

**GCE**

**Chemistry B (Salters)**

Unit **F334**: Chemistry of Materials

Advanced GCE

**Mark Scheme for June 2015**

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This mark scheme is published as an aid to teachers and students, to indicate the requirements of the examination. It shows the basis on which marks were awarded by examiners. It does not indicate the details of the discussions which took place at an examiners' meeting before marking commenced.

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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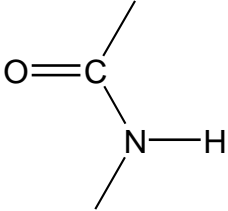
**Annotations** used in the detailed Mark Scheme (to include abbreviations and subject-specific conventions).

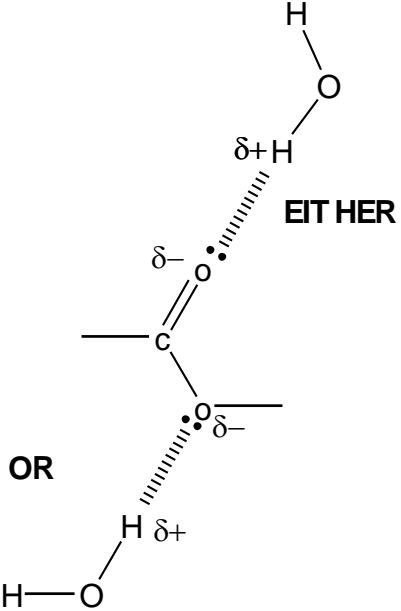
<b>Annotation</b>	<b>Meaning</b>
/	alternative and acceptable answers for the same marking point
✓	separates marking points
<b>not</b>	answers which are not worthy of credit and which will CON a correct answer
<b>ignore</b>	statements which are irrelevant and will NOT 'CON' a correct answer
<b>allow</b>	answers that can be accepted
( )	words which are not essential to gain credit
<u>    </u>	underlined words must be present in answer to score a mark
ecf	error carried forward
AW	alternative wording (replaces the old 'or words to that effect')
ora	or reverse argument

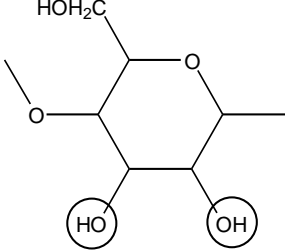
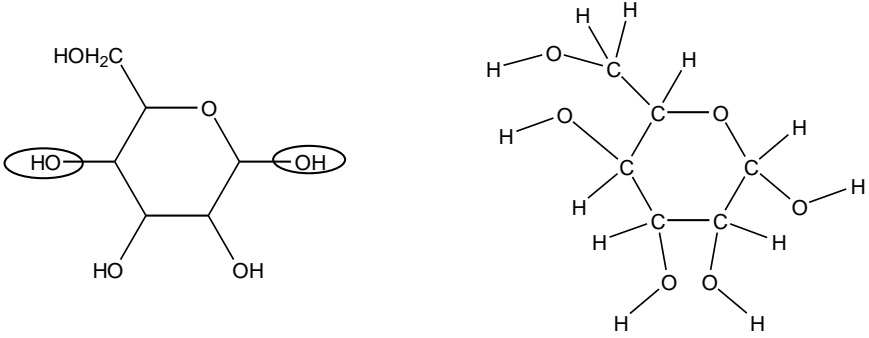
**Annotations** used in scoris:

<b>Annotation</b>	<b>Meaning</b>
✓	correct response
✗	incorrect response
bod	benefit of the doubt
nbod	benefit of the doubt <b>not</b> given
ECF	error carried forward
^	information omitted
I	Ignore
R	Reject

**Subject-specific Marking Instructions** that apply across the whole question paper to be included here.

Question	Answer	Mark	Guidance
1 a i	ammonia / NH <sub>3</sub> ✓	1	NH <sub>4</sub> <sup>+</sup> is a <b>CON</b> <b>DO NOT ALLOW</b> ammonium hydroxide, ammonium salt <b>IGNORE</b> conc., state
1 a ii	<b>N</b> (atom) has <b>lone pair</b> (of electrons) ✓ which can <b>accept a proton</b> / hydrogen ion /H <sup>+</sup> ✓	2	
1 b i	addition elimination ✓	1	<b>ALLOW</b> circles instead of underlining
1 b ii	 ✓	1	<b>IGNORE</b> additions to this linkage, but bonds must be shown on C and N.

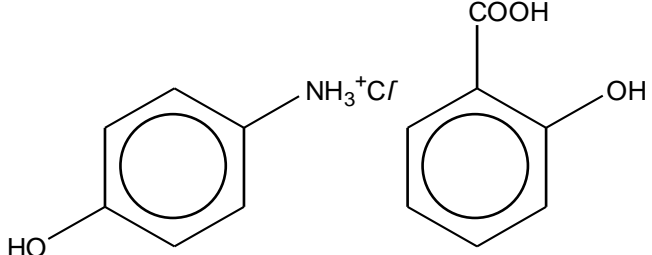
Question	Answer	Mark	Guidance
1 c	 <p>hydrogen bond between O and H ✓ lone pair AND partial charges ✓</p>	2	<p>Water must be shown as either H<sub>2</sub>O or HO<sub>2</sub> <b>DO NOT ALLOW</b> OH or H alone If water is given as <b>HO<sub>2</sub></b> AND hydrogen bond(s) and detail correct then award 1 mark</p> <p><b>IGNORE</b> bond angles</p> <p>If <b>two</b> examples given, <b>both</b> must be correct for 2 marks</p> <p>hydrogen bond must be as shown or dashed/dotted <b>NOT</b> a single line lone pair <b>MUST BE</b> in line with hydrogen bond</p>
1 d	more hydrogen bonds ✓	1	<p><b>ALLOW</b> more groups/sites/places which can form H bonds, more electronegative atoms/N and O atoms which can form H bonds <b>DO NOT ALLOW</b> more Os which can form H bonds, bonds more easily</p>
		8	

Question	Answer	Mark	Guidance
2 a i	ether ✓	1	
2 a ii	 <p>either or both circled as above ✓</p>	1	<b>ALLOW</b> if adjacent C is also circled
2 a iii	 <p>both circled OH groups present ✓ full structural formula correct ✓</p>	2	<b>AWARD</b> 1 mark if one of the circled OH groups is incorrect but rest of structure is correct
2 b i	ethanoic acid ✓	1	

Question	Answer	Mark	Guidance
2 b ii	<p><b>cellulose triethanoate:</b>  <u>instantaneous dipole-induced dipole</u> &amp; <u>permanent dipole-permanent dipole</u></p> <p><b>polyester:</b>  <u>instantaneous dipole-induced dipole</u> &amp; <u>permanent dipole-permanent dipole</u></p> <p>marking:  <u>instantaneous dipole-induced dipole</u> in both polymers ✓  <u>permanent dipole- permanent dipole</u> in both polymers ✓</p>	2	<p><b>ALLOW</b> van der Waals for <u>instantaneous dipole-induced dipole</u></p> <p><b>DO NOT ALLOW</b> abbreviations</p> <p><b>DO NOT ALLOW</b> instantaneous-induced dipole forces unless they have used <u>instantaneous dipole-induced dipole</u> once</p> <p><b>ALSO</b> applies to permanent dipoles</p> <p>hydrogen bonding is a <b>CON</b> so does not gain pd-pd mark</p>
2 b iii	<p>longer chains / length of chains ✓  more intermolecular bonds/forces in longer chains <b>ORA</b> ✓</p> <p><b>OR</b>  polymer molecules/chains closer ✓  intermolecular bonds stronger ✓ <b>ORA</b></p> <p><b>OR</b>  <b>shorter</b> monomer chains ✓  more intermolecular bonds between polymer chains /  intermolecular bonds more frequent along polymer chains ✓</p>	2	<p><b>NOTE</b>  amount of imb/fs per unit length  will get 1 mark, needs to relate 'more' to 'stronger' for 2nd mark</p> <p><b>NOT</b> how closely, more imbs</p> <p><b>may be inferred:</b> linkages closer together</p> <p><b>IGNORE</b> references to branch/side groups, crystallinity and any other factors</p>
2 c	<p>at temperatures <u>below</u> polymer's T<sub>g</sub> ✓  chains / molecules cannot move/slide over one another ✓  chains / molecules break when force applied ✓</p>	3	<p><b>IGNORE</b> references to polymers being brittle</p> <p><b>NOT</b> polymers break/shatter  chains/molecules may be implied by use of 'they' referring to polymer chains</p>

Question	Answer	Mark	Guidance
2 d	peak around 3500/3600 / in range 3200-3600 indicates O-H / hydroxyl bond in alcohol ✓  so <b>cellulose diethanoate</b> since only 2 of the 3 OH groups in repeating unit of cellulose have reacted / one OH / no OH groups in cellulose triethanoate ora AW ✓	2	<b>IGNORE</b> references to no O-H peak in range 2500-3200  <b>MUST</b> relate OH group to answer <b>DO NOT AWARD</b> this mark if a COOH group is also given as present
		14	




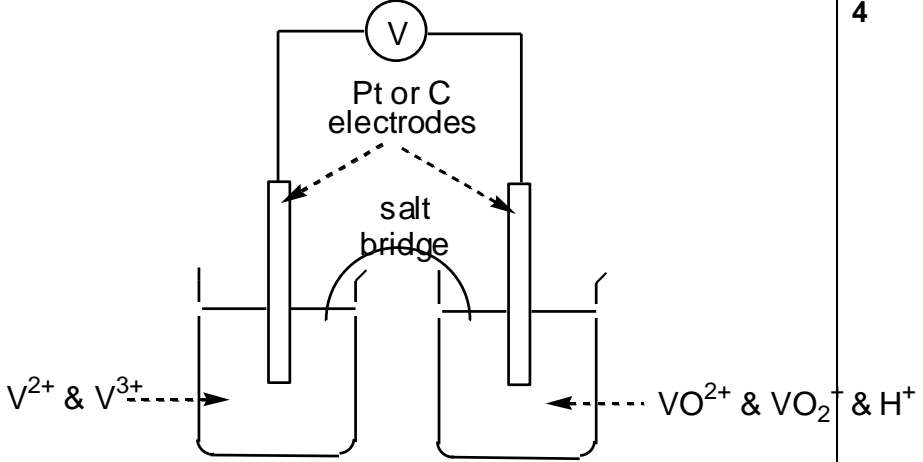
Question	Answer	Mark	Guidance						
3 a	<table border="1"> <thead> <tr> <th data-bbox="367 272 651 336">test</th> <th data-bbox="651 272 1140 336">observations</th> </tr> </thead> <tbody> <tr> <td data-bbox="367 336 651 536"> <b>For C</b>            (add neutral)            iron(III) chloride            (solution) / <math>\text{FeCl}_3</math> ✓         </td> <td data-bbox="651 336 1140 536">purple colour formed ✓</td> </tr> <tr> <td data-bbox="367 536 651 695"> <b>For D</b>            (add) <math>\text{Na}_2\text{CO}_3</math> (s)            or (aq) ✓         </td> <td data-bbox="651 536 1140 695">solution / mixture fizzes / bubbles / effervesces / (colourless) gas given off ✓</td> </tr> </tbody> </table>	test	observations	<b>For C</b> (add neutral) iron(III) chloride (solution) / $\text{FeCl}_3$ ✓	purple colour formed ✓	<b>For D</b> (add) $\text{Na}_2\text{CO}_3$ (s) or (aq) ✓	solution / mixture fizzes / bubbles / effervesces / (colourless) gas given off ✓	4	<p><b>DO NOT ALLOW</b> iron chloride  <b>ALLOW</b> violet <b>NOT</b> blue or red</p> <p><b>ALLOW</b> <math>\text{NaHCO}_3</math> / K for Na / Ca <math>\text{CO}_3</math> / <math>\text{MgCO}_3</math> / carbonate</p>
test	observations								
<b>For C</b> (add neutral) iron(III) chloride (solution) / $\text{FeCl}_3$ ✓	purple colour formed ✓								
<b>For D</b> (add) $\text{Na}_2\text{CO}_3$ (s) or (aq) ✓	solution / mixture fizzes / bubbles / effervesces / (colourless) gas given off ✓								
3 b i	 <p>1 mark for each correct structure ✓✓</p>	2	<p>Any clear structure acceptable  <b>ALLOW</b> <math>\text{NH}_3\text{Cl}</math>. <math>\text{NH}_3^+</math></p> <p><b>DO NOT ALLOW</b> an OH group to be bonded to the ring via -HO</p>						

Question	Answer	Mark	Guidance
3 b ii	<p>draw <b>pencil-line</b> near bottom of <b>plate/sheet</b> ✓</p> <p>place 1 <b>drop</b> (or similar word) <b>of mixture AND a drop of each</b> of the 2 compounds (on the line) AW ✓</p> <p>place plate in solvent, <b>line above solvent</b> level <b>AND add lid/cover</b> ✓</p> <p>when solvent nears <b>top of plate</b> AW, <b>remove/dry</b> plate ✓</p> <p>locate spots with/ expose to <b>UV light/iodine</b> ✓</p> <p><b>compare</b> heights/position of spots from mixture with the two compounds AW  <b>OR</b>  <b>calculate/measure/determine</b> <math>R_f</math> values of spots and <b>compare</b> with those of the two compounds AW ✓</p>	6	<p><b>please annotate marks given with ticks</b>  <b>ALL</b> marking points may be gained from labelled diagram(s)</p> <p>'near bottom' may be implied by what follows  <b>DO NOT ALLOW</b> paper for plate/sheet <b>BUT</b> ecf for further use  <b>ALLOW</b> draw base-line with pencil</p> <p>any other named locating agent is a <b>CON</b>  <b>DO NOT ALLOW</b> 'locating agent' alone  <b>DO NOT ALLOW</b> vague statements about comparing spots, <b>MUST</b> refer to positions</p> <p><b>DO NOT ALLOW</b> vague statements about <math>R_f</math> values  e.g. <math>R_f</math> values will identify compounds  <b>MUST</b> indicate that spot heights or <math>R_f</math> values have been measured <b>AND</b> compared</p>

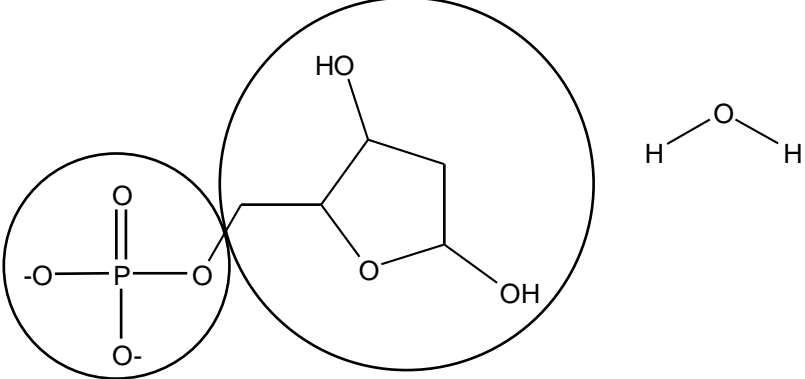
Question	Answer	Mark	Guidance
3 c i	for concentration to half: 1. $4.96 \times 10^{-3}$ to $2.48 \times 10^{-3} = 220-30 = \mathbf{190 \text{ hours}}$ ✓ 2. $3.83 \times 10^{-3}$ to $1.92 \times 10^{-3} = 290-100 = \mathbf{190 \text{ hours}}$ ✓ half-life constant (means 1 <sup>st</sup> order) ✓ <b>OR</b> every 70 hours, 1. from 30-100 hours = about <b>23%</b> of conc. Is used up ✓ 2. from 220-290 hours = about <b>23%</b> of conc. Is used up ✓ same proportion of starting conc. used up (means 1 <sup>st</sup> order) ✓ <b>OR</b> every 70 hours (a set time interval) for 1 <sup>st</sup> order the concentration drop will be a constant ratio ✓ 1. from 30-100 hours = about <b>1.295</b> ✓ 2. from 220-290 hours = about <b>1.292</b> ✓	3	units required for half-lives at least once <b>otherwise 1</b> mark only for both times correct  <b>DO NOT ALLOW</b> half-life <b>MARK</b> if no relevant data given  <b>IGNORE</b> units  <b>IGNORE</b> units  <b>NOTE</b> any data referring to actual rates cannot be relevant/meaningful
3 c ii	average rate of reaction = $(\mathbf{5.55 - 1.92}) \times 10^{-3} / 290$ ✓ $= 1.25 \times 10^{-5}$	1	Give mark for getting the working correct, may make an error with calculator <b>IGNORE</b> units for rate of reaction
3 c iii	$k = 4.96 \times 10^{-9} / 4.96 \times 10^{-3} = 1.0 \times 10^{-6}$ ✓ s <sup>-1</sup> ✓	2	<b>ALLOW</b> $1.00 \times 10^{-6}$ <b>or</b> $1 \times 10^{-6}$ <b>or</b> $10^{-6}$
3 d i	change/alter/different functional/side groups (in structure / formula) ✓	1	<b>ALLOW</b> add / remove group(s) <b>IGNORE</b> references to altering shape, specific named groups
3 d ii	make a large number of related compounds (together quickly) AW ✓	1	<b>DO NOT ALLOW</b> 'test' instead of 'make' <b>MUST</b> refer to a large number in some way e.g. many <b>NOT</b> just 'compounds'
		<b>20</b>	

Question	Answer	Mark	Guidance										
4 a i	sulfur dioxide / sulfur trioxide <b>OR</b> chlorine ✓	1	<b>IGNORE</b> formulae  <b>ALLOW</b> hydrogen chloride <b>NOT</b> hydrochloric acid or sulfur (di)chloride										
a ii	<table border="1" data-bbox="369 363 833 726"> <thead> <tr> <th>compound</th> <th>oxidation state of vanadium</th> </tr> </thead> <tbody> <tr> <td>VS<sub>4</sub></td> <td>+4</td> </tr> <tr> <td>NaVO<sub>3</sub></td> <td>+5</td> </tr> <tr> <td>Na<sub>2</sub>V<sub>6</sub>O<sub>16</sub></td> <td>+5</td> </tr> <tr> <td>V<sub>2</sub>O<sub>5</sub></td> <td>+5</td> </tr> </tbody> </table> VS <sub>4</sub> correct ✓ rest correct ✓	compound	oxidation state of vanadium	VS <sub>4</sub>	+4	NaVO <sub>3</sub>	+5	Na <sub>2</sub> V <sub>6</sub> O <sub>16</sub>	+5	V <sub>2</sub> O <sub>5</sub>	+5	2	<b>ONLY</b> penalise <b>ONCE</b> for lack of sign sign must be before number, ecf after first error
compound	oxidation state of vanadium												
VS <sub>4</sub>	+4												
NaVO <sub>3</sub>	+5												
Na <sub>2</sub> V <sub>6</sub> O <sub>16</sub>	+5												
V <sub>2</sub> O <sub>5</sub>	+5												
a iii	<b>step 1</b> ✓  oxidation state of V changes/increases ✓	2	more than step 1 is a <b>CON</b> but mark explanation separately  <b>ALLOW</b> electron loss by V <b>ALLOW ecf</b> for the 2 <sup>nd</sup> mark, from wrong oxidation state for VS <sub>4</sub> in table in <b>aii</b> , e.g. +8 (for VS <sub>4</sub> ) to +5, so oxidation state decreases <b>If two steps in first part</b> , both reasons must refer correctly to the data in <b>aii</b>										


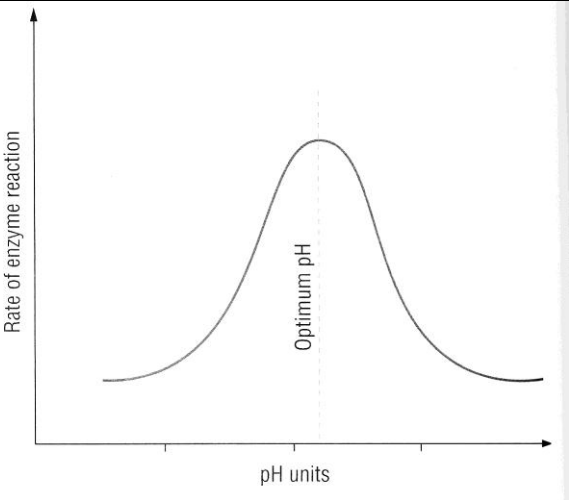
Question	Answer	Mark	Guidance
a iv	<p><b>absorb</b> (specific) <b>frequencies</b> (or wavelengths) in (specific) parts of the <b>visible</b> spectrum ✓</p> <p> <b>absorb</b> must be spelled correctly to gain this mark</p> <p><b>transmit</b> complementary colour / frequencies (or wavelengths) not absorbed / yellow light ✓</p>	2	<p><b>MUST</b> use frequency/frequencies/wavelength(s) for 1<sup>st</sup> mark  <b>IGNORE</b> any reference to energy levels and electrons  <b>ALLOW</b> 'light' for visible</p> <p><b>or</b> any of its variants e.g. absorbed, absorbing, absorption etc.</p> <p><b>ALLOW</b> only complementary colour / frequencies (or wavelengths) can be seen  <b>DO NOT ALLOW</b> reflect / emit / absorption</p>
a v	$\text{V}_2\text{O}_5 + 5\text{Ca} \rightarrow 2\text{V} + 5\text{CaO} \checkmark$	1	<p><b>IGNORE</b> state symbols</p>
b	<p>cooling / lowering / controlling temperature (of contents of furnace) ✓</p>	1	<p><b>ALLOW</b> <i>statements which infer cooling e.g.</i> absorbs heat, prevents furnace getting too hot / thermal shock  <b>IGNORE</b> references to cost, landfill</p>

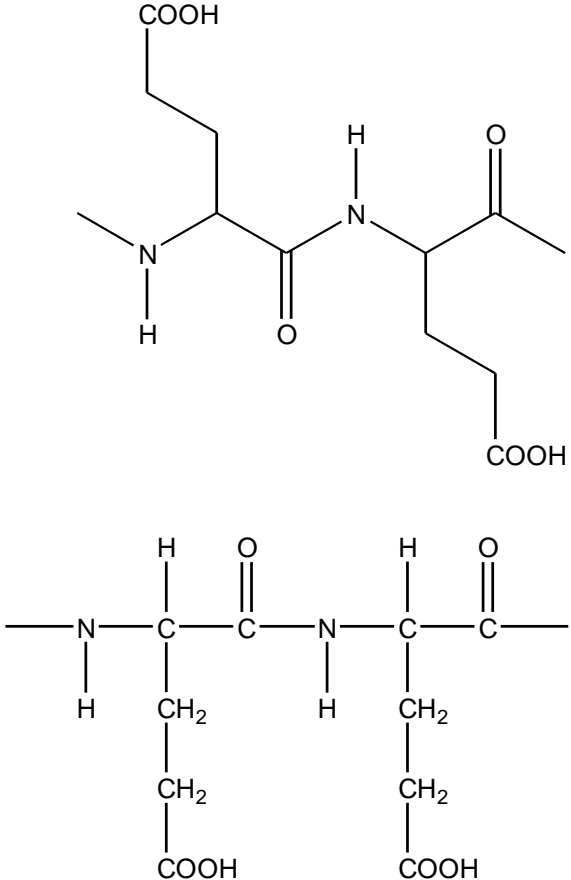
Question	Answer	Mark	Guidance
c i	 <p><i>diagram shows minimum requirements for 4 marks</i></p> <p>both electrodes correct ✓  correct vanadium species in each cell ✓  H<sup>+</sup> ions in VO<sup>2+</sup> and VO<sub>2</sub><sup>+</sup> solution ✓  salt bridge labelled &amp; dipping into solutions <b>AND</b>  voltmeter attached correctly <b>AND</b> circuit complete ✓</p>	4	<p><b>ALLOW</b> half-cells reversed</p> <p><b>IGNORE</b> anions, charges on electrodes, water, temperature and concentrations</p>
c ii	$E_{\text{cell}} = 1.26 \text{ V}$	1	<b>IGNORE</b> any sign
c iii	<p>temperature is not standard / 25°C ✓</p> <p>concentrations of ions in a half-cell are not equal ✓</p>	2	<p><b>DO NOT ALLOW</b> conditions not standard, must be specific to temperature and/or concentration</p> <p><b>ALLOW</b> concentration not 1 mol dm<sup>-3</sup></p> <p><b>DO NOT ALLOW</b> 1 mol(e) for concentration</p>
c iv	$\text{VO}_2^+ + 2\text{H}^+ + \text{V}^{2+} \rightarrow \text{VO}^{2+} + \text{H}_2\text{O} + \text{V}^{3+}$ vanadium species correct ✓ equation correct ✓	2	<b>GIVE 1 mark</b> if equation has species and balancing correct but is reversed

Question	Answer	Mark	Guidance
c v	$\text{VO}_2^+$ <b>AND</b> $E^\ominus(\text{I}_2/\text{I}^-)$ must be more negative/less positive than the V ion being reduced <b>ORA</b> ✓  $2\text{VO}_2^+ + 4\text{H}^+ + 2\text{I}^- \rightarrow 2\text{VO}^{2+} + 2\text{H}_2\text{O} + \text{I}_2$ ✓	2	<b>ORA</b> V half-cell is the only one with a more positive $E^\ominus$ than the $\text{I}_2/\text{I}^-$ half-cell  <b>ALLOW</b> balanced equation with $\frac{1}{2} \text{I}_2$ <b>IGNORE</b> state symbols
d i	moles of $\text{Cr}_2\text{O}_7^{2-}$ used = <b>0.02160 x 23.50 / 1000</b> = 0.0005076 (5.076 x $10^{-4}$ ) ✓  moles of $\text{Fe}^{2+}$ reacted = <b>6 x 0.0005076</b> ✓ = 0.0030456 (3.0456 x $10^{-3}$ )  mass of Fe in alloy = 0.0030456 x <b>55.8</b> ✓ = 0.16994448 g  % Fe = 0.16994448 x <b>100 / 0.1750</b> ✓  = <b>97.1</b> % ✓	5	<b>DO NOT award marks for random numbers without any explanation to what they refer to</b> marks are for working out shown in bold <b>OR</b> actual answers at each stage  <b>must be 3 sig. figs.</b> 97% gains 4 marks, 97.1% 5 marks, irrespective of working
d ii	<b>oxygen / air</b> (and water) will oxidise/change $\text{Fe}^{2+}$ (to $\text{Fe}^{3+}$ ) ✓  $\text{NaHCO}_3$ reacts with acid to form $\text{CO}_2$ ✓  air/gas is dispelled which cannot return AW ✓	3	
e i	<b>green solution</b> ✓ forms a <b>green precipitate</b> ✓	2	<b>IGNORE</b> any qualifying of green such as pale, dark, dirty, rust etc. <b>ALLOW</b> ppt <b>ALLOW</b> solid for precipitate
e ii	$\text{Fe}^{2+}(\text{aq}) + 2\text{OH}^-(\text{aq}) \rightarrow \text{Fe}(\text{OH})_2(\text{s})$ formulae correct & balanced ✓ state symbols correct for precipitation reaction ✓	2	
		32	

Question	Answer	Mark	Guidance
5 a	<p>1. to find out about their relatives / ancestors AW  <b>OR</b>  to diagnose illness / cure disease AW ✓</p> <p>2. they are police / criminal suspects AW ✓</p> <p>3. they have not been prosecuted  <b>OR</b> they are innocent  <b>OR</b> they have been found not guilty  <b>OR</b> infringes privacy  <b>OR</b> prevents access by other people AW ✓</p>	3	<p><b>NOTE</b> they may have answered in a different order to the questions in the stem  <b>DO NOT ALLOW</b> to be cleared of a crime AW</p> <p><b>IGNORE</b> references to human rights/liberty</p>
b i	nucleotide ✓	1	<b>IGNORE</b> any details of the constituents of nucleotides
b ii	 <p>phosphate structure (circled) correct ✓  deoxyribose structure (circled) correct ✓</p> <p>water structure ✓</p>	3	<p><b>ALLOW</b> an OH group for an O<sup>-</sup> on phosphate</p> <p>The phosphate <b>MUST</b> be joined at the primary OH</p> <p><b>ALLOW</b> H<sub>2</sub>O or any bond angle for water</p>



Question	Answer	Mark	Guidance
c	<p>base pairs held together by <u>hydrogen bonding/bonds</u> ✓  adenine-thymine 2 hydrogen bonds <b>AND</b>  guanine-cytosine 3 hydrogen bonds ✓</p> <p><b>hydrogen bonds</b> between base pairs <b>break</b> ✓</p> <p>(two single) <b>helices / strands</b> are formed ✓</p> <p><b>each base</b> (on these helices/strands) forms <b>hydrogen bonds</b>  to a <b>new (correct/complementary) base</b> AW ✓</p> <p> to gain this last point replication must be linked to  breaking <b>AND</b> forming hydrogen bonds</p>	5	<p>May be implied by stating the number of hydrogen bonds  in each case  different numbers of H bonds is <b>NGE</b></p> <p>May refer to just <b>one strand</b></p> <p><b>IGNORE</b> references to phosphodiester bonds</p> <p><i>i.e.</i> both forming and breaking H bonds have to be  mentioned in the answer</p>
d	 <p>curve correct shape ✓  label for <b>optimum pH</b> in line with <b>fastest rate of reaction</b> ✓</p>	2	<p><b>ALLOW</b> any curve with a peak, not necessary to be  symmetrical</p> <p>Optimum pH should indicate the <b>peak</b> of graph <b>AND</b> be  labelled as such</p>

Question	Answer	Mark	Guidance
e i	<p>structure correct - 2 examples shown below</p>  <p>✓</p>	1	<p><b>ALLOW</b> any correct structure  <b>ALLOW</b> without 'spare bonds'  <b>ALLOW</b> dipeptide structure which must have its secondary amide/peptide between two chiral C atoms</p>
e ii	<p>((side) <b>-COOH/carboxyl groups</b> will (lose protons and))  form <b>-COO<sup>-</sup>/carboxylate ions</b> ✓</p>	1	<p>also forming <math>\text{NH}_2^+</math> / <math>\text{NH}_3^+</math> is a <b>CON</b>  <b>ALLOW</b> carboxyl/COOH becomes deprotonated AW</p>
		16	

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