (as this is stated in the question) |
| | Y is CH ₃ COOCH ₂ CH ₂ CH ₃ OR CH ₃ COOCH(CH ₃) ₂ OR either CH ₃ COOCH ₂ CH ₂ CH ₃ or CH ₃ COOCH(CH ₃) ₂ ✓ Must be consistent with a structure of alcohol X | | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) If no structure of X is provided one mark can be awarded for |
| | | | a correct structure of CH ₃ COOCH ₂ CH ₂ CH ₃ OR CH ₃ COOCH(CH ₃) ₂ |
| | $m/z = 31$ is $CH_2OH^+ \checkmark$ | | DO NOT ALLOW CH ₃ O ⁺ |
| | QWC $m/z = 31$ or CH_2OH indicates that X must be $CH_3CH_2CH_2OH$ OR cannot be $CH_3CH(OH)CH_3$ OR shows that X is the primary alcohol \checkmark | | QWC must link the evidence to the structure of propan-1-ol. |
| | $C_3H_8O + C_2H_4O_2 \rightarrow C_5H_{10}O_2 + H_2O \checkmark$ | | In equation ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above |
| | Total | 10 | |