# Chemistry B (Salters) 

Advanced GCE A2 H435
Advanced Subsidiary GCE AS H035

## Mark Schemes for the Units

## January 2010

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## F331 Chemistry for Life

| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (a) | (i) | skeletal $\checkmark$ | 1 | ALLOW recognisable spellings |
|  |  | (ii) | 2,2,4-trimethylpentane $\checkmark \checkmark$ | 2 | IGNORE gaps, dashes, hyphens, commas <br> pentane $\checkmark$ <br> rest $\checkmark$ |
|  |  | (iii) | ring structure / arene / cyclic OR short(er) molecule $\checkmark$ | 1 | ALLOW small |
|  | (b) | (i) | burn measured mass / amount of fuel / octane <br> measure temp rise <br> of a fixed volume / mass / amount of water $\downarrow$ <br> use - energy transferred = mass of water x specific heat capacity (of water) $x$ temp rise <br> scale up to one mole of fuel / octane used / AW | 5 | ALLOW measure starting and finishing temperature / mass of octane / fuel DO NOT ALLOW just 'final' temp. recorded <br> IGNORE reference to solution <br> ALLOW q / energy $=m c \Delta T$ or mc $\theta$ allow ' $m$ ' as mass of water unless conned eg ALLOW answer divided by moles burnt |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
|  | (ii) | any two from 4: <br> heat loss to surroundings / air / effect of draughts; etc <br> heat losses to calorimeter / apparatus; <br> incomplete combustion of fuel / lack of (enough) oxygen; <br> evaporation of fuel (from wick); | 2 | DO NOT ALLOW 'not standard conditions' / reference to data book values / AW <br> DO NOT ALLOW 'enthalpy may escape’ <br> IGNORE evaporation of water / measurement error / human error |
| (c) | (i) | $\Delta \mathrm{H}_{1}=$ enthalpy (change) of formation (of octane) <br> $\Delta \mathrm{H}_{2}=$ enthalpy (change) of combustion of eight moles of carbon / (enthalpy (change) of formation of eight moles of carbon dioxide) <br> $\Delta \mathrm{H}_{3}=$ enthalpy (change) of combustion of nine moles of hydrogen / (enthalpy (change) of formation of nine moles of water) <br> $\Delta \mathrm{H}_{4}=$ enthalpy (change) of combustion of octane $\checkmark$ | 4 | ALLOW omission of the words 'enthalpy change of...' <br> IGNORE references to oxygen <br> ALLOW appropriate symbols eg $\Delta \mathrm{H}_{\mathrm{f}}$ <br> ALLOW $\Delta H_{2}$ and $\Delta H_{3}$ in either order. <br> Score one out of two if numbers of moles not mentioned <br> ALLOW $\Delta \mathrm{H}_{2} / \Delta \mathrm{H}_{3}$ in terms of enthalpy changed of formation of 8 moles $\mathrm{CO}_{2}$ and 9 moles of $\mathrm{H}_{2} \mathrm{O}$. <br> DO NOT ALLOW any rearrangement of $\Delta \mathrm{H}_{1}$ etc |
|  | (ii) | answer $=-248 \checkmark$ | 1 |  |
|  |  | Total | 16 |  |



| Question |  | Expected Answers | Marks | Additional Guidance |
| :--- | :--- | :--- | :---: | :--- |
|  | (c) | (i) | $\begin{array}{l}\text { difficult to detect very small amounts of Ar-40 formed K-40 } \\ \text { decayed / dating errors very large when so little decay has } \\ \text { taken place / AW } \checkmark\end{array}$ | $\mathbf{1}$ |
| (ii) | $\begin{array}{l}\text { Ar }{ }^{+} \text {(allow Ar } r^{2+} \text { ) } \checkmark \\ \text { 'not even fLLOW answers that talk only in terms of one half life' }\end{array}$ |  |  |  |
|  | (iii) | $\begin{array}{l}\text { peak / bar / line at (mass numbers) 36, 38 and 40 } \checkmark \\ \text { size / height of peak related to abundance } \checkmark\end{array}$ | $\mathbf{1}$ | $\begin{array}{l}\text { ALLOW with correct mass / atomic numbers } \\ \text { (iv) }\end{array}$ |
| $\begin{array}{l}\text { (energy lost as) electrons go from higher to lower levels } \checkmark \\ \text { relationship of energy to frequency / wavelength } \checkmark \\ \text { gives a (specific) line(s) } \checkmark \\ \text { energy gaps / levels different for different elements } \checkmark \\ \text { QWC - wavelength / frequency / frequencies must be } \\ \text { spelled correctly }\end{array}$ | $\mathbf{4}$ | mass numbers needed to score |  |  |$]$| eg E = hf or in words |
| :--- |
| mention of lines scores a mark |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| Q | (a) | (i) | (hydrocarbon) contains no benzene rings / not an arene $\checkmark$ | 1 | DO NOT ALLOW contains no rings |
|  |  | (ii) | fractional distillation $\checkmark$ | 1 |  |
|  |  | (iii) | $\mathrm{C}_{25} \mathrm{H}_{52}+38 \mathrm{O}_{2} \rightarrow 25 \mathrm{CO}_{2}+26 \mathrm{H}_{2} \mathrm{O} \checkmark$ | 1 |  |
|  | (b) | (i) | unburnt hydrocarbon / $\mathrm{C}_{25} \mathrm{H}_{52} \checkmark$ | 1 | ALLOW paraffin wax ALLOW CO <br> ALLOW smaller hydrocarbon |
|  |  | (ii) | carbon monoxide $\checkmark$ carbon / soot | 2 | ALLOW water IGNORE oxides of nitrogen |
|  | (c) | (i) | $\mathrm{C}_{3} \mathrm{H}_{6} \checkmark$ | 1 | order of elements immaterial |
|  |  | (ii) | $110-130^{\circ}$ <br> 3 areas of electron density <br> around central C <br> areas of electron density / pairs repel as far apart as possible / minimize energy $\downarrow$ | 4 | DO NOT ALLOW 3 'atoms' or 'electron pairs' ALLOW names or descriptions of electron groups eg double bond <br> ALLOW clear diagram or description <br> DO NOT ALLOW repel as much as possible <br> TAKE CARE repel and 'as far apart' run together for only one mark <br> ALLOW bonds (but not atoms) repel |
|  |  | (iii) | catalysts and reactants in different (physical) states $\checkmark$ | 1 |  |
|  |  | (iv) | contain hole(s) / channels / porous / gaps / rings $\checkmark$ can trap branched / let through straight isomers $\checkmark$ | 2 |  |
|  |  |  | Total | 14 |  |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 4 (a) |  | $\begin{aligned} & \text { mass number }=1 \mathrm{v} \\ & \text { atomic number }=0 \end{aligned}$ | 2 |  |
| (b) | (i) | $\begin{aligned} & \text { moles of } \mathrm{Be}=1.75 / 9(0.19) \checkmark \\ & \text { moles of } \mathrm{Cu}=98.25 / 63.5(1.55) \end{aligned}$ | 2 | all usual ecf's apply (allow working to more / less sig. figs.) <br> Max 1 if unit other than moles put in |
|  | (ii) | 11 scores all three total no. of moles $=1.74$ $\% \mathrm{Be}=0.19 / 1.74 \times 100 \checkmark(=10.919)$ <br> Sig. figs. separate mark based on a followable calculation | 3 | ALLOW ecf's from (b)(i) <br> ALLOW sig. figs. mark for a (wrong) calculation based on some given figures |
| (c) |  | Delocalised electrons <br> Regular array of cations / positive ions / residues $\checkmark$ <br> Labels but any used must be correct $\checkmark$ | 3 | First two points can be on diagram or labels minimum of five cations shown (can touch) <br> ALLOW positive atoms DO NOT ALLOW positive nucleus or positive metal |
| (d) |  |  <br> 'correct' pairs on Be $\checkmark$ <br> 3 pairs on $\mathrm{Cl} \checkmark$ | 2 | DO NOT ALLOW ionic structure |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :--- | :---: | :--- |
| (e) | melting point (is different) $\checkmark$ <br> (melting point) is higher in ionic compounds $\checkmark$ | $\mathbf{3}$ | ALLOW ORA throughout <br> AND <br> Any one of: <br> ionic compounds conduct electricity when in solution / molten $\checkmark$ <br> OR <br> ionic compounds (generally) water soluble / ora / AW $\checkmark$ <br> ALLOW bases liquids <br> Must be a comparison for $2^{\text {nd }}$ mark |  |

## F332 Chemistry of Natural Resources

| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (a) |  | Incomplete combustion $\checkmark$ <br> of hydrocarbons | 2 | ALLOW not enough oxygen or air linked to the idea of combustion / uncomplete combustion <br> Second mark depends on the first. ALLOW fossil fuel or named fossil fuel / carbon in the fuel / organic fuel <br> DO NOT ALLOW just 'fuel' or carbon as the fuel |
|  | (b) |  | Toxic / poisonous / reduces the capacity of blood to carry oxygen around the body / AW <br> AND <br> Any one from: <br> causes (photochemical) smog <br> oxidised to $\mathrm{CO}_{2}$ which is a greenhouse gas / reacts with $\mathrm{O}_{2}$ to form $\mathrm{CO}_{2}$ which is a greenhouse gas | 2 | ALLOW respiratory problems, but not breathing problems. IGNORE harmful / dangerous <br> Answer must have the $\mathrm{CO}_{2}$ AND the greenhouse gas for this alternative. <br> ALLOW global warming instead of greenhouse gas. |
|  | (c) | (i) | Homolytic (fission) / homolysis $\checkmark$ | 1 | IGNORE 'photochemical dissociation' |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
|  | (ii) | $464 \times 1000$ <br> Energy value/6.02 $\times 10^{23}$ AND a correct evaluation (= $\left.7.71 \times 10^{-19} \mathrm{~J}\right) \checkmark$ | 2 | One mark is for converting from kJ to J (ie: multiplying by 1000) <br> The other is for dividing their energy value by $6.02 \times 10^{23}$ (the Avogadro constant) <br> ALLOW 2 or more sig. figs. but rounding must be correct. <br> In order to score the second mark, there must be a correct evaluation of their expression. <br> A completely correct answer on its own scores both marks. |
|  | (iii) | Answer to (c)(ii)/6.63 $\times 10^{-34} \checkmark$ $=1.16 \times 10^{15}$ <br> 3 sig. fig. | 3 | DO NOT ALLOW the second mark for evaluating any other expression (eg: answer to (c)(ii) $\times 6.63 \times 10^{-34}$ ) <br> ALLOW sig. fig. mark for any 3 sig. fig. answer that follows from any calculation (even if their evaluation of their calculation is incorrect). <br> A completely correct answer on its own scores all marks, including the sig. fig. mark. |
| (d) | (i) | (A particle) with one (or more) unpaired electron(s). $\checkmark$ | 1 | Answer must be in the context of an electron as part of some sort of particle. <br> IGNORE 'free' or 'lone' or single electron. |
|  | (ii) | bond electrons $\checkmark$ <br> rest of structure | 2 | Any symbols can be used to represent the electrons (including the same symbol for all electrons). <br> Candidate does not have to draw circles for electron shells. <br> Non-bonding electrons do not have to be shown in pairs. <br> It MUST be clear that a pair of electrons (with any symbols) is being shared between the H and the O for the first mark. <br> IGNORE any inner electron shells. |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
|  | (iii) | propagation <br> one radical is used and replaced by another / AW $\checkmark$ | 2 | ALLOW there is a radical on both sides of the equation. <br> Mark independently. |
| (e) |  | $\mathrm{SiO}_{2}$ : giant covalent / network solid / lattice / whole structure held together by covalent bonds / diagram <br> $\mathrm{CO}_{2}$ : simple molecular / molecules / $\mathrm{O}=\mathrm{C}=\mathrm{O} / \mathrm{AW} \checkmark$ <br> comparison of forces: weak intermolecular bonds (or forces) in $\mathrm{CO}_{2}$ / less energy needed to separate molecules / bonds in $\mathrm{SiO}_{2}$ are stronger than $\mathrm{CO}_{2}$ intermolecular bonds (or forces) | 3 | IGNORE 'intermolecular bonds' in $\mathrm{SiO}_{2}$ / giant molecule / giant structure Marks can be given for a labelled / annotated diagram <br> IGNORE 'covalent'. <br> Any type of intermolecular bonds can be named and can be abbreviated. <br> It must be clear that the intermolecular bonds in $\mathrm{CO}_{2}$ are being discussed, not the covalent bonds. |
| (f) | (i) | $0.008 / 8 \times 10^{-3} \checkmark$ | 1 |  |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
|  | (ii) | Any four points from: <br> 1 Sun emits UV <br> 2 Earth absorbs some of the energy (from the Sun) / heats up $\checkmark$ <br> 3 Earth radiates emits / re-emits IR <br> $4 \quad\left(\mathrm{CO}_{2}\right)$ absorbs IR radiation $\checkmark$ <br> 5 making bonds vibrate (more) <br> 6 turned into kinetic energy that raises the temperature / transfers kinetic energy to thermal energy or heat or it warms the atmosphere or Earth. <br> 7 some $\mathrm{CO}_{2}$ molecules radiate IR (which warms Earth) $\checkmark$ <br> AND <br> more $\mathrm{CO}_{2}$ molecules means more radiation is absorbed / more $\mathrm{CO}_{2}$ means greater temperature increase / enhancing the greenhouse effect / causing global warming / warming the atmosphere / Earth / planet more $\checkmark$ <br> QWC - mark for connection of ideas: idea of linking IR absorbtion to vibrations of bonds / increase in temperature (marking point 4 linked to 5 or 6) $\checkmark$ | 6 | IGNORE other types of radiation from the Sun. <br> DO NOT ALLOW Earth reflects IR in point 3. <br> Award marks for points 5 and 6 if the wrong frequency range of radiation is given as being absorbed in 4 . (eg candidate states $\mathrm{CO}_{2}$ absorbs UV). |
| (g) | (i) | aldehyde(s) $\checkmark$ | 1 | ALLOW alkanal(s) |
|  | (ii) | $\mathrm{CO}+\mathrm{C}_{2} \mathrm{H}_{4}+\mathrm{H}_{2} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CHO} \checkmark \checkmark$ | 2 | ALLOW $\mathrm{C}_{3} \mathrm{H}_{6} \mathrm{O}$ or full structural formula for propanal. <br> Completely correct scores both marks. <br> Correct formula for ethane / correctly identifies $\mathrm{H}_{2}$ as the additional reagent scores one mark. |




| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
| (f) |  | Any two from: <br> 1. chloromethane is not broken down / unreactive in the troposphere / lower atmosphere <br> 2. but is broken down / photodissociated (in the stratosphere) / AW by $\checkmark$ <br> 3. high energy UV / high frequency UV <br> 4. (breakdown of chloromethane) producing chlorine atoms / chlorine radicals <br> AND <br> (products of chloromethane) catalyse ozone breakdown / AW $\checkmark$ <br> $\mathrm{C}-\mathrm{Br}$ bond is weaker (than $\mathrm{C}-\mathrm{C} /$ ) ORA <br> so can be broken in the troposphere / molecule reacts in the troposphere / reacts before reaching the stratosphere $\checkmark$ | 5 | ALLOW 'radiation' for 'UV' <br> Points 2 and 4 can be scored from a reaction equation. <br> QWC: To gain this mark, candidate must use the word catalyst or a derivative of it, spelled correctly and used in a grammatically correct way (eg: do not award for 'it catalyse the breakdown of ozone'). <br> ALLOW 'catalyze’. |
| (g) | (i) | (concentration) values were low $\checkmark$ | 1 | Answers need to show that values were less and not just different from the expected ones. |
|  |  | Total | 15 |  |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (a) | (i) | addition $\checkmark$ | 1 | DO NOT ALLOW additional. |
|  |  | (ii) | propene $\checkmark$ | 2 | ALLOW prop-1-ene DO NOT ALLOW prop-2-ene <br> Mark independently. No ecf for the second mark. |
|  | (b) | (i) | bromine (water) $\checkmark$ | 1 | ALLOW Br 2 |
|  |  | (ii) | (from) brown / orange / yellow $\checkmark$ <br> (to) colourless | 2 | ALLOW any combination of these colours, but no others for the first mark (eg no mark for red / brown). <br> DO NOT ALLOW clear for the second answer |
|  | (c) |  |   <br> (Z) <br> (E) | 2 | Name and structure required for the mark in each case <br> Correct structures with names swapped round scores 1 mark. <br> Diagrams do not have to show correct bond angles. <br> A correct representation of but-2-ene scores 1. |
|  | (d) |  | instantaneous (dipole) - induced dipole $\checkmark$ | 1 | ALLOW temporary dipole-induced temporary dipole / van der Waals forces |
|  | (e) | (i) | Iow flexibility / resistant to chemical attack / does not react with water / unreactive / not prone to stress fractures / high tensile strength / abrasion resistant / impermeable / insoluble / rigid $\checkmark$ | 1 | IGNORE strong, hard, durable, tough, malleable, dense, high melting point, can be moulded or remoulded. <br> ALLOW waterproof or 'will not wear away'. |
|  |  | (ii) | bags $\checkmark$ | 1 | IGNORE food wrap / cling film / packaging. |
|  |  |  | Total | 11 |  |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (a) | (i) |  <br> hydrogen bond between correct atoms <br> lone pair on relevant O in line with H bond partial charges shown, $\delta$ - on each O and $\delta+$ on each H $\mathrm{O}-\mathrm{H}-\mathrm{O}$ straight $\checkmark$ | 4 | Hydrogen bond can be shown in other forms, but not as a solid line. <br> Second mark, but NOT third mark, can be scored if the hydrogen bond is between incorrect atoms. |
|  |  | (ii) | Any three from: <br> 1. intermolecular bond in propene is instantaneous dipole-induced dipole <br> 2. hydrogen bonds / intermolecular bonds (in propan-1ol) are stronger than those in propene (ORA) <br> 3. intermolecular bonds must be broken for the liquid to boil <br> 4. more energy is needed to break them (ORA) $\checkmark$ <br> AND <br> QWC - mark for connection of ideas: idea of linking strength of intermolecular bonds to amount of energy needed to break them | 4 | ALLOW van der Waals' <br> DO NOT ALLOW harder / easier <br> DO NOT ALLOW 'higher temperature' for 'more energy'. |


| Question |  | Expected Answers |  | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| (b) |  | Elimination $\checkmark$ |  | 1 | ALLOW any indication of chosen answer (eg: circling). <br> DO NOT ALLOW the mark if more than one answer has been chosen. |
| (c) |  | reagent conditions <br> sulfuric / phosphoric acid $\checkmark$ heat / reflux $\checkmark$ <br>  concentrated $\checkmark$$\|$OR heat $\checkmark$ <br> alumina / silica / pumice / <br> porous pot $\checkmark$ with (propan-1-ol) vapour $\checkmark$ |  | 3 | ALLOW correct formula for reagent. <br> ALLOW temperatures over $100^{\circ} \mathrm{C}$ for the heat mark <br> Sulfuric acid AND alumina: CON reagent mark (but can still score condition marks). <br> Clear alternatives (ie: sulfuric acid OR alumina) scores the mark. <br> ALLOW c. for concentrated. <br> Aqueous / water CONs the concentrated mark. <br> The conditions marks may only be awarded if candidate has written an appropriate reagent, even if they have made a small mistake, eg: sulfuric without acid, or wrong formula (like AlO) <br> (Concentrated) sulfuric acid with dichromate and heat scores zero. <br> IGNORE references to pressure conditions. |
| (d) |  | rate of forward reaction = rat <br> concentrations of reactants / closed system | of back reaction d products remain constant | 2 | IGNORE references to steady state. |
| (e) | (i) | amount of propene produced (increased pressure) pushes left/to the reactants / side with | decreases <br> (position of) equilibrium to the fewest molecules | 2 | MUST mention equilibrium for the second mark. <br> Mark independently. |
|  | (ii) | amount of propene produced (increased temperature) push the endothermic direction / to | ncreases $\checkmark$ <br> (position of) equilibrium in he right / to the products $\checkmark$ | 2 | MUST mention equilibrium for the second mark. <br> Mark independently. |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
| (f) |  | Any three from: <br> 1. increased pressure increases number of particles per unit of volume <br> 2. more collisions occur <br> 3. (more collisions) per unit of time <br> 4. rate increases/gets faster | 3 | ALLOW 'particles are closer together' for the first point DO NOT ALLOW 'reactants are closer together'. <br> More frequent collisions / collisions occur more often covers two points IGNORE more likely to collide / greater chance of collisions in point 2. |
| (g) | (i) |  | 1 | ALLOW any clear representations of a structural formula, eg: $\mathrm{CH}_{3} \mathrm{CHBrCH}_{3}$ |
|  | (ii) |  | 1 | ALLOW CH3 $\mathrm{CH}_{2} \mathrm{CH}_{3}$ |
| (h) |  | platinum $\checkmark$ | 1 | ALLOW Pt. |
|  |  | Total | 10 |  |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 5 | (a) | Chemical that: causes another chemical to be oxidised / is itself reduced / decreases in oxidation state / is an electron acceptor / removes electrons from another chemical <br> $\mathrm{O}_{2} /$ (potassium) manganate(VII) <br> AND either <br> (Oxidises) iron from oxidation state $+2 / \mathrm{Fe}(\mathrm{II})$ <br> to $+3 / \mathrm{Fe}(\mathrm{III})$ <br> OR <br> (Manganese reduced) from $\mathrm{Mn}(\mathrm{VII}) /+7 /$ manganate(VII) $\downarrow$ <br> to $\mathrm{Mn}(\mathrm{IV}) /+4 /$ manganese(IV) oxide $\checkmark$ | 4 | ALLOW 'chemical that oxidises another chemical' / oxidising agent. <br> IGNORE references to change in pH . <br> ALLOW permanganate / $\mathrm{MnO}_{4}{ }^{-} / \mathrm{KMnO}_{4}$ <br> $\mathrm{Fe}^{2+}$ to $\mathrm{Fe}^{3+}$ scores 1. |
|  | (b) | $\mathrm{Al}^{3+}(\mathrm{aq})+3 \mathrm{HCO}_{3}^{-}(\mathrm{aq}) \rightarrow \mathrm{Al}(\mathrm{OH})_{3}(\mathrm{~s})+3 \mathrm{CO}_{2}(\mathrm{~g} / \mathrm{aq})$ <br> Correct species <br> Balanced <br> State symbols | 3 | Second and third marks depend on the first. <br> ALLOW Al ${ }^{3+}(\mathrm{aq})+\mathrm{HCO}_{3}^{-}(\mathrm{aq}) \rightarrow \mathrm{Al}(\mathrm{OH})_{3}(\mathrm{~s})$ for one mark, if no other mark is scored (IGNORE any other chemicals) |
|  | (c) | Calcium hydroxide / calcium oxide $\checkmark$ <br> Sodium carbonate <br> Hydrogencarbonate | 3 | IGNORE a correct oxidation state for Ca and Na <br> ALLOW 'hydrogen carbonate' but NOT 'bicarbonate' |
|  | (d) | Calcium ions more highly charged or more positive (than sodium ions) / mention of $\mathrm{Ca}^{2+}$ and $\mathrm{Na}^{+}$ <br> so are more strongly attracted to the negative charge on the resin / (R)COO ${ }^{-} /$anion groups $\checkmark$ | 2 | IGNORE references to reactivity. <br> DO NOT ALLOW just 'attracted to the resin'. |


| Questi | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| (e) | $\begin{aligned} & \text { moles } \mathrm{Ca}^{2+}=(800 / 1000) \times 0.002(=0.0016) \checkmark \\ & \text { moles } \mathrm{Na}^{+}=2 \times \text { moles } \mathrm{Ca}^{2+}(=0.0032) \checkmark \\ & \text { mass } \mathrm{Na}^{+}=\text {moles } \mathrm{Na}^{+} \times 23=(0.0032 \times 23=0.0736 / 0.074) \end{aligned}$ $(\mathrm{g})^{\checkmark}$ | 3 | Mass $\mathrm{Na}^{+}=0.0368 / 0.037$ scores 2. |
| (f) | Any five points from: <br> 1. kills bacteria / kills pathogens / disinfectant <br> 2. cheap compared to other water treatment chemicals. <br> 3. $\mathrm{Cl}_{2}$ or chlorine is a gas, making it difficult to contain / it spreads easily. <br> 4. toxic / poisonous <br> 5. causes respiratory problems / breathing problems <br> 6. forms by-products / THMs that are suspected carcinogens <br> 7. dissolves in rivers / local water supplies <br> 8. forming bleach and acid <br> 9. (bleach and acid) kill life forms in the water | 5 | DO NOT ALLOW just 'cheap'. <br> Answer must have 'gas' and either 'difficult to contain' or 'spreads easily' to gain the mark. <br> ALLOW $\mathrm{Cl}_{2}$ / chlorine is a gas so needs a strong container. IGNORE 'difficult to store / difficult to transport' <br> DO NOT ALLOW harmful / irritant / dangerous instead of toxic. <br> Answer must have 'by-products / THMs' and 'suspected carcinogens' to gain the mark. |
|  | Total | 20 |  |

## F334 Chemistry of Materials

| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| 1 (a) | 2-hydroxypropanoic acid $\checkmark \checkmark$ | 2 | mark independently <br> 2-hydroxy <br> DO NOT ALLOW hydroxyl <br> propanoic acid <br> ALLOW if propan- and -oic are separated. |
| (b) (i) | (enantiomers are) isomers whose structures are mirror images of one another $\checkmark$ <br> and are non-superimposable | 2 | mark independently <br> mirror images <br> non-superimposable <br> IGNORE references to 4 different groups around a C atom, optical isomerism, various chiral words \& rotation of plane polarised light |
| (ii) |  | 2 | 3D structure correct for one isomer $\checkmark$ <br> DO NOT ALLOW 90 or 180 degree angles between the two bonds in the plane of the paper <br> mirror-image correct (must have 4 bonds around the C) $\checkmark$ ALLOW ecf for non 3D structure with four different groups only <br> IGNORE the way the groups are bonded to carbon eg -OH or - HO , same for $\mathrm{COOH} \& \mathrm{CH}_{3}$ |
| (c) (i) | $\begin{aligned} & P=C \\ & Q=A \\ & R=B \end{aligned}$ | 1 |  |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| (ii) | $\mathrm{P}=(\mathbf{C})$ does not fizz / does not react / with $\mathrm{Na}_{2} \mathrm{CO}_{3}$ AND so no - COOH group present / AW $\checkmark$ <br> Q = no phenol group (in A), so $\mathrm{FeCl}_{3}$ remains yellow / AW $\checkmark$ <br> $R=(B$ has) both phenol \& carboxylic acid (COOH) (so will turn $\mathrm{FeCl}_{3}$ purple and will fizz with $\mathrm{Na}_{2} \mathrm{CO}_{3}$ ) / AW $\checkmark$ The words in brackets are only needed if tests not discussed for $P$ \& $Q$. | 3 | if answers to (i) are incorrect/no response award 1 mark for correct answers for having both tests for phenol and carboxylic acid: <br> ie: <br> purple solution $=$ phenol <br> AND acids fizz with carbonate $\checkmark$ <br> IGNORE $P$ is $\mathbf{C}$ as it turns $\mathrm{FeCl}_{3}$ purple <br> IGNORE $Q$ is $\mathbf{A}$ as it fizzes \& is therefore a carboxylic acid |
| (d) | C <br> because: <br> (broad) peak at around $3250\left(\mathrm{~cm}^{-1}\right)$ indicates alcohol or phenol / OH / hydroxyl group <br> no peak at 1700-1725 ( $\mathrm{cm}^{-1}$ ) so no $\underline{\mathrm{C}=\mathrm{O}}$ (in -COOH ) present | 3 | IGNORE any reference to carboxylic acid for the 3250 peak <br> ALLOW a range around 3250 <br> Peaks may be identified on the diagram <br> IGNORE all other peaks |
| (e) (i) | nucleophilic addition $\checkmark \checkmark$ | 2 | nucleophilic <br> addition <br> mark independently |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| (ii) |  | 4 | 'curly' arrow showing attack by ${ }^{-} \mathrm{CN}$ at $\mathrm{C}=\mathrm{O}$ carbon DO NOT ALLOW arrow starting from N of ${ }^{-} \mathrm{CN} /$ singleheaded arrows but give 1 mark if both are single headed but otherwise correct <br> $\mathrm{C}=\mathrm{O}$ bond polarised correctly $\checkmark$ <br> curly arrow showing movement of double bond <br> final structure correct ALLOW any correct structural formula not just skeletal <br> O MUST be -ve <br> IGNORE any further reaction showing $\mathrm{O}^{-}$gaining $\mathrm{H}^{+}$ |
| (iii) | the rate determining step (slow step) does not involve water $\checkmark$ <br> (since) water does not appear in the rate equation / water is zero order $\checkmark$ <br> Since water / H+ required to form product it must react in a subsequent (fast) step/there must be at least 2 steps in the reaction / AW $\checkmark$ <br> OR <br> the rate determining step (slow step) only involves ethanal \& cyanide <br> (because only) ethanal \& cyanide appear in the rate equation <br> and so water must react in a subsequent step $\checkmark$ | 3 |  |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :--- | :---: | :--- |
| (f) (i) | reaction 1.1 has a higher atom economy than <br> reaction $1.2 \checkmark$ | $\mathbf{2}$ | ALLOW comparison of percentage atom economy <br> eg reaction 1.1 has 100\% economy, reaction 1.2 does not. <br> because it is an addition reaction / only one product is <br> formed <br> whereas <br> in reaction 1.2 hydrolysis / condensation occurs / atoms are <br> wasted / lost / two 'products' are formed / co-products are <br> also formed / AW $\checkmark$ |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| 2 (a) | $\mathrm{Fe}^{3+}$ will oxidise Cu / ORA <br> OR <br> Cu loses electrons to form $\mathrm{Cu}^{2+} /$ ORA $\checkmark$ <br> because electrode potential of $\mathrm{Fe}^{3+} I\left(\mathrm{Fe}^{2+}\right)$ is more positive / ORA (involves the Copper half-cell) OR <br> Uses $\mathrm{E}_{\text {cell }}$ calculation to show reaction is feasible $\checkmark$ $\mathrm{Cu}+2 \mathrm{Fe}^{3+} \rightarrow \mathrm{Cu}^{2+}+2 \mathrm{Fe}^{2+} \checkmark$ | 3 | ALLOW Fe(III) and $\mathrm{Cu}(\mathrm{II})$ <br> DO NOT ALLOW electronegativity or higher / lower or larger / smaller <br> IGNORE state symbols |
| (b) (i) | Pt electrode for $\mathrm{Fe}^{3+} / \mathrm{Fe}^{2+}$ half cell in $\mathrm{Fe}^{3+} / \mathrm{Fe}^{2+}$ <br> Cu electrode in $\mathrm{Cu}^{2+}$ <br> salt bridge labelled and in solutions <br> conditions: $1 \mathrm{~mol} \mathrm{dm}^{-3}$ and $298 \mathrm{~K} / 25^{\circ} \mathrm{C}$ | 5 | ALLOW CuSO 4 etc. instead of $\mathrm{Cu}^{2+}$ <br> ALLOW 1 molar / 1M |
| (ii) | $0.43 \mathrm{~V} \checkmark$ | 1 | IGNORE any sign |
| (c) (i) | Copper(I) iodide $\checkmark$ | 1 | DO NOT ALLOW copper iodide ALLOW Copper I iodide |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| (ii) | 1. moles of thiosulfate $=0.200 \times(20.5 / 1000)=\mathbf{0 . 0 0 4 1}$ <br> 2. (moles of iodine $\left(\mathrm{I}_{2}\right)=0.5 \times$ answer from 1 (0.0041) and moles of $\mathrm{Cu}^{2+}(\mathrm{aq})$ in $25.0 \mathrm{~cm}^{3}=2 \times 0.5 \mathrm{x}$ answer from $1(0.0041)$ ) $=0.0041$ <br> 3. moles of $\mathrm{Cu}^{2+}(\mathrm{aq})$ in $250 \mathrm{~cm}^{-3}=10 \mathrm{x}$ answer from $1(0.0041)=0.0410 \checkmark$ <br> 4. mass of Cu in coin = answer from 3 (0.0410) $x 63.5=2.6035 \mathrm{~g} \checkmark$ <br> 5. \% of Cu in coin $=($ answer from $4(2.6035) / 3.47)$ $\times 100=75.0$ <br> 6. (3 sig. figs.) | 6 | 75.0 with no / incomplete working scores 6 marks. <br> The marks are awarded for the working out given in bold: <br> If final answer is incorrect please annotate marks given with ticks AND crosses where errors have occurred eg missing out step 3. <br> 1. moles of thiosulfate $=0.0041$ moles <br> 2. correct moles of $\mathrm{Cu}^{2+}(\mathrm{aq})$ in $25.0 \mathrm{~cm}^{3}$ ecf from 1 <br> 3. moles of $\mathrm{Cu}^{2+}(\mathrm{aq})$ in $250 \mathrm{~cm}^{-3}=10 \times$ moles of thiosulfate <br> 4. mass of Cu in coin $=$ moles of $\mathrm{Cu}^{2+}(\mathrm{aq})$ in $250 \mathrm{~cm}^{-3}$ x 63.5 <br> 5. \% of Cu in coin $=$ mass of Cu in coin $\times 100 \checkmark$ <br> 6. correct/incorrect answer MUST be given to 3 sig figs. <br> Steps 3 \& 4 may be in a different order |
| (d) (i) | (transmits) Blue <br> $\mathrm{Cu}^{2+}(\mathrm{aq})$ absorb red / orange light <br> QWC - absorb(s) / absorbing / absorption / absorbance / absorbed <br> One of these words has to be used to gain the second mark and spelling must be correct | 2 | IGNORE pale, deep or light etc. referring to blue, reflects DO NOT ALLOW green <br> ALLOW complementary colour / specific frequencies / wavelengths of light are absorbed <br> DO NOT ALLOW all other frequencies / colour / AW <br> Use of 'emit' is a CON for the $2^{\text {nd }}$ mark |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| (i) | Ligand substitution $\left[\mathrm{CuCl}_{4}\right]^{2-} \checkmark$ | 2 | ALLOW ligand replacement / displacement / exchange or complex formation <br> The charge is required <br> ALLOW $\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2} \mathrm{Cl}_{4}\right]^{2-}$ |
| (iii) | Any five from a-f: <br> a. make up standard solutions / of known concentrations (of $\mathrm{Cu}^{2+}(\mathrm{aq})$ ) <br> b. choose a suitable filter / set the colorimeter to a suitable wavelength <br> c. zero colorimeter with water / solvent / AW <br> d. measure absorbance / absorption of standard solutions / AW <br> e. plot calibration curve <br> f. measure unknown <br> AND <br> g. QWC read off concentration from calibration curve / AW $\checkmark$ | 6 | IGNORE transmittance |
|  | Total | 26 |  |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| $3 \text { (a) (i) }$ |  | 2 | the link between phosphate group and any -OH group on the sugar <br> ie $\mathrm{P}-\mathrm{O}-\mathrm{CH}_{2}-\mathrm{C}$ (in ring next to -O-) <br> or P-O-C (in ring) depending which -OH group is used $\checkmark$ correct linkage ( $\mathrm{P}-\mathrm{O}-\mathrm{CH}_{2}-\mathrm{C}$ ) and rest of structure correct $\checkmark$ |
| (ii) | Water / $\mathrm{H}_{2} \mathrm{O}$ and condensation (reaction) $\checkmark$ | 1 |  |
| (iii) |  | 1 | ALLOW <br> ALLOW if the $\mathbf{N}-\mathbf{H}$ group is circled |
| (iv) | lone pair (of electrons) on $\mathrm{N} /$ amine (group) $\checkmark$ accepts a proton / $\mathrm{H}^{+}$ | 2 |  |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| (b) |  | 3 | 2 hydrogen bonds correct $\checkmark$ <br> DO NOT ALLOW more than 2 bonds <br> correct polarities (all 3) on one group of atoms $\checkmark$ <br> lone pair on the NORO $\checkmark$ |
| (c) | two strands (not three) <br> phosphate groups on the outside (not on the inside) <br> bases face into the centre (not the outside) | 3 | ALLOW double helix <br> IGNORE phosphate backbone |
| (d) | for removal: <br> infringement of personal liberty / AW / <br> 'fingerprint' not unique only probability / <br> techniques used not foolproof / <br> law / type of government might change changing accessibility / AW <br> against removal: <br> helps to solve many crimes, particularly 'cold' crimes / <br> 'innocent until proved guilty' / AW <br> future research into disease | 2 | IGNORE hacking into database / leakage of data (NOTE: probability may only be 1 in 20 for some population groups) |
|  | Total | 14 |  |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| 4 (a) |   | 2 | 1 mark for each structure correct <br> DO NOT ALLOW missing H atoms if structural formulae are drawn |
| (b) (i) | there are (two) different groups on each carbon of a $\mathrm{C}=\mathrm{C}$ in neoprene <br> these groups can not rotate about the double bond | 2 | May be shown in a diagram <br> ALLOW restricted / limited rotation about the double bond |
| (ii) | (less trans linkages) will make the chains less linear / less regular / less ordered / ORA $\checkmark$ <br> so they can not line up / be arranged so regularly / closely and the crystallinity will be reduced | 2 | ALLOW chains have a kink / are more randomly arranged <br> Note: <br> $1^{\text {st }}$ mark is for shape of chain, <br> $2^{\text {nd }}$ mark is for relative arrangement of chains. |
| (c) | -CONH group / NH group / amide group allows nylon to form hydrogen bonds with water molecules <br> no hydrogen bonding in neoprene <br> water molecules will force chains further apart / chains will not fit as closely together / AW $\checkmark$ <br> chains less crystalline / weaker intermolecular forces so $T_{\mathrm{g}}$ will be lowered $\checkmark$ | 4 | IGNORE chains sliding over each other |

\(\left.\begin{array}{|c|c|c|l|}\hline Question \& Expected Answers \& Marks \& Additional Guidance <br>
\hline (d) \& \& \mathbf{2} \& 6 carbon atoms and - \mathrm{NH}_{2} group \checkmark <br>

-\mathrm{COO}^{-} \checkmark\end{array}\right]\)| ALLOW skeletal or any equivalent/alternative structural |
| :--- |
| formula |
| DO NOT ALLOW structures with missing H atoms |

| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| 5 (a) (i) | $3 \mathrm{H}_{2} \mathrm{~S}+\underset{\checkmark}{2 \mathrm{MnO}_{4}^{-}}+2 \mathrm{H}^{+} \rightarrow 2 \mathrm{MnO}_{2}+3 \mathrm{~S}+4 \mathrm{H}_{2} \mathrm{O}$ | 1 | all 4 numbers MUST be correct |
| (ii) | oxidation state $=+4 \checkmark$ | 1 | DO NOT ALLOW 4+ OR 4 |
| (b) (i) | iron(III) hydroxide $\checkmark$ | 1 | DO NOT ALLOW iron hydroxide / Fe(OH) ${ }_{3}$ |
| (ii) | $\mathrm{Fe}^{3+}(\mathrm{aq})+3 \mathrm{OH}^{-}(\mathrm{aq}) \rightarrow \mathrm{Fe}(\mathrm{OH})_{3}(\mathrm{~s})$ | 2 | equation correct <br> state symbols correct <br> ALLOW ecf for iron(II) hydroxide in (i) |
| (c) (i) | The large excess of ethanedioate and acid means that their concentrations were virtually constant during the reaction / concentrations hardly changed / concentrations were high so little effect on rate / AW $\checkmark$ | 1 | 'A large excess of ethanedioate and acid' by itself does not get the mark <br> IGNORE 'excess ethanedioate and acid not limit the rate of reaction' / AW <br> Look for concentration in answer |
| (ii) | One of the following: <br> Method 1 half-lives determination of at least two half-lives, 13-15 s half-life constant $\checkmark$ first order $\checkmark$ <br> OR <br> Method 2 finding rate at different concentrations calculation of at least two rates rate is proportional to concentration $\checkmark$ first order $\checkmark$ | 3 | Two values for half-life MUST be given Working must be shown as either a calculation or by lines on graph. |


| Question | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| (iii) | 3d     <br> 4s     <br> $A$ $\uparrow$ $\uparrow$ $\uparrow$ $\perp$ | 1 |  |
| (d) | One from the following: <br> loss of $\mathrm{CO}_{2} / \mathrm{CO}_{2}$ produced $\checkmark$ by weighing / gas collection / measuring milkiness of lime-water $\checkmark$ <br> OR <br> titration of $\mathrm{MnO}_{4}^{-} \checkmark$ with (standard) $\mathrm{Fe}^{2+}(\mathrm{aq}) \checkmark$ <br> OR <br> titration of $\mathrm{H}^{+}(\mathrm{aq})$ decrease $\checkmark$ with $\mathrm{OH}^{-} / \mathrm{CO}_{3}{ }^{2-}(\mathrm{aq}) \checkmark$ <br> OR <br> measure pH change $\checkmark \mathrm{H}^{+}$ions used in the reaction | 2 | The two parts are marked independently <br> ALLOW gas for $\mathrm{CO}_{2}$ and measuring volume for gas collection |
|  | Total | 12 |  |

## Grade Thresholds

Advanced GCE Chemistry B (Salters) (H035 H435) January 2010 Examination Series

Unit Threshold Marks

| Unit |  | Maximum <br> Mark | A | B | C | D | E | U |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| F331 | Raw | 60 | 41 | 36 | 31 | 26 | 21 | 0 |
|  | UMS | 90 | 72 | 63 | 54 | 45 | 36 | 0 |
| F332 | Raw | 100 | 74 | 67 | 60 | 54 | 48 | 0 |
|  | UMS | 150 | 120 | 105 | 90 | 75 | 60 | 0 |
| F334 | Raw | 90 | 65 | 58 | 51 | 44 | 37 | 0 |
|  | UMS | 90 | 72 | 63 | 54 | 45 | 36 | 0 |

## Specification Aggregation Results

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

|  | Maximum <br> Mark | A | B | C | D | E | U |
| :--- | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| H035 | 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 <br> Candidates |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\mathbf{H 0 3 5}$ | 12.9 | 37.1 | 61.7 | 83.6 | 97.4 | 100.0 | 823 |

## 823 candidates aggregated this series

For a description of how UMS marks are calculated see:
http://www.ocr.org.uk/learners/ums/index.html
Statistics are correct at the time of publication.

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