## Chemistry A

## Advanced Subsidiary GCE

## Mark Scheme for January 2012

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All examiners are instructed that alternative correct answers and unexpected approaches in candidates' scripts must be given marks that fairly reflect the relevant knowledge and skills demonstrated.

Mark schemes should be read in conjunction with the published question papers and the report on the examination.

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## Annotations

| Annotation | Meaning |
| :---: | :---: |
| [-Tom | Benefit of doubt given |
| [ $¢ \cdot 0]$ | Contradiction |
| 3 | Incorrect response |
| [-4 | Error carried forward |
| $\square$ | Ignore |
| [W0] | Not answered question |
| - | Benefit of doubt not given |
| Wir | Power of 10 error |
| -1. | Omission mark |
| [1:7 | Rounding error |
| $\Gamma \mathrm{F}$ | Error in number of significant figures |
|  | Correct response |

Abbreviations, annotations and conventions used in the detailed Mark Scheme (to include abbreviations and subject-specific conventions).

| Annotation | Meaning |
| :--- | :--- |
| DO NOT ALLOW | Answers which are not worthy of credit |
| IGNORE | Statements which are irrelevant |
| ALLOW | Answers that can be accepted |
| () | Words which are not essential to gain credit |
| ECF | Underlined words must be present in answer to score a mark |
| AW | Error carried forward |
| ORA | Alternative wording |

## Subject-specific Marking Instructions

The following questions should be annotated with ticks to show where marks have been awarded in the body of the text:

| Q2 | (b)(i) |
| :--- | :--- |
| Q4 | (c) |
| Q8 | (d) |
| Q8 | (e) |

## All questions where an ECF has been applied.

## Checking additional pages

All the Additional Pages in the examination script must be checked to see if any candidates include any answers.

- When you open question $\mathbf{1 ( a )}$ you will see a view of page 23 one of the Additional Pages.
- If the page is blank then, using the marking mode, annotate the page with an omission mark, ^.
- Scroll down to page 24 and annotate with $\mathrm{a}^{\wedge}$ if the page is blank.
- If pages 23 or 24 are not blank then use the paper clip icon to link the pages to the correct questions.
- You may need to contact your Team Leader if you do not know how to do this.


## Generic comments

## ORGANIC STRUCTURES

For a 'structure' or 'structural formula',

- ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous)

For an alkyl group shown within a structure,

- ALLOW bond drawn to C or H ,
e.g.ALLOW $\mathrm{CH}_{3}-, \mathrm{CH}_{2}-, \mathrm{C}_{3} \mathrm{H}_{7}-$, etc
- ALLOW vertical 'bond' to any part of an alkyl group

For an OH group shown within a structure,

- DO NOT ALLOW formula with horizontal -HO OR OH -
- ALLOW vertical 'bond' to any part of the OH group

For a CHO group shown within a structure,

- DO NOT ALLOW COH

For a 3D structure,

- For bond in the plane of paper, a solid line is expected:
- For bond out of plane of paper, a solid wedge is expected:
- For bond into plane of paper, ALLOW:
- ALLOW a hollow wedge for 'in bond' OR an 'out bond', provided it is different from the other in or out wedge e.g.:



## NAMES

Names including alkyl groups:

- ALLOW alkanyl, e.g. ethanyl (i.e. IGNORE 'an')
- DO NOT ALLOW alkol, e.g. ethol (ie 'an' is essential)

Names of esters:

- Two words are expected, e.g. ethyl ethanoate
- ALLOW one word, e.g. ethylethanoate

Names with multiple numbers and hyphens:
Use of 'e'

- ALLOW superfluous 'e' , e.g. propane-1-ol ('e' is kept if followed by consonant)
- ALLOW absence of 'e', e.g. propan-1,2-diol ('e' is omitted if followed by vowel)

Hyphens separate name from numbers:

- ALLOW absence of hyphens, e.g. propane 1,2 diol

Multiple locant numbers must be clearly separated

- ALLOW full stops: e.g. 1.2 OR spaces: 12
- DO NOT ALLOW e.g. 12

Locant numbers in formula must be correct

- DO NOT ALLOW propan-3-ol

Order of substituents should be alphabetical:

- ALLOW any order (as long as unambiguous), e.g. 2-chloro-3-bromobutane


## ABBREVIATIONS

van der Waal's forces
ALLOW vdw forces OR VDW forces (and any combination of upper and lower cases)

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


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (d) | (ii) | $\mathrm{C}_{7} \mathrm{H}_{16} \rightarrow \mathrm{C}_{7} \mathrm{H}_{14}+\mathrm{H}_{2} \checkmark$ | 1 | ALLOW molecular formulae OR correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) <br> DO NOT ALLOW cycloheptane structure in equation |
|  | (e) |  | D has more surface (area of) contact OR D is a bigger molecule <br> D has more van der Waals' forces OR C have fewer van der Waals' forces $\checkmark$ | 2 | Both answers need to be comparisons <br> Assume 'it' refers to D <br> ALLOW has more electrons OR longer (carbon) chain OR higher molecular mass <br> IGNORE surface area <br> ALLOW ORA <br> ALLOW D has stronger van der Waals' forces / larger VDW / greater VDW <br> OR C has weaker van der Waals' forces OR C has smaller VDW <br> ALLOW more VDW forces <br> More intermolecular forces is not sufficient <br> DO NOT ALLOW reference to bonds breaking or more bonds present unless it is clear that that the bonds are VDW |
|  | (f) |  | Same structural formula <br> Different arrangement of groups around a double bond OR different arrangement (of atoms) in space $\checkmark$ | 2 | ALLOW have the same structure / displayed formula / skeletal formula <br> Stereoisomers have the same formula or molecular formula is not sufficient <br> ALLOW different three dimensional arrangement |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (g) |  | $\mathrm{C}_{7} \mathrm{H}_{16}+11 \mathrm{O}_{2} \rightarrow 7 \mathrm{CO}_{2}+8 \mathrm{H}_{2} \mathrm{O}$ <br> Correct reactants and products $\checkmark$ Balancing $\checkmark$ | 2 | ALLOW molecular formulae OR correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) <br> ALLOW any correct multiple <br> IGNORE state symbols <br> Balancing is dependent on correct formulae |
|  | (h) |  | $\mathrm{C}_{16} \mathrm{H}_{34} \rightarrow \mathrm{C}_{8} \mathrm{H}_{18}+2 \mathrm{C}_{4} \mathrm{H}_{8} \checkmark$ | 1 | ALLOW molecular formulae OR correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) <br> ALLOW any correct multiple <br> ALLOW structural OR displayed OR skeletal formulae in equation <br> ALLOW but-1-ene <br> IGNORE state symbols |
|  | (i) | (i) | Group of atoms (in a molecule or compound) that is responsible for the reactions $\checkmark$ | 1 | ALLOW the 'part' (of the molecule or compound) that reacts ALLOW the group of atoms that gives the chemical properties <br> ALLOW group of atoms which indicates the homologous series |
|  |  | (ii) | $8 \checkmark$ | 1 |  |
|  |  | (iii) | has an unpaired electron $\checkmark$ | 1 | ALLOW plural i.e. unpaired electrons has a lone OR single OR free electron is not sufficient |
|  |  |  | Total | 16 |  |



| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 2 | (b) | (i) |  | 5 | ANNOTATE ANSWER WITH TICKS AND CROSSES ETC |
|  |  |  | Contains C=O bond because of absorption between 1700 and $1740 \mathrm{~cm}^{-1}$ (from the spectrum) $\checkmark$ |  | ALLOW contains a carbonyl group because of absorption within range $1640-1750 \mathrm{~cm}^{-1}$ OR contains an aldehyde, ketone or carboxylic acid because of absorption within range $1640-1750 \mathrm{~cm}^{-1}$ <br> Mention of only an aldehyde or a ketone is not sufficient it needs reference to the wavenumber <br> LOOK FOR THIS MARK ON THE SPECTRUM |
|  |  |  | does not contain an $\mathrm{O}-\mathrm{H}$ bond $\checkmark$ |  | ALLOW not a carboxylic acid ALLOW does not have any other characteristic absorbance due to other functional groups |
|  |  |  | (So was a) ketone OR aldehyde $\checkmark$ |  | ALLOW (so was a) carbonyl compound ALLOW this mark if a structure of an aldehyde or a ketone is given even if the structure has an incorrect number of carbon atoms |
|  |  |  | $M_{r}=86 \checkmark$ |  |  |
|  |  |  | Correct structure $\checkmark$ |  | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) |
|  |  |  |  |  | LOOK FOR AN ALDEHYDE or KETONE with FIVE carbon atoms OR a DIALDEHYDE, DIONE OR an OXOALDEHYDE with FOUR carbon atoms - a comprehensive list of correct structures is shown on page 34 IGNORE incorrect name |
|  |  |  |  |  | DO NOT ALLOW COH for an aldehyde |


| Question |  | Garks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |


| Question |  |  | Answer | Guidance |
| :---: | :---: | :---: | :---: | :---: | :--- |
| (b) | (ii) | Correct structure $\checkmark$ | 2 | $\begin{array}{l}\text { ALLOW correct structural OR displayed OR skeletal formula } \\ \text { OR mixture of the above (as long as unambiguous) }\end{array}$ |
| All bonds and all hydrogen atoms must be shown in a |  |  |  |  |
| displayed formula within this question |  |  |  |  |$]$| Name must correspond to the correct structure for two marks |
| :--- |
| ALLOW butanoic acid or 2-methylpropanoic acid if the |
| structure drawn is incorrect |
| There is no ECF in this question |


| Question |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 2 | (c) | Use of propan-1-ol $\mathrm{CH}_{3} \mathrm{COOH}+\mathrm{C}_{3} \mathrm{H}_{7} \mathrm{OH} \rightarrow \mathrm{CH}_{3} \mathrm{COOCH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{3}+\mathrm{H}_{2} \mathrm{O}$ <br> Correct formulae for the ester $\checkmark$ Correctly balanced equation <br> Add $\mathrm{H}_{2} \mathrm{SO}_{4} \mathrm{OR}$ acid catalyst $\mathrm{OR} \mathrm{H}^{+} \checkmark$ | 4 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) ALLOW from the equation propanol $\mathrm{OR}_{3} \mathrm{H}_{7} \mathrm{OH}$ is not sufficient <br> ALLOW molecular formula OR correct structural OR displayed OR skeletal formula OR mixture of the above ALLOW propan-2-ol in the equation <br> ALLOW conditions mark over the arrow in the equation |
|  |  | Total | 14 |  |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (a) | (i) | FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer = 2.68 (kJ) award 2 marks $q=m c \Delta T \text { OR }=50.0 \times 4.18 \times 12.8 \checkmark$ $=2.68(\mathrm{~kJ}) \checkmark$ | 2 | IF there is an alternative answer, check to see if there is any ECF credit possible using working below <br> IF ECF, ANNOTATE WITH TICKS AND CROSSES, etc <br> ALLOW ecf only from using mass of 50.486 <br> ALLOW 2675.2 J <br> IGNORE sign <br> If mass used is 50.486 answer is 2701.202944 <br> ALLOW 2.7 OR 2.675 OR 2.6752 <br> DO NOT ALLOW 3 <br> IGNORE sign <br> If mass used is 50.486 answer is 2.7, 2.70, 2.701 up to calculated value of 2.701202944 correctly rounded <br> ALLOW one mark for using 4.2 and correctly calculating q in kJ to at least 2 sig figs |
|  |  | (ii) | amount $=0.02(00)(\mathrm{mol})^{\checkmark}$ | 1 | ALLOW 1/50 <br> IGNORE trailing zeroes |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (a) | (iii) | FIRST, CHECK THE ANSWER ON ANSWER LINE <br> IF answer = -134 (kJ) award 3 marks <br> IF answer = +134 (kJ) award 2 marks $2.68 \div 0.02 \checkmark$ <br> Correctly calculates the value to 3 sig figs | 3 | IF there is an alternative answer, check to see if there is any ECF credit possible using working below <br> IF ECF, ANNOTATE WITH TICKS AND CROSSES, etc <br> ALLOW ECF i.e. $\frac{(i)}{(i i)}$ <br> This is dependant on the previous mark <br> ALLOW ECF <br> If $2.68,2.675$ or 2.6752 and moles of 0.02answer is $(-) 134$ <br> If mass of magnesium included answer is (-)135 <br> If 2.7 kJ and moles of 0.02 used answer is (-)135 <br> ALLOW only answers to three significant figures <br> - sign is independent of answer |
|  | (b) | (i) | (Enthalpy change) when one mole of a compound <br> is formed from its elements <br> at $25^{\circ} \mathrm{C} / 298 \mathrm{~K}$ AND 1 atmosphere/101 kPa | 3 | ALLOW energy required OR energy released <br> ALLOW (energy change) when one mole of a substance/molecule/product <br> DO NOT ALLOW enthalpy change for one mole of products DO NOT ALLOW one mole of reactants <br> ALLOW any stated temperature and 1 bar/1000/mb/100kPa/100000Pa/101000Pa/101000 $\mathrm{Nm}^{-2}$ etc IGNORE reference to concentration |



| Question |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 4 | (a) | FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer $=431.5\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ award 2 marks <br> Energy required to break bonds $=(+) 679 \mathrm{~kJ} \checkmark$ <br> so bond enthalpy $=(+) 431.5 \checkmark$ | 2 | IF there is an alternative answer, check to see if there is any ECF credit possible using working below <br> IF ECF, ANNOTATE WITH TICKS AND CROSSES, etc <br> ALLOW (+)432 <br> ALLOW one mark in this question for -431.5 OR (+)863 ALLOW ecf for bond enthalpy $=0.5 \times(-184+$ energy required to break bonds) |
|  | (b) | more concentrated (particles) OR more particles per (unit) volume $\checkmark$ <br> more collisions per second OR more frequent collisions $\checkmark$ | 2 | Must state somewhere in the answer that the rate is faster for full marks <br> ALLOW ORA if lower pressure is specified <br> ALLOW particles are closer together <br> OR more crowded particles <br> OR more particles in the same space <br> OR same number of particles in a smaller volume <br> ALLOW molecules for particles but DO NOT ALLOW atoms <br> DO NOT ALLOW 'area' instead of 'volume' <br> ALLOW collisions more often <br> OR increased rate of collision <br> OR collisions are more likely <br> OR there is a greater chance of collisions <br> 'More collisions' is not sufficient <br> IGNORE successful |



| Question |  | Marss | Guidance |
| :--- | :--- | :--- | :--- | :--- | :--- |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | (a) |  | Only one (desired) product formed $\checkmark$ | 1 | ALLOW no waste products OR no co-product OR all atoms on left hand side are in the desired product OR sulfuric acid is the only product IGNORE it is an addition reaction |
|  | (b) |  | FIRST, CHECK THE ANSWER ON ANSWER LINE <br> IF answer = 94\% award 3 marks <br> Moles of sulfur reacted or theoretical moles of $\mathrm{H}_{2} \mathrm{SO}_{4}=$ $1.60 \times 10^{6} \checkmark$ <br> Actual moles of $\mathrm{H}_{2} \mathrm{SO}_{4}=1.50 \times 10^{6} \checkmark$ <br> $\%$ yield $=94 \checkmark$ | 3 | IF there is an alternative answer, check to see if there is any ECF credit possible using working below <br> ALLOW $1.6 \times 10^{6}$ to the calculator value $1.601246106 \times 10^{6}$ correctly rounded <br> ALLOW 1.60 up to calculator value 1.601246106 correctly rounded <br> ALLOW $1.5 \times 10^{6}$ to the calculator value $1.498470948 \times 10^{6}$ correctly rounded <br> ALLOW 1.5 up to calculator value 1.498470948 correctly rounded <br> ALLOW theoretical mass of $\mathrm{H}_{2} \mathrm{SO}_{4}=157$ (tonnes) up to the calculator value of 157.0822430 correctly rounded for two marks <br> ALLOW ECF for a percentage yield from wrong moles above but answer must have two significant figures |
|  | (c) | (i) | Position of equilibrium - unchanged <br> Rate of backward reaction - decreases $\checkmark$ | 2 |  |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | (c) | (ii) | (equilibrium position shifts) to the left because (forward) reaction is exothermic <br> OR equilibrium position shifts) to the left because reverse reaction is endothermic | 1 | Both position of equilibrium AND explanation needed for one mark <br> Note: ALLOW suitable alternatives for 'to left', e.g. towards $\mathrm{SO}_{2}$ or $\mathrm{O}_{2}$ / towards reactants OR in backward direction OR in reverse direction OR decreases yield of $\mathrm{SO}_{3} /$ products <br> ALLOW 'favours the left', as alternative for 'shifts equilibrium to left' <br> ALLOW reaction gives out heat for exothermic ALLOW reaction takes in heat for endothermic ALLOW moves to the left in the endothermic direction <br> ALLOW ORA if specified IGNORE responses in terms of rate |
|  |  | (iii) | (equilibrium position shifts) to the left because there are more moles (of gas) on the reactant side <br> OR <br> (equilibrium position shifts) to the left because there are fewer moles (of gas) on product side | 1 | Both position of equilibrium AND explanation needed for one mark <br> Note: ALLOW suitable alternatives for 'to left', e.g.: towards $\mathrm{SO}_{2}$ or $\mathrm{O}_{2}$ / towards reactants OR in backward direction OR in reverse direction OR decreases yield of $\mathrm{SO}_{3} /$ products <br> ALLOW 'favours the left', as alternative for 'shifts equilibrium to left' <br> ALLOW correct reference to volume of gases e.g. shifts to the left because there is a smaller volume of gas on the product side <br> ALLOW ORA if specified <br> IGNORE responses in terms of rate |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | (d) | (i) | Correct structure $\checkmark$ <br> OR <br> OR | 1 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) <br> ALLOW bonds going to any part of the $\mathrm{CH}_{3}, \mathrm{CH}_{2}$ and CH bonds <br> ALLOW vertical 'bond' to any part of the OH group DO NOT ALLOW horizontal -HO in the formula <br> ALLOW as a slip one stick with no H on in a displayed formula <br> IGNORE name |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | (d) | (ii) | Correct structure for $\mathbf{L}$ <br> Correct structure for $\mathbf{M} \checkmark$ <br> Correct structure for $\mathbf{N} \checkmark$ | 3 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) for $\mathbf{L}, \mathbf{M}$ and $\mathbf{N}$ <br> e.g.  <br> L or M  <br> L or M $\mathrm{N}-\mathrm{CH}_{2} \mathrm{CHCH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{3}$ <br> Answers to $L$ and $M$ are interchangeable <br> IGNORE cis/trans OR E/Z labels <br> ALLOW as a slip one stick with no H on in a displayed formula <br> ALLOW 2 marks if three correct structures are drawn but some are in the wrong boxes <br> ALLOW 1 mark if two correct structures are drawn but in the wrong boxes |


| Question |  |  | Answer |  | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | (d) | (iii) |  |  | 1 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) <br> ALLOW vertical 'bond' to any part of the OH group DO NOT ALLOW horizontal -HO in the formula <br> ALLOW as a slip one stick with no H on in a displayed formula |
|  |  |  |  | Total | 13 |  |


| Question |  |  | Answer |  | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 6 | (a) | (i) | $(m / z=) 46 \checkmark$ |  | 1 |  |
|  |  | (ii) | $\mathrm{CH}_{3} \mathrm{O}^{+} \mathrm{OR} \mathrm{CH} \mathrm{OH}^{+} \checkmark$ |  | 1 | MUST show '+' |
|  |  | (iii) | $\mathrm{C}_{2} \mathrm{H}_{6} \mathrm{O} \checkmark$ |  | 1 | ALLOW $\mathrm{H}_{2} \mathrm{CO}_{2}$ |
|  | (b) |  | $\frac{63 \times 72.2+65 \times 27.8}{100}$ OR 63.556 OR $63.56 \checkmark$ $A_{\mathrm{r}}=63.6 \mathrm{~V}$ <br> Copper / Cu $\checkmark$ |  | 3 | ALLOW two marks for 63.6 with no working out |
|  |  |  |  | Total | 6 |  |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 7 | (a) |  | Shape - tetrahedral $\checkmark$ <br> Bond angle $109.5^{\circ} \checkmark$ | 2 | ALLOW 109-110 ${ }^{\circ}$ |
|  | (b) | (i) | Volatile OR non-toxic OR non-flammable OR easily vaporised | 1 | ALLOW not carcinogenic / not an irritant / not harmful / not hazardous <br> IGNORE cheap / not dangerous / gas / low boiling point DO NOT ALLOW inflammable |
|  |  | (ii) | (C-F or $\mathrm{C}-\mathrm{Cl}$ ) bonds need a large amount of energy to break $\checkmark$ | 1 | ALLOW (the C-F or C-Cl) bonds are strong / bonds have a large bond enthalpy <br> ALLOW the molecule is not polar enough / non-polar molecule is not sufficient ALLOW the activation energy is too high DO NOT ALLOW dissolves IGNORE references to hydrogen bonding |
|  | (c) |  | $\mathrm{CF}_{2} \mathrm{Cl}_{2} \rightarrow \mathrm{CF}_{2} \mathrm{Cl}+\mathrm{Cl} \checkmark$ <br> AND ANY TWO FROM <br> Cl catalyses the decomposition of ozone $\checkmark$ $\begin{aligned} & \mathrm{Cl}+\mathrm{O}_{3} \rightarrow \mathrm{ClO}+\mathrm{O}_{2} \checkmark \\ & \mathrm{ClO}+\mathrm{O} \rightarrow \mathrm{Cl}+\mathrm{O}_{2} \checkmark \end{aligned}$ | 3 | ALLOW $\mathrm{CF}_{2} \mathrm{Cl}_{2}$ (breaks down to) produces chlorine atoms/radicals <br> ALLOW equation with any CFC <br> ALLOW ClO $+\mathrm{O}_{3} \rightarrow \mathrm{Cl}+2 \mathrm{O}_{2}$ <br> ALLOW $\mathrm{O}_{3}+\mathrm{O} \rightarrow 2 \mathrm{O}_{2} \mathrm{OR}_{3} \mathrm{O}_{2} \rightarrow 2 \mathrm{O}_{3}$ for one mark if the two equations for the steps have not been given <br> IGNORE other propagation equations |


| Question |  | Answer | Marks | Guidance |
| :--- | :--- | :--- | :---: | :--- |
| 7 | (d) | Because (more) $\underline{\text { UV will reach the Earth's surface and risk }}$of (skin) cancer increased/risk of cataracts/crop mutation <br> increased $\checkmark$ <br> (e)Ideas related to uses <br> CFCs are still entering the atmosphere (from disused <br> items) OR CFCs are still used (for some purposes and by <br> some countries) $\checkmark$ | 2 | DO NOT ALLOW global warming <br> ALLOW protects from UV which causes skin cancer etc |
| Ideas relating to lifetime within the atmosphere <br> CFCCs have a long lifetime in the atmosphere OR it takes a <br> long time for CFCs to reach upper atmosphere OR CFCs <br> are inert $\checkmark$ | ALLOW 'stratosphere' for 'upper atmosphere' <br> ALLOW CFCCs are still entering the ozone layer |  |  |  |
|  |  | Total | 10 |  |


|  | uestion | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 8 | (a) |  <br> compound $P \checkmark$ <br> compound $Q \checkmark$ <br> compound $R \checkmark$ | 3 | ALLOW structures with missing hydrogen atoms on the carbon atoms that do not take part in the reaction. i.e. all hydrogen atoms must be shown in $\mathbf{Q}$ but not in $\mathbf{P}$ and $\mathbf{R}$ <br> For example for the structures of $\mathbf{P}$ and $\mathbf{R}$ |


|  | est | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 8 | (b) | Orange OR brown to colourless $\checkmark$ | 1 | ALLOW shades of orange OR yellow OR brown DO NOT ALLOW red alone DO NOT ALLOW any response that includes precipitate OR solid, irrespective of colour |
|  | (c) | Two or more repeat units $\checkmark$ | 1 | ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) <br> Must have at least two repeat units and the free bonds at the end <br> ALLOW free bonds with dotted lines <br> All carbon-carbon bonds in the polymer chain must be shown <br> IGNORE any brackets drawn IGNORE any missing hydrogen atoms on the $\mathrm{CH}_{2}$ groups <br> ALLOW skeletal formula |


|  | esti | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 8 | (d) | Curly arrow from double bond to attack hydrogen of $\mathrm{H}-\mathrm{Cl}$ and breaking of $\mathrm{H}-\mathrm{Cl}$ bond $\checkmark$ <br> Correct dipole shown on $\mathrm{H}-\mathrm{Cl} \checkmark$ <br> Correct carbonium ion drawn $\checkmark$ <br> Curly arrow from $\mathrm{Cl}^{-}$to the carbonium ion $\checkmark$ <br> Correct product $\checkmark$ | 5 | ANNOTATE ANSWER WITH TICKS AND CROSSES ETC <br> Curly arrow must start from the double bond and not a carbon atom; other curly arrow must start from $\mathrm{H}-\mathrm{Cl}$ bond <br> DO NOT ALLOW dipoles on double bond <br> Dipole must be partial charge and not full charge <br> Carbocation needs a full charge and not a partial charge (charges do not need to be in a circle) <br> C $/$ curly arrow must come from one lone pair on $\mathrm{Cl}^{-}$ ion OR from minus sign on $\mathrm{Cl}^{-}$ion Lone pair does not need to be shown on Cl ion <br> ALLOW structures with missing hydrogen atoms on the $\mathrm{CH}_{2}$ groups |


| Question |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 8 | (e) | Nucleophilic substitution <br> Heterolytic (fission) spelt correctly <br> dipole shown on $\mathrm{C}-\mathrm{Cl}$ bond, $\mathrm{C}^{\delta+}$ and $\mathrm{Cl}{ }^{\delta-} \downarrow$ <br> curly arrow from $\mathrm{HO}^{-}$to carbon atom of $\mathrm{C}-\mathrm{Cl}$ bond <br> curly arrow from $\mathrm{C}-\mathrm{Cl}$ bond to the chlorine atom and formation of Cl | 5 | ANNOTATE ANSWER WITH TICKS AND CROSSES ETC <br> Dipole must be partial charge and not full charge <br> $\mathrm{HO}^{-}$curly arrow must come from one lone pair on O of $\mathrm{HO}^{-}$ion OR from minus sign on $\mathrm{HO}^{-}$ion <br> curly arrow must start from $\mathrm{C}-\mathrm{Cl}$ bond and not from C atom <br> ALLOW structures with missing hydrogen atoms on the $\mathrm{CH}_{2}$ groups <br> ALLOW SN 1 mechanism <br> dipole shown on $\mathrm{C}-\mathrm{Cl}$ bond, $\mathrm{C}^{\delta+}$ and $\mathrm{Cl}^{\delta-} \checkmark$ curly arrow from $\mathrm{C}-\mathrm{Cl}$ bond to the Cl atom and Cl shown $\checkmark$ <br> curly arrow from $\mathrm{HO}^{-}$to correct carbonium ion $\checkmark$ |
|  |  | Total | 15 |  |

APPENDIX 1













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