

Chemistry A

Advanced GCE A2 H434

Advanced Subsidiary GCE AS H034

Mark Schemes for the Units

June 2009

H034/H434/MS/R/09

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Advanced Subsidiary GCE Chemistry (H034)

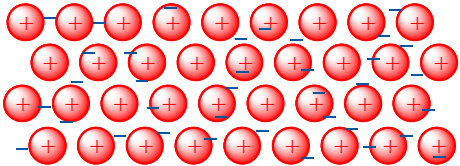
MARK SCHEME FOR THE UNITS

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F321 Atoms, Bonds and Groups

| Question | | | Expected Answers | Marks | Additional Guidance | | | | | | | | | | | | |
|------------------|---------|----------|--|-------|---|----------|-----------|------------------|----|----|----|------------------|----|----|----|---|-------------|
| 1 | (a) | (i) | <table border="1"> <thead> <tr> <th></th> <th>protons</th> <th>neutrons</th> <th>electrons</th> </tr> </thead> <tbody> <tr> <td>²⁴Mg</td> <td>12</td> <td>12</td> <td>12</td> </tr> <tr> <td>²⁵Mg</td> <td>12</td> <td>13</td> <td>12</td> </tr> </tbody> </table> <p>²⁴Mg line correct ✓ ²⁵Mg line correct ✓</p> | | protons | neutrons | electrons | ²⁴ Mg | 12 | 12 | 12 | ²⁵ Mg | 12 | 13 | 12 | 2 | mark by row |
| | protons | neutrons | electrons | | | | | | | | | | | | | | |
| ²⁴ Mg | 12 | 12 | 12 | | | | | | | | | | | | | | |
| ²⁵ Mg | 12 | 13 | 12 | | | | | | | | | | | | | | |
| | | (ii) | $\frac{24 \times 78.60 + 25 \times 10.11 + 26 \times 11.29}{100}$ <p>OR $18.8640 + 2.5275 + 2.9354$</p> <p>OR 24.3269 ✓</p> <p>$A_r = 24.33$ (to 4 sig figs) ✓</p> | 2 | <p>ALLOW two marks for $A_r = 24.33$ with no working out</p> <p>ALLOW one mark for ecf from incorrect sum provided final answer is between 24 and 26 and is to 4 significant figures, e.g. 24.3235 ✗ gives ecf of 24.32 ✓</p> | | | | | | | | | | | | |
| | | (iii) | <p>The (weighted) mean mass of an atom OR (weighted) average mass of an atom ✓</p> <p>relative to 1/12th (the mass) ✓</p> <p>of (one atom of) ¹²C ✓</p> | 3 | <p>ALLOW The (weighted) mean mass OR (weighted) average mass of an atom OR average atomic mass ✓ compared with (the mass of) carbon-12 ✓ which is 12 ✓</p> <p>For 1st marking point, ALLOW mean mass of the isotopes OR average mass of the isotopes Do NOT ALLOW the singular: isotope</p> <p>ALLOW mass of one mole of atoms ✓ compared to 1/12th ✓ (the mass) of one mole / 12 g of carbon-12 ✓</p> | | | | | | | | | | | | |

| Question | | Expected Answers | Marks | Additional Guidance |
|----------|-------|--|-----------|--|
| | | | | $\frac{\text{mass of one mole of atoms}}{12}$ ✓ 1/12th ✓ the mass of one mole / 12 g of carbon-12 ✓ |
| (b) | (i) | Mg ✓ oxidation number changes from 0 to (+)2 OR oxidation number increases by 2 ✓ | 2 | ALLOW correct oxidation numbers shown in equation 2nd mark is dependent on identification of Mg IGNORE electrons |
| | (ii) | Mg/solid dissolves OR Mg/solid disappears OR (Mg/solid) forms a solution ✓ bubbles OR fizzes OR effervesces OR gas produced ✓ | 2 | IGNORE metal reacts IGNORE temperature change IGNORE steam produced DO NOT ALLOW carbon dioxide gas produced DO NOT ALLOW hydrogen produced without gas |
| (c) | (i) | $M(\text{MgSO}_4) = 120.4 \text{ OR } 120 \text{ (g mol}^{-1}\text{)} \checkmark$ $\text{mol MgSO}_4 = \frac{1.51}{120.4} = 0.0125 \text{ mol } \checkmark$ | 2 | ALLOW 0.013 up to calculator value of 0.012541528 correctly rounded (from $M = 120.4 \text{ g mol}^{-1}$) ALLOW 0.013 up to calculator value of 0.012583333 correctly rounded (from $M = 120 \text{ g mol}^{-1}$) ALLOW ecf from incorrect M i.e. $1.51 \div M$ |
| | (ii) | $\frac{1.57}{18.0} = 0.0872(2) \text{ (mol)} \checkmark$ | 1 | ALLOW 0.09 up to calculator value of 0.08722222 |
| | (iii) | $x = 7 \checkmark$ | 1 | ALLOW ecf i.e. answer to (ii) \div answer to (i) ALLOW correctly calculated answer from 1 significant figure up to calculator value, ie, x does not have to be a whole number. Likely response = 6.95 ✓ |
| | | Total | 15 | |

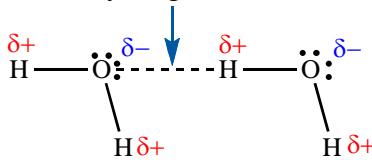
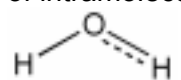
| Question | Expected Answers | Marks | Additional Guidance |
|----------|---|-------|---|
| 2 (a) |  <p>regular arrangement of labelled + ions with some attempt to show electrons ✓</p> <p>scattering of labelled electrons between other species OR a statement anywhere of delocalised electrons (can be in text below) ✓</p> <p>metallic bond as (electrostatic) attraction between the electrons and the positive ions ✓</p> | 3 | <p>Lattice must have at least 2 rows of positive ions If a metal ion is shown (e.g. Na⁺), it must have the correct charge</p> <p>ALLOW for labels: + ions, positive ions, cations If '+' is unlabelled in diagram, award the label for '+' from a statement of 'positive ions' in text below DO NOT ALLOW as label or text positive atom OR protons OR nuclei</p> <p>ALLOW e⁻ OR e as label for electron DO NOT ALLOW '- ' as label for electron</p> |
| (b) (i) | $4 \text{ Na} + \text{O}_2 \longrightarrow 2 \text{ Na}_2\text{O}$ <p>OR</p> $2 \text{ Na} + \frac{1}{2} \text{ O}_2 \longrightarrow \text{Na}_2\text{O} \checkmark$ | 1 | <p>ALLOW correct multiples including fractions IGNORE state symbols</p> |
| (ii) | (electrostatic) attraction between oppositely charged ions ✓ | 1 | |
| | | | |

| Question | Expected Answers | Marks | Additional Guidance |
|----------|--|-----------|--|
| (iii) | <p>Na shown with either 8 or 0 electrons AND O shown with 8 electrons with 6 crosses and 2 dots (or vice versa) ✓ Correct charges on both ions ✓</p> | 2 | <p>For 1st mark, if 8 electrons shown around cation then 'extra' electron(s) around anion must match symbol chosen for electrons in cation Shell circles not required</p> <p>IGNORE inner shell electrons</p> <p>ALLOW: 2[Na⁺] 2[Na]⁺ [Na⁺]₂ (brackets not required) DO NOT ALLOW [Na₂]²⁺ / [Na₂]⁺ / [2Na]²⁺ DO NOT ALLOW: [Na₂]²⁺ [Na₂]⁺ [2Na]²⁺ [Na]₂⁺</p> |
| (c) | <p>sodium is a (good) conductor because it has mobile electrons OR delocalised electrons OR electrons can move ✓</p> <p>sodium oxide does not conduct as a solid ✓</p> <p>sodium oxide conducts when it is a liquid ✓</p> <p>ions cannot move in a solid ✓</p> <p>ions can move OR are mobile when liquid ✓</p> | 5 | <p>Throughout this question, 'conducts' and 'carries charge' are treated as equivalent terms.</p> <p>DO NOT ALLOW 'free electrons' for mobile electrons</p> <p>ALLOW poor conductor OR bad conductor 'Sodium oxide only conducts when liquid' is insufficient to award 'solid conductivity' mark</p> <p>ALLOW ions are fixed in place IGNORE electrons IGNORE charge carriers</p> <p>IGNORE 'delocalised ions' or 'free ions' for mobile ions Any mention of electrons moving is a CON</p> |
| | Total | 12 | |

| Question | | | Expected Answers | Marks | Additional Guidance |
|----------|-----|------|---|----------|---|
| 3 | (a) | (i) | mol HCl = 1.50×10^{-2} ✓ volume HCl(aq) = 75.0 ✓ | 2 | ALLOW answers to 2 significant figures ALLOW ecf from wrong number of moles i.e. $\frac{\text{moles of HCl} \times 1000}{0.200}$ ALLOW one mark for 37.5 (from incorrect 1:1 ratio) |
| | | (ii) | 180 ✓ | 1 | No other acceptable answer |
| | (b) | | $\text{CaCO}_3(\text{s}) \longrightarrow \text{CaO}(\text{s}) + \text{CO}_2(\text{g})$ equation ✓ state symbols ✓ | 2 | state symbols are dependent on correct formulae of CaCO_3 , CaO and CO_2 DO NOT ALLOW the 'equation mark' if O_2 is seen on both sides (but note that the 'state symbol mark' may still be accessible) |
| | (c) | (i) | $\text{Ca}(\text{OH})_2$ ✓ | 1 | IGNORE charges, even if wrong |
| | | (ii) | $\text{Ca}(\text{NO}_3)_2$ ✓ | 1 | IGNORE charges, even if wrong |
| | | | Total | 7 | |

| Question | | Expected Answers | Marks | Additional Guidance |
|----------|---------|---|-------|---|
| 4 | (a) (i) | the energy required to remove one electron ✓ from each atom in one mole ✓ of gaseous atoms ✓ | 3 | <p>ALLOW 3 marks for: the energy required to remove one mole of electrons ✓ from one mole of atoms ✓ atoms in the gaseous state ✓</p> <p>If no definition, ALLOW one mark for the equation below, including state symbols. $X(g) \rightarrow X^+(g) + e^-$ / $X(g) - e^- \rightarrow X^+(g)$ ALLOW e for electron IGNORE state symbol for electron</p> |
| | (b) (i) | <p>outer electrons closer to nucleus OR radii decreases ✓</p> <p>nuclear charge increases OR protons increase ✓</p> <p>electrons added to the same shell OR screening OR shielding remains the same ✓</p> | 3 | <p>IGNORE 'atomic number increases' IGNORE 'nucleus gets bigger' 'charge increases' is not sufficient ALLOW 'effective nuclear charge increases' OR 'shielded nuclear charge increases'</p> <p>ALLOW shielding is similar</p> |
| | (ii) | <p>atomic radii increase OR there are more shells ✓</p> <p>there is more shielding OR more screening ✓</p> | 3 | <p>ALLOW electrons in higher energy level ALLOW electrons are further from the nucleus DO NOT ALLOW more orbitals OR more sub-shells DO NOT ALLOW different shell or new shell</p> <p>There must be a clear comparison: e.g. 'more shielding', 'increased shielding'. i.e. DO NOT ALLOW just 'shielding'. ALLOW 'more electron repulsion from inner shells'</p> |

| Question | | Expected Answers | Marks | Additional Guidance |
|----------|----------------|---|-----------|---|
| | | the nuclear attraction decreases OR Increased shielding / distance outweigh the increased nuclear charge ✓ | | Nuclear OR proton(s) OR nucleus spelt correctly ONCE ALLOW 'nuclear pull' IGNORE any reference to 'effective nuclear charge' |
| | (c) (i) | $O^+(g) \longrightarrow O^{2+}(g) + e^-$ ✓ | 1 | answer must have state symbols ALLOW e for electron ALLOW $O^+(g) - e^- \rightarrow O^{2+}(g)$ DO NOT ALLOW $O^+(g) + e^- \longrightarrow O^{2+}(g) + 2e^-$ IGNORE state symbol for electron |
| | (ii) | the O^+ ion, is smaller than the O atom OR the electron repulsion/shielding is smaller OR the proton : electron ratio in the 2+ ion is greater than in the 1+ ion ✓ | 1 | ALLOW the outer electrons in an O^+ ion are closer to the nucleus than an O atom DO NOT ALLOW 'removed from next shell down' |
| | | Total | 11 | |

| Question | | Expected Answers | Marks | Additional Guidance |
|----------|-----|---|-------|--|
| 5 | (a) | (i) number of protons (in the nucleus) ✓ | 1 | ALLOW proton number ALLOW number of protons in an atom IGNORE reference to electrons |
| | | (ii) $(1s^2)2s^22p^63s^23p^63d^24s^2$ ✓ | 1 | ALLOW $1s^2$ written twice ALLOW subscripts ALLOW $4s^2$ before $3d^2$ |
| | | (iii) Mn / manganese and d ✓ | 1 | ALLOW D |
| | (b) | (i) <p style="text-align: center;">Hydrogen bond</p>  <p>Shape of water with at least one H with $\delta+$ and at least one O with $\delta-$ ✓</p> <p>H-bond between H in one water molecule and a lone pair of an O in another water molecule ✓</p> <p>hydrogen bond labelled OR H_2O has hydrogen bonding ✓</p> | 3 | all marks can be awarded from a labelled diagram If HO_2 shown then DO NOT ALLOW 1st mark Dipole could be described in words so it does not need to be part of diagram. At least one hydrogen bond must clearly hit a lone pair Lone pair interaction could be described in words so it does not need to be part of diagram. DO NOT ALLOW hydrogen bonding if described in context of intramolecular bonding, <i>ie</i>  |
| | | (ii) no hydrogen bonding OR weaker intermolecular forces ✓ | 1 | DO NOT ALLOW 'weaker' / 'weak' hydrogen bonding ALLOW weaker van der Waals' forces ALLOW weaker dipole-dipole interactions DO NOT ALLOW 'weak intermolecular forces' (<i>ie</i> comparison essential here) DO NOT ALLOW 'no intermolecular forces' |

| Question | | Expected Answers | Marks | Additional Guidance |
|--------------|---------|--|-----------|---|
| | (c) | <p>van der Waals' forces OR induced dipole interactions ✓</p> <p>number of electrons increases ✓</p> <p>Down the group, intermolecular forces / van der Waals' forces increase OR Down the group, more energy needed to break intermolecular / van der Waals' forces ✓</p> | 3 | <p>electron(s) must be seen and spelt correctly ONCE ALLOW number of electron shells increases ALLOW iodine has most electrons ALLOW chlorine has the least electrons</p> <p>For 'Down the group' ALLOW 'Increase in boiling points' or 'Molecules get bigger'</p> |
| | (d) (i) | goes brown ✓ | 1 | ALLOW yellow OR orange OR any shade of yellow, orange and brown, e.g. reddish-brown IGNORE precipitate |
| | (ii) | <p>iodine and (potassium) chloride ✓</p> $\text{Cl}_2 + 2\text{I}^- \longrightarrow \text{I}_2 + 2\text{Cl}^- \quad \checkmark$ | 2 | DO NOT ALLOW formulae (<i>i.e.</i> names essential) ALLOW any correct multiple including fractions IGNORE state symbols |
| | (iii) | <p>chlorine / Cl_2 is more reactive (than iodine) OR chlorine / Cl_2 is a more powerful oxidising agent ✓</p> | 1 | <p>ALLOW chlorine is better at electron capture OR chlorine attracts electrons more</p> <p>ALLOW iodine is less reactive (than chlorine) ALLOW iodide (ion) / I^- is a stronger reducing agent</p> <p>DO NOT ALLOW Cl is more reactive DO NOT ALLOW explanation in terms of displacement DO NOT ALLOW chlorine is more electronegative</p> |
| | (iv) | goes purple / violet / lilac / pink ✓ | 1 | ALLOW pink OR any combination of purple, violet, lilac and pink |
| Total | | | 15 | |

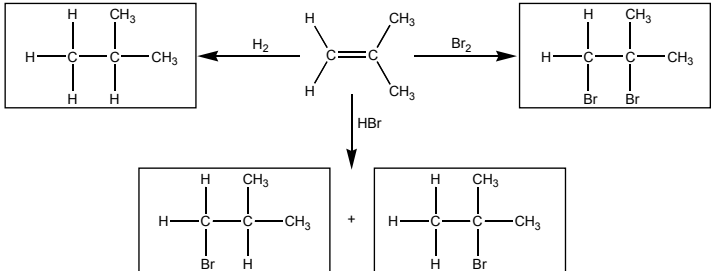
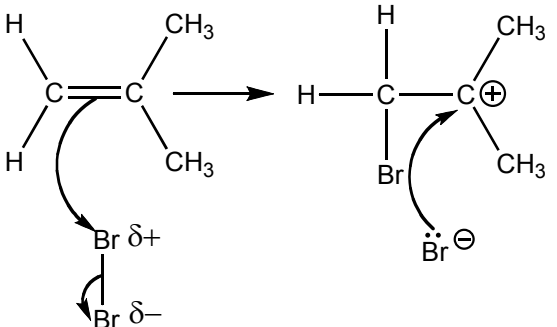
F322 Chains, Energy and Resources

| Question | | Expected Answers | Marks | Additional Guidance |
|----------|-----|--|-------|---|
| 1 | (a) | C_nH_{2n+2} ✓ | 1 | ALLOW $C_nH_{2(n+1)}$ ✓ IGNORE size of subscripts |
| | (b) | (i) $C_8H_{18} + 8\frac{1}{2}O_2 \rightarrow 8CO + 9H_2O$ ✓ | 1 | ALLOW any correct multiples IGNORE state symbols |
| | | (ii) limited supply of air OR not enough O_2 ✓ | 1 | ALLOW use of air or oxygen IGNORE it is not completely oxidised |
| | (c) | (i) $2CO + 2NO \rightarrow 2CO_2 + N_2$ ✓ | 1 | ALLOW any correct multiples including fractions IGNORE state symbols |
| | (c) | (ii) CO and NO are adsorbed (onto surface) OR reactants are adsorbed (onto surface) ✓ weakening of bonds OR lowers activation energy ✓ CO_2 and N_2 desorbs (from the surface) OR products desorbs (from the surface) ✓ | 3 | ALLOW CO and NO stick onto surface OR CO and NO form weak attractions to the surface OR gases are adsorbed onto surface NOT absorb but allow ecf for deabsorb later on IGNORE alternative pathway Requires less energy is not sufficient ALLOW products leave the surface OR products diffuse away from surface OR weak attraction to surface is broken ALLOW deadsorb |
| | (d) | skeletal formula of a branched isomer of C_8H_{18} ✓ skeletal formula of a cyclic hydrocarbon OR skeletal formula of substituted arene of C_8H_{10} ✓ | 2 | ALLOW any ring between C_3 and C_8 with 8 carbon atoms per molecule IGNORE wrong names If two correct structural or displayed formulae drawn award one mark |

| Question | | Expected Answers | Marks | Additional Guidance |
|----------|-----|--|-----------|--|
| | (e) | <p>Any TWO from: atmospheric concentration ✓ ability to absorb infrared radiation ✓ residence time ✓</p> | 2 | <p>ALLOW the amount of the gas OR abundance of gas ALLOW how much IR it absorbs OR ability to absorb heat IGNORE global warming potential / heat reflected / how much is produced ALLOW how long it stays in the atmosphere</p> |
| | | <p>Any TWO from: deep in the oceans OR on the sea-bed ✓ storage in geological formations OR under the sea-bed ✓ by reaction (with metal oxides) to form carbonates ✓</p> | 2 | <p>ALLOW piped into disused or partially filled oil wells ALLOW stored as a carbonate OR equation to show formation of suitable carbonate from an oxide IGNORE mineral storage IGNORE reforestation</p> |
| | | Total | 13 | |

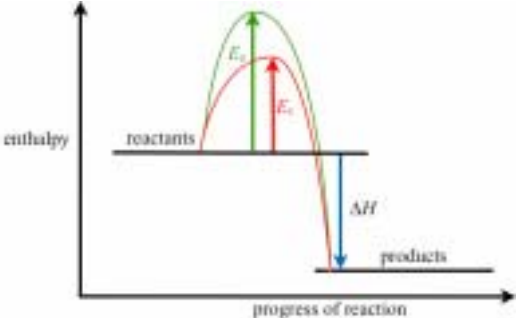
| Question | | Expected Answers | Marks | Additional Guidance |
|----------|---------|--|-------|--|
| 2 | (a) (i) | The enthalpy change for the complete combustion ✓ of 1 mol (of a substance) ✓ | 2 | ALLOW energy change for combustion in excess oxygen OR energy released during complete combustion OR energy change for combustion in excess air NOT energy required This mark is not stand alone but must relate to statement about an enthalpy change even if the statement was not awarded a mark |
| | (b) (i) | 56.430 (kJ) ✓ | 1 | ALLOW 56.43 (kJ) OR 56.4 kJ ✓ OR 56 kJ ALLOW -56.43 i.e. ignore sign |
| | (ii) | M_r [CH ₃ (CH ₂) ₄ OH] = 88.0 ✓ n = 0.0200 mol ✓ | 2 | ALLOW 88 ALLOW 0.02 OR ecf from wrong M_r ALLOW full marks for 0.02 with no working out |
| | (iii) | (-)2821.5 ✓ = (-)2820 (3 SF) ✓ correct minus sign ✓ | 3 | ALLOW correct substitution into formula (b)(i) ÷ (b)(ii) e.g. 56.4 ÷ 0.02 this is essentially a mark for the working ALLOW ecf from i.e. answer from (b)(i) ÷ (b)(ii) The minus mark is stand alone and is independent of the numerical answer |
| | (c) (i) | pressure: 100 kPa OR 101 kPa AND temperature: 298 K OR 25 °C ✓ | 1 | units needed ALLOW 1 bar OR 1 atm OR 760 mmHg ALLOW any stated temperature so for example 100kPa and 40°C would be credited with a mark IGNORE any reference to moles or concentration |
| | (ii) | 6C(s) + 7H ₂ (g) → C ₆ H ₁₄ (l) ✓ | 1 | ALLOW graphite / gr |
| | (iii) | many different hydrocarbons would form OR activation energy too high OR reaction too slow OR they don't react together ✓ | 1 | ALLOW can form different isomers OR can form different structures IGNORE reaction may be reversible |

| Question | | Expected Answers | Marks | Additional Guidance |
|----------|------|--|-----------|---|
| | (iv) | $6 \times -394 + 7 \times -286$ shown OR calculated as -4366 ✓ -4366 and -4163 added OR subtracted ✓ correct answer $-4366 - (-4163) = -203$ ✓ | 3 | ALLOW THREE marks for -203 on its own with no working out or written on the answer line ALLOW TWO marks for $+203, +3483, +1513, +1767$ or -8529 on its own with no working out ALLOW ONE mark for or $-3483, -1513, -1767$ or $+8529$ on its own with no working out units NOT needed Positive sign not needed for endothermic answers |
| | | Total | 14 | |

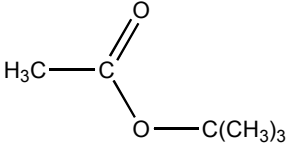
| Question | Expected Answers | Marks | Rationale |
|----------|--|-------|--|
| 3 (a) |  <p>one mark for each correct structure ✓ ✓ ✓ ✓</p> | 4 | <p>ALLOW skeletal formula OR displayed formulae IGNORE molecular formulae IF two answers given e.g. name and structure then both must be correct to be given a mark</p> <p>ALLOW methylpropane OR $(\text{CH}_3)_3\text{CH}$ ✓</p> <p>ALLOW 1,2-dibromo-methylpropane OR $\text{CH}_2\text{BrCBr}(\text{CH}_3)_2$ ✓</p> <p>ALLOW 1-bromo-methylpropane OR $\text{CH}_2\text{BrCH}(\text{CH}_3)_2$ ✓</p> <p>ALLOW 2-bromo-methylpropane OR $\text{CH}_3\text{CBr}(\text{CH}_3)_2$ ✓</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> |
| (b) | <p>curly arrow from double bond to $\text{Br}^{\delta+}$ and curly arrow from $\text{Br}-\text{Br}$ bond pair to $\text{Br}^{\delta-}$ in 1st step ✓</p> <p>curly arrow in 2nd step from bromide ion ✓</p> <p>correct dipole shown on Br_2 ✓</p> <p>correct carbocation shown ✓</p>  | 4 | <p>Curly arrow must start from the double bond and not a carbon atom, other curly arrow must start from $\text{Br}-\text{Br}$ bond</p> <p>ALLOW curly arrow from any part of bromide ion The bromide 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. $^+\text{CH}_2\text{CBr}(\text{CH}_3)_2$</p> |

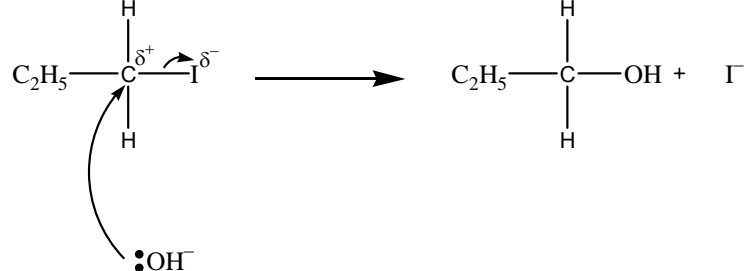
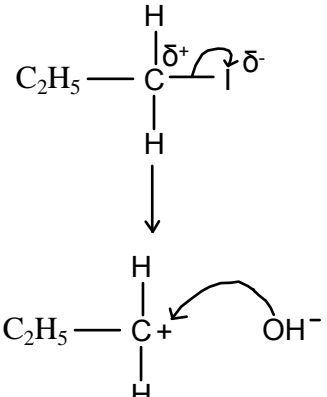
| Question | | Expected Answers | Marks | Rationale |
|----------|---------|---|-----------|---|
| | (c) (i) | C_6H_{10} ✓ | 1 | |
| | (ii) | $M_r(\text{cyclohexanol}) = 100$ ✓ amount of cyclohexanol = 0.0765 mol ✓ percentage yield = 35.0% ✓ | 3 | ALLOW full marks for correct answer with no or limited working out ALLOW ecf from wrong molar mass i.e. $7.65 \div$ molar mass ALLOW ecf from wrong amount in moles i.e. $[0.0268 \div \text{moles}] \times 100$ ALLOW 35% ALLOW two marks for 0.35% If M_r of 82 is used then % yield will be 28.7 or 29 and this is worth two marks |
| | (d) (i) | (sum of) the molecular masses of the desired product \div sum of molecular masses of all products $\times 100$ ✓ | 1 | ALLOW (sum of) the molecular masses of the desired product \div sum of molecular masses of all reactants $\times 100$ ✓ |
| | (ii) | this preparation is addition OR has 100% atom economy OR there is only one product ✓ preparation from cyclohexanol has less than 100% atom economy OR H_2O is produced as well OR calculated atom economy = 82% ✓ | 2 | ALLOW no by products formed ALLOW other substances formed OR cyclohexene is not the only product |
| | | Total | 15 | |

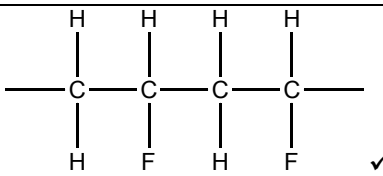
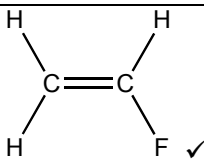
| Question | | Expected Answers | Marks | Additional Guidance |
|----------|---------|---|-------|--|
| 4 | (a) | <p>high pressure as fewer moles (of gas) on right-hand side OR high pressure as volume of products less than that of reactants ✓</p> <p>low temperature as (forward) reaction is exothermic ✓</p> | 2 | <p>ALLOW ora ALLOW fewer particles OR fewer molecules</p> <p>ALLOW ora</p> |
| | (b) | <p>Too expensive to use a high pressure ✓</p> <p>Too slow to use a low temperature ✓</p> | 2 | <p>ALLOW high pressures provide a safety risk OR high pressure is too dangerous</p> <p>ALLOW with low temperature molecules cannot overcome activation barrier</p> |
| | (c) (i) | <p>Cl + O₃ → ClO + O₂ ✓ ClO + O → Cl + O₂ ✓ overall: O₃ + O → 2O₂ ✓</p> <p>OR</p> <p>Cl + CH₄ → CH₃ + HCl ✓ CH₃ + Cl₂ → CH₃Cl + Cl ✓ overall: CH₄ + Cl₂ → CH₃Cl + HCl ✓</p> | 3 | <p>Marks must come from one or other of the radical process and not from both of them. If two processes are described then an incorrect step in one process will contradict a correct step in the other process.</p> <p>ALLOW overall equation mark even if the steps are wrong the radicals do NOT need a single dot IGNORE any state symbols</p> <p>ALLOW Cl + O₃ → ClO + O₂ ✓ ClO + O₃ → Cl + 2O₂ ✓ overall: 2O₃ → 3O₂ ✓</p> <p>ALLOW any saturated hydrocarbon including cyclic ALLOW ecf for second step and overall reaction if wrong hydrocarbon used e.g. C₂H₄ is used in first step</p> |

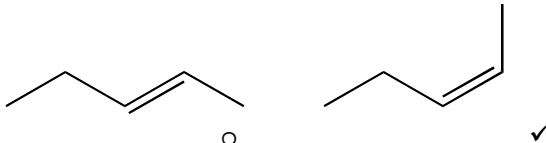
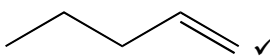
| Question | Expected Answers | Marks | Additional Guidance |
|----------|---|-----------|---|
| (ii) | <p>ΔH shown and products below reactants ✓</p> <p>E_a shown ✓</p> <p>E_c shown $< E_a$ ✓</p>  | 3 | <p>NOT double headed arrows but apply ecf for more than one double headed arrow</p> <p>ALLOW one mark if two correctly labelled curves are drawn but the arrows are not shown or are incorrectly drawn</p> <p>The arrows must be positioned as closely as possible to the maximum height of the curves but allow some degree of bod</p> |
| (d) | <p>Any FOUR from:</p> <p>catalyst not used up in reaction ✓</p> <p>reactions take place at lower temperatures ✓</p> <p>with lower energy demand OR lower activation energy OR use less fuel ✓</p> <p>so less carbon dioxide emitted into atmosphere OR so fossil fuels last longer ✓</p> <p>different reactions can be used ✓</p> <p>with better atom economy OR less waste ✓</p> <p>less hazardous chemicals ✓</p> <p>catalysts or enzymes can generate specific products ✓</p> | 4 | <p>ALLOW catalysts can work at room temperature OR enzymes work at room temperature</p> <p>IGNORE cheaper</p> |
| | Total | 14 | |

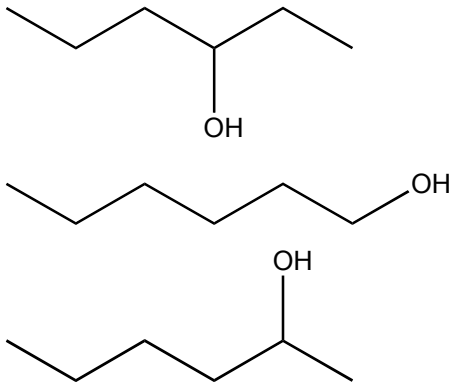
| Question | | Expected Answers | Marks | Additional Guidance |
|----------|-----|---|-------|--|
| 5 | (a) | <p>method 1: fermentation of sugars or carbohydrates OR reaction with yeast with sugar or carbohydrates ✓ $C_6H_{12}O_6 \rightarrow 2C_2H_5OH + 2CO_2$ ✓</p> <p>method 2: hydration of ethene OR reaction of ethene with water OR reaction of steam with ethene ✓ $C_2H_4 + H_2O \rightarrow C_2H_5OH$ ✓</p> | 4 | <p>ALLOW sugar from equation</p> <p>ALLOW C_2H_6O in equation ALLOW correct multiples IGNORE state symbols</p> <p>ALLOW ethene from the equation IGNORE mention of any catalyst ALLOW C_2H_6O in equation OR H_2O over the arrow ALLOW correct multiples IGNORE state symbols</p> |
| | (b) | (i) | 2 | <p>If name and formula given both need to be correct ALLOW propanone OR acetone IGNORE propone NOT incorrect named compound</p> <p>ALLOW $C_3H_8O + [O] \rightarrow C_3H_6O + H_2O$ ALLOW O instead of [O] ALLOW correct multiples IGNORE state symbols</p> |
| | | (ii) | 3 | <p>ALLOW $C=O$ and $O-H$ marks independent of compound identified i.e. stand alone marks ALLOW correct bonds shown by the appropriate absorption on the IR spectrum IGNORE reference to $C-O$ bond</p> |
| | (c) | (i) | 1 | ALLOW methylpropan-2-ol OR tertiarybutanol |

| Question | Expected Answers | Marks | Additional Guidance |
|----------|--|-----------|--|
| | (ii) ester ✓ | 1 | |
| | (iii) $\text{CH}_3\text{CO}_2\text{C}(\text{CH}_3)_3$ OR $\text{CH}_3\text{COOC}(\text{CH}_3)_3$ OR  ester group shown ✓ rest of molecule ✓ | 2 | ALLOW skeletal formula OR displayed formula ALLOW ester linkage even if rest of structure is wrong |
| | Total | 13 | |

| Question | Expected Answers | Marks | Additional Guidance |
|-----------|---|-------|--|
| 6 (a) (i) |  <p>C-I curly arrow from the bond not from carbon atom ✓</p> <p>curly arrow from the OH⁻ ✓</p> <p>correct partial charges on C—I ✓</p> | 3 | <p>no need to show any lone pairs on oxygen but must have a clear negative sign rather than partial negative charge IGNORE lone pairs IGNORE products of this reaction</p> <p>ALLOW curly arrow from a negative charge or from any part of hydroxide ion</p> <p>If S_N1 mechanism is given then use the mark scheme below</p> <p>correct partial charges on C—I ✓</p> <p>C—I curly arrow from the bond not from carbon atom ✓</p> <p>curly arrow from the OH⁻ to the correct carbocation ✓</p>  |
| | (ii) nucleophilic substitution ✓ | 1 | |
| (b) | <p>C—I bonds broken more easily ✓</p> <p>C—I bonds are weaker OR have less bond enthalpy OR C—I bonds are longer ✓</p> | 2 | <p>ALLOW ora e.g. C—Br bonds are stronger OR broken less easily</p> |

| Question | | Expected Answers | Marks | Additional Guidance |
|----------|---------|---|-----------|--|
| | (c) | <p>Any TWO from: CFCs take many years to reach the ozone layer OR long residence time ✓</p> <p>CFCs are still being used ✓</p> <p>there are other ozone depleting substances ✓</p> | 2 | <p>IGNORE because chlorine radicals stay in the stratosphere</p> <p>ALLOW other named ozone depleting substances e.g. NO and HFCs</p> |
| | (d) (i) |  | 1 | <p>Free bonds at bond ends must be present</p> <p>ALLOW minor slip e.g. missing one hydrogen and left as a stick</p> <p>ALLOW more than two repeat units but must be a whole number of repeat units</p> <p>IGNORE brackets, use of numbers and n in the drawn structure</p> |
| | (ii) |  | 1 | <p>ALLOW skeletal formula</p> <p>ALLOW CH₂CHF</p> |
| | (e) | <p>Any two from: separation into types and recycling OR sort plastics, melt and remould ✓</p> <p>combustion for energy generation ✓</p> <p>used for cracking OR feedstock for plastics or chemicals ✓</p> | 2 | <p>IGNORE biodegradable</p> <p>used as a fuel is insufficient releases energy is insufficient</p> <p>ALLOW burning plastics to release energy</p> <p>ALLOW organic feedstock / raw materials to make organic compounds</p> |
| | | Total | 12 | |

| Question | Expected Answers | Marks | Additional Guidance |
|----------|---|-------|--|
| 7 (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 ✓ 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), M_r = 102 ✓</p> <p>evidence that it has been shown that the empirical formula is the molecular formulae e.g. M_r 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 | |

Grade Thresholds

Advanced GCE (Chemistry A) (H034 H434)
June 2009 Examination Series

Unit Threshold Marks

| Unit | | Maximum Mark | a | b | c | d | e | u |
|------|-----|--------------|-----|-----|----|----|----|---|
| F321 | Raw | 60 | 50 | 43 | 37 | 31 | 25 | 0 |
| | UMS | 90 | 72 | 63 | 54 | 45 | 36 | 0 |
| F322 | Raw | 100 | 75 | 65 | 55 | 46 | 37 | 0 |
| | UMS | 150 | 120 | 105 | 90 | 75 | 60 | 0 |
| F323 | Raw | 40 | 34 | 31 | 28 | 25 | 22 | 0 |
| | UMS | 60 | 48 | 42 | 36 | 30 | 24 | 0 |

Specification Aggregation Results

Overall threshold marks in UMS (ie after conversion of raw marks to uniform marks)

| | Maximum Mark | A | B | C | D | E | U |
|------|--------------|-----|-----|-----|-----|-----|---|
| H034 | 300 | 240 | 210 | 180 | 150 | 120 | 0 |

The cumulative percentage of candidates awarded each grade was as follows:

| | A | B | C | D | E | U | Total Number of Candidates |
|------|------|------|------|------|------|-------|----------------------------|
| H034 | 17.6 | 35.1 | 52.8 | 68.8 | 82.2 | 100.0 | 16327 |

16327 candidates aggregated this series

For a description of how UMS marks are calculated see:

http://www.ocr.org.uk/learners/ums_results.html

Statistics are correct at the time of publication.

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