

**Chemistry B (Salters)**

Advanced GCE

Unit **F334**: Chemistry of Materials

**Mark Scheme for January 2013**

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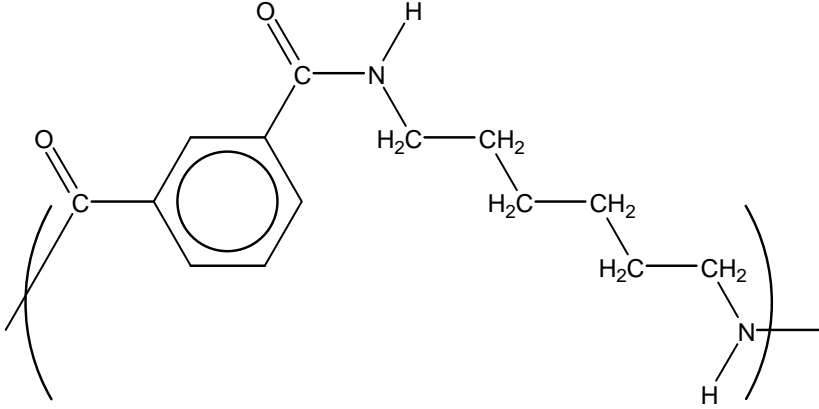
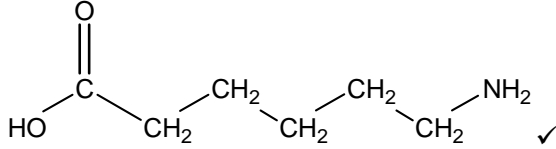
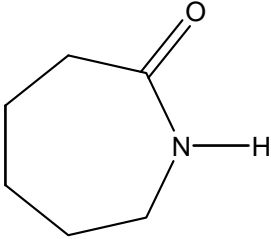
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**Annotations** used in the detailed Mark Scheme (to include abbreviations and subject-specific conventions).

Annotation	Meaning
/	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
—	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:

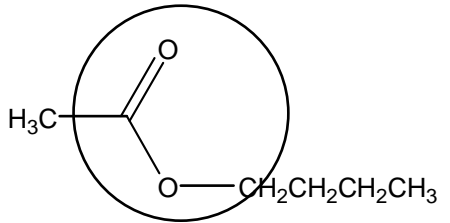
Annotation	Meaning
	correct response
	incorrect response
	benefit of the doubt
	benefit of the doubt <b>not</b> given
	error carried forward
	information omitted
	Ignore
	Reject

Question		Answer	Marks	Guidance
1	(a) (i)	 <p>all correct 2 marks ✓✓ amide link only ✓</p>	2	<p><b>ALLOW</b></p> <ul style="list-style-type: none"> <li>any correct structural formula including skeletal</li> <li>mixtures of structural and skeletal</li> <li>-CONH-</li> <li>Cs and Hs on ring</li> </ul> <p><b>IGNORE</b> brackets and n etc</p>
	(ii)	-CONH- circled ✓	1	<b>ALLOW</b> adjacent C atoms in circle
	(iii)	1,6-diamino ✓ hexane ✓ <b>OR</b> hexane ✓ -1,6-diamine ✓	2	If butane <b>ALLOW</b> 1,4-diamino for ecf mark <b>IGNORE</b> commas and dashes <b>ALLOW</b> 1,6-hexanediamine
	(b) (i)	 <p>✓</p>	1	<p><b>ALLOW</b> any formula that makes structure clear</p> <p><b>ALLOW</b> cyclic amide</p> 

Question		Answer	Marks	Guidance
	(ii)	condensation <b>AND</b> water is eliminated/formed/lost ✓  <b>IF</b> cyclic amide in 1b(i) <b>THEN</b> addition <b>AND</b> hydrolysis (of ring) / water added	1	'addition' or 'addition polymerisation' is a <b>CON</b> <b>ALLOW</b> 'small molecule' instead of water any other named small molecule is a <b>CON</b>
(c)	(i)	(polymer) <u>chains/molecules</u> are (highly) ordered/aligned <b>AW</b> ✓	1	<b>ALLOW</b> labelled diagram that shows alignment of chains by using parallel lines  <b>AW</b> means other suitable phrases eg 'stacked closely and neatly' 'arranged regularly'/'regularity of chains'
	(ii)	PPA chains are <b>closer</b> together <b>ORA</b> ✓  so <b>intermolecular bonds</b> in PPA will be <b>stronger</b> <b>OR</b> more hydrogen bonds in PPA <b>OR</b> more intermolecular bonds in PPA <b>ORA</b> ✓  <b>more energy/heat</b> will be required to break the <b>intermolecular bonds</b> <b>OR</b> melt the polymer <b>OR</b> allow chains to move over each other <b>ORA</b> ✓	3	<b>AS ALWAYS</b> intermolecular <b>bonds</b> is synonymous with intermolecular <b>forces</b>  <b>NOT</b> more areas of contact <b>NOT</b> just compacted together <b>NOT</b> just closely packed <b>NOTE</b> ALL 3 marking points are <b>comparative</b>  <b>ALLOW</b> any named intermolecular bond  <b>NOT</b> 'higher temperature' for 'energy'
	(iii)	cold-drawing ✓	1	<b>NOT</b> co-polymerising <b>IGNORE</b> 'necking'

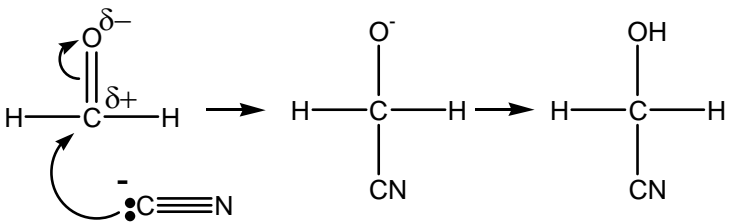
Question		Answer	Marks	Guidance
	(d) (i)	1. elimination ✓ 2. addition ✓ 3. addition ✓	3	<b>IGNORE</b> 'nucleophilic'
	(ii)	can be used in step 3 <b>AW</b> ✓	1	<b>NOT</b> step 2 <b>ALLOW</b> <ul style="list-style-type: none"> <li>• can be sold</li> <li>• can be used as a source of energy</li> <li>• can be recycled</li> </ul>
	(e)	acidified dichromate ✓	1	<b>IGNORE</b> <ul style="list-style-type: none"> <li>• metal cation</li> <li>• name of acid</li> <li>• wrong formulae if name given</li> <li>• oxidation state of 'dichromate'</li> </ul> <b>ALLOW</b> H <sub>2</sub> SO <sub>4</sub> /H <sup>+</sup> and Cr <sub>2</sub> O <sub>7</sub> <sup>2-</sup> <b>DO NOT ALLOW</b> just dichromate
		<b>Total</b>	<b>17</b>	

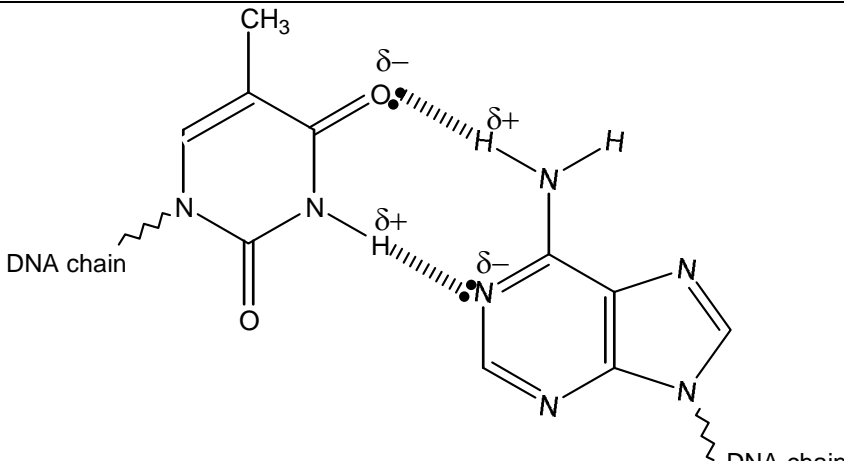
Question		Answer	Marks	Guidance
2	(a)	-COOH / carboxyl / carboxylic acid group ✓ is a proton/H <sup>+</sup> donor / loses H <sup>+</sup> ✓	2	correct equation showing dissociation gains both marks <b>ALLOW</b> 'gives H <sup>+</sup> ' <i>Mark separately</i>
	(b) (i)	CH <sub>2</sub> OHCOOH + NaOH → CH <sub>2</sub> OHCOONa + H <sub>2</sub> O ✓	1	<b>ALLOW</b> CH <sub>2</sub> OHCOO <sup>-</sup> Na <sup>+</sup> , Na <sup>+</sup> CH <sub>2</sub> OHCOO <sup>-</sup> <b>OR</b> Na(CH <sub>2</sub> OHCOO) <b>IGNORE</b> state symbols <b>NOT</b> CH <sub>2</sub> OHNaCOO
	(ii)	1. moles of NaOH used in titration = <b>16.00/1000</b> x <b>0.250</b> ✓ = 0.00400  2. moles of glycolic acid used in titration = answer from <b>1</b> , <b>scaled by ratio in equation in 2bi</b> ✓ 0.00400  <b>CHECK equation in 2(b)(i)</b>  3A. moles of glycolic acid in 250 cm <sup>3</sup> = (answer from <b>2</b> ) x <b>10</b> ✓ = 0.0400 mol <b>OR</b> 3B. mass of glycolic in 25 cm <sup>3</sup> = (answer from <b>2</b> )x <b>M<sub>r</sub></b> of glycolic acid ✓ 0.304 g <b>OR</b> 3C. concentration of glycolic acid = (answer from <b>2</b> ) x <b>1000/25</b> ✓ 0.16 mol dm <sup>-3</sup>  4. <b>M<sub>r</sub></b> of glycolic acid = <b>76.(0)</b> ✓	6	<i>There are several possible routes through this question after point 2, the 'mole route' A, the 'mass route' B and the 'concentration route' C</i>  <b>If final answer is incorrect please annotate with ticks where the marks are awarded</b>  2. <b>ALLOW</b> by implication if 0.004 used subsequently  <b>The marks are awarded for the working out given in bold OR</b> the correctly calculated answer to that working (but no mark if calculated answer is shown and is wrong)  4. Award if 76 used correctly anywhere

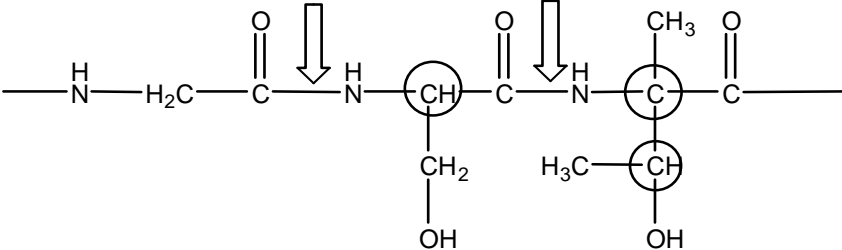
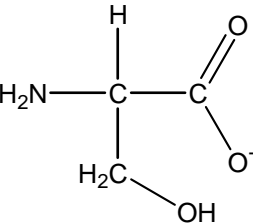
Question	Answer	Marks	Guidance
	<p>5A. moles of glycolic acid in 100 cm<sup>3</sup> = (answer from <b>3A</b>) x 100/14 ✓ 0.286 mol</p> <p><b>OR</b></p> <p>5Bi. mass of glycolic acid in 250 cm<sup>3</sup> = (answer from <b>3A</b>) x <i>M<sub>r</sub></i> of glycolic acid ✓ 3.04 g</p> <p><b>OR</b></p> <p>5Bii. mass of glycolic in 250 cm<sup>3</sup> = (answer from <b>3B</b>) x <b>10</b> = ✓ 3.04 g</p> <p><b>OR</b></p> <p>5C. concentration of undiluted glycolic acid = (answer from <b>3C</b>) x <b>250/14</b> ✓ 2.86 mol dm<sup>-3</sup></p> <hr/> <p>6. mass of glycolic acid in 100 cm<sup>3</sup> undiluted = (answer from <b>5A</b>) x <i>M<sub>r</sub></i> of glycolic acid</p> <p><b>OR</b> = (answer from <b>5B</b>) x <b>100/14</b></p> <p><b>OR</b> = (answer from <b>5C</b> /10) x <i>M<sub>r</sub></i> of glycolic acid</p> <p>= <b>21.7</b> (3 sf) <b>AND</b> correct comment ✓</p>		<p><b>ALLOW</b> ecf for incorrect equation <b>AND</b> between each step</p> <p><b>ALLOW</b> 'Acnegone' for 'glycolic acid'</p> <p><b>ALLOW</b> answers in standard form</p> <p><i>The following on the answer line with correct corresponding comment, score as follows, irrespective of working or lack of it:</i></p> <p>21.7 scores 6</p> <p>1.22 scores 5 (error in 5C)</p> <p>12.2 scores 5 (error in 6C)</p> <p>2.17 scores 5 (error in 3A or 5Bii)</p> <p>3.04 scores 5 (error in 6)</p> <p><i>these to other sf OR with incorrect comment score one mark less</i></p> <p><b>If one of the answers above applies place correct number of ticks by answer</b></p>
(c) (i)	 <p>butyl ethanoate ✓ ester link correct ✓</p>	2	<p><b>IGNORE</b> where the circle cuts the bond as long COO is included</p> <p><b>ALLOW</b> adjacent C atoms in circle</p> <p><b>ALLOW</b> butylethanoate without gap</p>
(ii)	<p>butan-1-ol ✓</p> <p><u>concentrated</u> sulfuric acid/hydrochloric acid ✓</p>	2	<p><b>MUST HAVE</b> number 1</p> <p><b>DO NOT ALLOW</b> ecf for alcohol in 2(c)(i)</p> <p><b>ALLOW</b> formula for acid <b>ONLY</b></p> <p><b>IGNORE</b> spelling of name for H<sub>2</sub>SO<sub>4</sub> as long as it is clear</p>

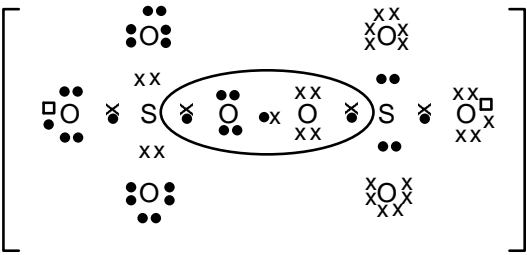


Question		Answer	Marks	Guidance
	(iii)	<p><b>ethanoic acid:</b> hydrogen bonding ✓ <b>IGNORE</b> pd-pd</p> <p><b>compound D:</b> pd-pd ✓ hydrogen bonding is <b>CON</b></p> <p>instantaneous dipole – induced dipole / id-id in both ✓</p>	3	<p><b>ALLOW</b> 'pd-pd' / id-id abbreviations here</p> <p><b>ALLOW</b> Van der Waals for id-id</p>
(d)	(i)	<p><i>The answer requires a comment for a carboxylic acid AND an alcohol so award ONE mark from each section below:</i></p> <p><b>CARBOXYLIC ACID</b>            (strong) peak at 1743 (<math>\text{cm}^{-1}</math>) shows C=O in ester not acid  <b>OR</b> no peak 1700 – 1725 (<math>\text{cm}^{-1}</math>) shows no C=O in acid  <b>OR</b> no <u>broad</u> peak at 2500 – 3200 (<math>\text{cm}^{-1}</math>) shows no O–H in acid ✓</p> <p><b>ALCOHOL</b>            no peak greater than 3000 / in range 3200 – 3600 (allow 3640) so            no O–H in alcohol ✓</p>	2	<p>answers may be given on spectrum</p> <p>to score each point, <b>range</b>, <b>bond</b> and <b>group</b> in which it is found must be given</p> <p><b>ALLOW</b> carboxyl or carboxylic acid or COOH or ethanoic acid (or formula) for '<b>acid</b>'</p> <p>FOR O–H <b>ALLOW</b> OH / hydroxyl            FOR C=O <b>ALLOW</b> carbonyl <b>NOT</b> CO</p>
	(ii)	<p><b>peak at <math>m/z</math> 73:</b>  <math>\text{CH}_3\text{COOCH}_2</math> / <math>\text{C}_3\text{H}_5\text{O}_2</math> ✓            positive charge on any formula ✓</p> <p><b>species lost:</b>  <math>\text{CH}_2\text{CH}_2\text{CH}_3</math> / <math>\text{C}_3\text{H}_7</math> (NO charge) ✓</p>	3	<p><b>ALLOW</b> any correct structural or molecular formula for both answers</p> <p><b>ALLOW</b> <math>\text{C}_4\text{H}_9\text{O}^+</math></p> <p><b>IF</b> <math>\text{C}_4\text{H}_9\text{O}^+</math> given above  <b>THEN species lost must be</b> <math>\text{C}_2\text{H}_3\text{O}</math> / <math>\text{CH}_3\text{CO}</math></p>
(e)	(i)	nucleophilic ✓ addition ✓	2	

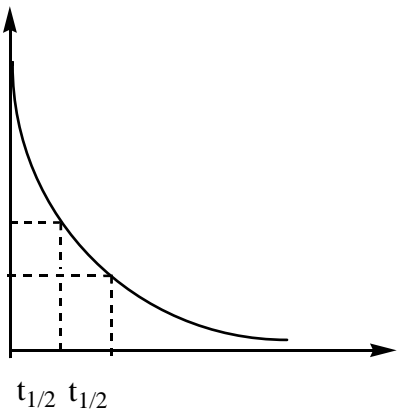
Question	Answer	Marks	Guidance
(ii)	<p style="text-align: center;"><math>\text{H}^+</math> or <math>\text{H}_2\text{O}</math> or <math>\text{HCN}</math></p>  <p>HCHO with bond polarity correct ✓  CN<sup>-</sup> with lone pair of electrons on C ✓</p> <p>both curly arrows correct (lower one must be from C of CN) ✓  intermediate ion correct ✓  gain of H<sup>+</sup> from <b>water, HCN</b> or <b>direct</b> to form the cyanohydrin ✓</p>	5	<p><b>ALLOW</b> CN<sup>-</sup> without triple bond  <b>ALLOW</b> -ve charge on N of CN<sup>-</sup></p> <p><b>IGNORE</b> any arrow used to add H<sup>+</sup> etc to the intermediate</p>
	<b>Total</b>	<b>28</b>	

Question		Answer	Marks	Guidance
3	(a) (i)	repeating unit in DNA is a <b>nucleotide</b> ✓ formed from pentose/ <b>sugar</b> /deoxyribose, <b>phosphate</b> and a <b>base</b> AW ✓	2	<i>mark separately</i> <b>ALLOW</b> • names for the four bases (T,A,C,G) instead of 'bases' • a named example of a base <b>IGNORE</b> ribose for <b>two marks</b> it must be clear that the sugar etc makes up the nucleotide
	(ii)	proteins are formed from <b>amino acid(s)</b> ✓	1	
	(iii)	proteins have <b>more</b> amino acids than the <b>bases/nucleotides</b> in DNA ✓ from which to construct many unique/different/more structures/arrangements/combinations to carry genetic data AW ✓	2	<b>ALLOW</b> actual numbers eg there are 20/21 amino acids in proteins and only 4 bases/nucleotides in DNA
	(b)	 <p>correct formula completed for adenine <b>AND</b> both H bonds between correct atoms ✓ correct polarities for one hydrogen bond ✓ lone pair for one hydrogen bond ✓</p>	3	

Question		Answer	Marks	Guidance
(c)	(i)	 <p>ANY TWO of the THREE CORRECT chiral Cs ONLY ✓</p>	1	
	(ii)	arrows as in (d) (i) ✓	1	<b>IGNORE</b> extra arrows pointing to C–N bonds at ends of chain. Any other arrows are a <b>CON</b>
	(iii)	 <p>correct amino acid (see opposite) ✓</p> <p>correct formula of ion ✓</p>	2	<b>IGNORE</b> any cations <b>ALLOW</b> any correct structural formula <b>IGNORE</b> species formed from amino acids to right and left <b>IGNORE</b> <ul style="list-style-type: none"> <li>added H<sup>+</sup> to –NH<sub>2</sub></li> <li>negative charge on alcohol i.e. –O<sup>–</sup></li> <li>–COOH instead</li> </ul> <b>NO</b> ecf <b>IF</b> anion formed at CH <sub>2</sub> O <sup>–</sup> then 1 mark max for amino acid
	(iv)	secondary: folding/twisting of polypeptide/amino acid chains/primary structure ✓  tertiary: further/final folding <b>OR</b> 3D shape/structure ✓	2	<b>ALLOW</b> (alpha)-helix/coiled <b>OR</b> (beta)-sheets / pleated sheets <b>ALLOW</b> overall/global structure
<b>Total</b>			<b>14</b>	

Question		Answer	Marks	Guidance
4	(a) (i)	 <p>correct inside oval (around peroxy Os) ✓ rest correct ✓</p>	2	<p><b>ALLOW</b> another symbol for S electrons</p> <p>second mark depends on first</p>
	(ii)	$(\text{NH}_4)_2 \text{S}_2\text{O}_8$ ✓	1	<p><b>ALLOW</b></p> <ul style="list-style-type: none"> <li>with correct charges</li> <li><math>\text{S}_2\text{O}_8(\text{NH}_4)_2</math></li> </ul>
	(b)	<p><math>\text{S}_2\text{O}_8^{2-}</math> (is the stronger oxidising agent) because <math>\text{S}_2\text{O}_8^{2-}</math> has a <b>more positive</b> <math>E^\ominus</math> value ✓ indicates a <b>greater</b> tendency AW to gain/attract/accept electrons/to be reduced ✓</p> <p><b>ORA ie:</b> <math>\text{S}_2\text{O}_8^{2-}</math> (is the stronger oxidising agent) because <math>\text{Cr}_2\text{O}_7^{2-}</math> has a <b>less positive / more negative</b> <math>E^\ominus</math> value ✓</p> <p>indicates a <b>smaller</b> tendency AW to gain/attract/accept electrons/to be reduced <b>OR</b> indicates a <b>greater</b> tendency AW to lose electrons/to be oxidised ✓</p>	2	<p>without reference to oxidising strength of <math>\text{S}_2\text{O}_8^{2-}</math> only scores 1 as it doesn't answer the question</p> <p><b>NOT</b> 'higher' for 'more positive' both statements <b>MUST BE</b> comparative</p>
	(c) (i)	$\text{S}_2\text{O}_8^{2-} + 2\text{I}^- \rightarrow 2\text{SO}_4^{2-} + \text{I}_2$ correct equation balanced ✓	1	<b>IGNORE</b> state symbols

Question	Answer	Marks	Guidance
	<p>(ii)</p> <ol style="list-style-type: none"> <li>1. use filter of complementary colour to iodine/solution ✓</li> <li>2. zero colorimeter with water ✓</li> <li>3. measure absorbance readings of standard solutions/solutions of known concentration (of iodine) ✓</li> <li>4. plot calibration graph ✓</li> <li>5. take absorbance readings of the reacting mixture at known/certain times AW (must refer to <b>time</b> – may state units of time) ✓</li> <li>6. convert absorbance readings to iodine concentrations using the calibration curve ✓</li> </ol> <p><b>QWC:</b> In order to gain the mark for point 3, 5 or 6, absorbance must be used <b>AND</b> spelled correctly at least once</p>	6	<p>Using starch ANYWHERE is a CON and CANNOT gain marking point 3 <b>ALLOW</b></p> <ul style="list-style-type: none"> <li>• filter/wavelength giving maximum absorption/absorbance</li> <li>• green/blue filter</li> <li>• suitable filter</li> </ul> <p>orange/yellow/brown is <b>CON</b> <b>NOT</b> 'solvent' instead of 'water'</p> <p><b>IGNORE</b> references to quenching procedures</p>

Question	Answer	Marks	Guidance
(iii)	 <p>suitable graph sketched ✓</p> <p>construction lines to determine half-lives from graph ✓</p> <p>constant half-life (means first order) ✓</p>	3	<p>graph should show a curve (<b>NOT</b> straight line) <b>decreasing</b> from left to right (ANY decreasing curve will do)</p> <p>at least <b>two</b> sets of construction lines should be shown on graph</p> <p>dependent on showing half-lives on graph (numbers on scale <i>or</i> by labelling with <math>t_{1/2}</math>)</p>
(d) (i)	homogeneous, reactants and catalyst/ $\text{Fe}^{3+}$ are in the same phase / state ✓	1	<p><b>MUST</b> mention 'reactants <b>and</b> catalyst'</p> <p><b>IGNORE</b> 'redox'</p> <p><b>NOT</b> 'substrate'</p> <p><b>ALLOW</b> 'all in solution' for 'same phase'</p>
(ii)	<p><math>\text{Fe}^{2+}</math>      <math>1s^2 2s^2 2p^6 3s^2 3p^6 3d^6</math></p> <p><math>\text{Fe}^{3+}</math>      <math>1s^2 2s^2 2p^6 3s^2 3p^6 3d^5</math></p> <p>correct number of <b>EXTRA</b> electrons added</p> <p><math>14(\text{Fe}^{2+})</math> and <math>13(\text{Fe}^{3+})</math> ✓</p> <p>both fully correct ✓</p>	2	<b>IGNORE</b> $4s^0$

Question	Answer	Marks	Guidance
	<p>(iii) 1. <math>\text{Fe}^{3+}</math> reacts with/oxidises <math>\text{I}^-</math> <b>ORA</b> ✓</p> $2\text{Fe}^{3+} + 2\text{I}^- \rightarrow 2\text{Fe}^{2+} + \text{I}_2 \quad \checkmark$ <p><i>explanation:</i> the <math>E^\ominus</math> of <math>\text{Fe}^{3+}/\text{Fe}^{2+}</math> half-cell is <b>more positive / less negative</b> than that of the <math>\text{I}^-/\text{I}_2</math> half-cell <b>ORA</b> ✓ AW</p> <p>2. <math>\text{Fe}^{2+}</math> reacts with/reduces <math>\text{S}_2\text{O}_8^{2-}</math> <b>ORA</b> ✓</p> $\text{S}_2\text{O}_8^{2-} + 2\text{Fe}^{2+} \rightarrow 2\text{SO}_4^{2-} + 2\text{Fe}^{3+} \quad \checkmark$ <p><i>explanation:</i> the <math>E^\ominus</math> of <math>\text{S}_2\text{O}_8^{2-}/\text{SO}_4^{2-}</math> half-cell is <b>more positive</b> than that of the <math>\text{Fe}^{3+}/\text{Fe}^{2+}</math> half-cell <b>ORA</b> ✓ AW</p> <p><b>QWC:</b> to gain the explanation mark for either <b>1</b> or <b>2</b>, the data has to be linked correctly to the reaction</p>	6	<p><b>IGNORE</b> references to activation enthalpy <b>OR</b> variable valency/oxidation states</p> <p><b>ALLOW</b> <math>E^\ominus</math> of <math>\text{Fe}^{3+}</math> is more positive etc</p> <p><b>ALLOW</b> <math>E^\ominus</math> of <math>\text{S}_2\text{O}_8^{2-}</math> is more positive etc</p> <p><b>NOTE</b> If answer starts with <math>\text{Fe}^{2+}</math> rather than <math>\text{Fe}^{3+}</math> lose 1<sup>st</sup> mark but ecf since not answering question so can get 5 marks</p>
(e)	(i) Rate = $k \times [\text{S}_2\text{O}_8^{2-}] \checkmark \times [\text{I}^-] \checkmark$	2	<b>IGNORE</b> state symbols
	(ii) uses one of the 3 sets of results: 1. $k = 2.0 \times 10^{-5} / (0.075 \times 0.040)$ 2. $k = 4.0 \times 10^{-5} / (0.150 \times 0.040)$ 3. $k = 1.0 \times 10^{-5} / (0.075 \times 0.020) \checkmark$  $k = \mathbf{0.0067} \checkmark$  $\text{mol}^{-1} \text{dm}^3 \text{s}^{-1} \checkmark$	3	<p>ecf from part e(i) provided working is shown <b>ALLOW</b> standard form (eg <math>6.7 \times 10^{-3}</math>) Correct numerical answer without working scores 2 for calculation part</p> <p><b>NO</b> ecf from first mark to second</p> <p><b>ALLOW</b> any number of sig figs e.g. 0.007, 0.00667 <b>NOT</b> 0.006 <b>NOR</b> 0.0066</p> <p><b>ALLOW</b> units in any order, e.g. <math>\text{dm}^3 \text{s}^{-1} \text{mol}^{-1}</math></p>
	(iii) $2.0 \times 10^{-5} \checkmark$ $\text{mol dm}^{-3} \text{s}^{-1} \checkmark$	2	<i>Mark separately</i>
<b>Total</b>		<b>31</b>	



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