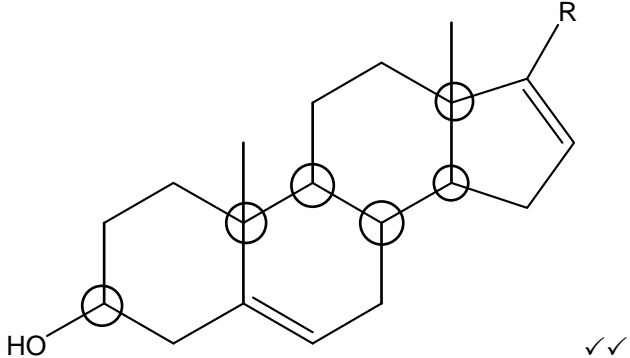
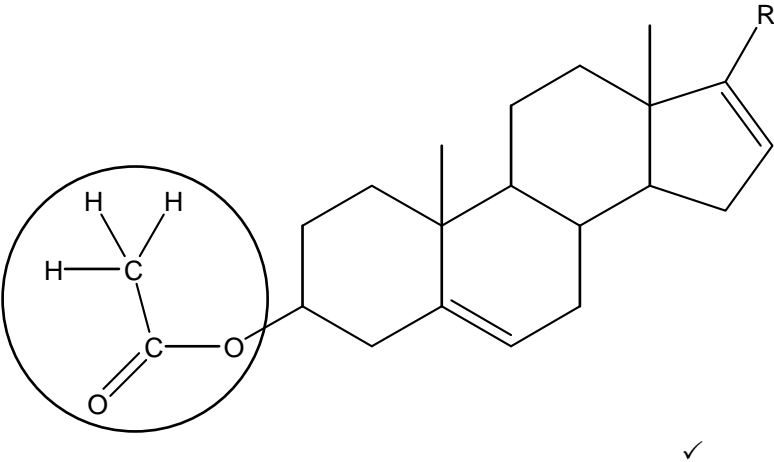


Question	Expected Answers	Marks	Additional Guidance
1 (a)	Alkene / carbon-carbon double bond / C=C ✓ alcohol / hydroxyl / hydroxy ✓	2	Double bond alone does <b>NOT</b> score. <b>ALLOW</b> secondary alcohol but not primary or tertiary. Do <b>NOT</b> allow hydroxide.
1 (b) (i)		2	<p>If more than 3 circles, any surplus <b>INCORRECT</b> ones are <b>CON IGNORE</b> surplus <b>CORRECT</b> circles</p> <p>2 marks for 3 correct 1 mark for 2 correct 0 marks for only 1 correct</p>
1 (b) (ii)	(The mirror images / molecules / structures / enantiomers / they) are non-superimposable / cannot be superimposed ✓	1	
1 (c) (i)	<p>The <b>masses</b> of the different types of atom present are <b>not integers</b> / masses are measured relative to carbon-12 (12.00000) ✓</p> <p><b>OR</b></p> <p>different compounds with the same whole number molecular mass will have <b>different <math>M_r</math> values</b> from high resolution spectra / AW ✓</p> <p><b>AND</b></p> <p>Comparison of <b><math>M_r</math></b> with database / list of formulae/<math>M_r</math> values ✓</p>	2	<p><i>mark independently</i> <b>ALLOW</b> high resolution MS gives accurate <math>M_r</math> to 4 decimal places</p> <p><b>ALLOW</b> calculate molecular formula by using masses of atoms involved</p>
1 (c) (ii)	<p>Peak: CH<sub>3</sub><sup>+</sup> ✓</p> <p>species lost = OH ✓</p>	2	<p><b>MUST</b> have correct charge for first mark</p> <p><b>MUST</b> be neutral for the second mark</p>

Question	Expected Answers	Marks	Additional Guidance
1 (d) (i)		1	<b>ALLOW</b> this, though not full structural! <b>DO NOT ALLOW</b> skeletal formula
1 (d) (ii)	ethanoic acid ✓ <u>concentrated</u> sulfuric acid / <u>concentrated</u> hydrochloric acid ✓	2	<i>Mark separately</i> <b>IGNORE</b> conc./ dil. / aq. for ethanoic acid Moderately is CON for acid <b>ALLOW</b> correct formula. e.g. CH <sub>3</sub> COOH and conc H <sub>2</sub> SO <sub>4</sub>
1 (e)	$\text{C}_{24}\text{H}_{31}\text{NO} + 2\text{Br}_2 \rightarrow \text{C}_{24}\text{H}_{31}\text{NOBr}_4$ ✓	1	<b>ALLOW</b> correct formula to be given in any order of atoms

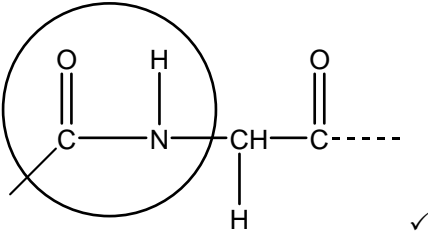
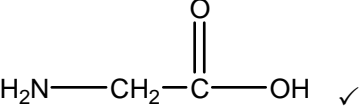
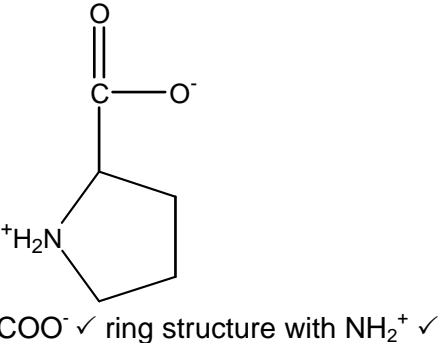
Question	Expected Answers	Marks	Additional Guidance
1 (f)	<p>Any <b>four</b> from:</p> <ol style="list-style-type: none"> <li>1. Heat the (impure) sample with solvent / use hot/warm solvent ✓</li> <li>2. with a minimum amount (of solvent) / add solvent to solid until just dissolves ✓</li> <li>3. filter ✓</li> <li>4. leave filtrate/solution/mixture to cool / leave to crystallise ✓</li> <li>5. filter off crystals, wash and dry ✓</li> </ol> <p><b>AND for QWC mark:</b></p> <p><b>EITHER</b> First filtration / filtration of hot solution removes <b><u>insoluble</u></b> impurities</p> <p><b>OR</b> after crystallisation <b><u>soluble</u></b> impurities stay in solution / AW</p> <p>(QWC) ✓</p>	5	<p><b>Please annotate with ticks to show where ALL marks are awarded</b></p> <p><b>IGNORE</b> <i>extra points after 4 have been scored</i></p> <p><b>IGNORE</b> name of solvent</p> <p>If no solvent is used then no marks can be scored similarly if <b>4.</b> is <b>incorrect</b> then <b>5.</b> cannot be scored</p> <p><b>QWC</b> only one statement required for this mark</p> <p><b>ALLOW</b> 'impurities which do <b>NOT</b> dissolve' / solid impurities</p> <p><b>ALLOW</b> 'impurities which dissolve'</p>

Question	Expected Answers	Marks	Additional Guidance
1 (g)	<p>1. Is the drug safe (to be used in humans)? ✓</p> <p>2 Does it do the job it is designed to do? ✓</p> <p>3 Is it better than the standard treatment being used? ✓</p>	3	<p><i>Alternative answers to the 3 answers on the left:</i></p> <p><b>IGNORE</b> incorrect answers, including 'is it toxic/harmful?', 'is it cost-effective?' 'can it be modified?'</p> <p><b>NOTE</b> only one mark can be scored for each of the candidate's questions (e.g. 'effective and better' in one question only scores 1)</p> <p>1. Are there any <b>side-effects</b>? / AW</p> <p>1. What is the <b>safe-dose</b>? / AW</p> <p>2. Is it <b>effective</b>? / does it work AW</p> <p>3. Is it an <b>improvement</b> on other drugs? / AW</p> <p>3. Can it be used to treat <b>other</b> symptoms/health problems / diseases? / AW <b>ALLOW</b> specific examples e.g. <i>can it be used to treat cancer?</i></p>
<b>Total</b>		<b>21</b>	

Question	Expected Answers	Marks	Additional Guidance
2 (a)	$\text{Fe} + 2\text{CH}_3\text{COOH} \rightarrow \text{Fe}(\text{CH}_3\text{COO})_2 + \text{H}_2$ ✓✓	2	Correct formula for $\text{H}_2$ or formula of salt ✓ correct formulae <b>AND</b> balanced ✓ <b>ALLOW</b> $\text{Fe}(\text{C}_2\text{H}_3\text{O}_2)_2$ or $\text{Fe}^{2+}(\text{CH}_3\text{COO}^-)_2$ <b>DO NOT ALLOW</b> $\text{Fe}(\text{II})(\text{CH}_3\text{COO})_2$
2 (b) (i)	Fe atom: .... $3d^6 4s^2$ <b>OR</b> $4s^2 3d^6$ ✓ Fe(II) ion: .... $3d^6$ Fe(III) ion: .... $3d^5$ ✓	2	<b>Fe atom</b> correct ✓ <b>BOTH</b> ions correct ✓ <b>IGNORE</b> ' $4s^0$ ' for ions
2 (b) (ii)	Half-filled d-shell / Half-filled d-orbitals (more stable) ✓ <b>OR</b> $3d^5$ configuration / $3d^5$ arrangement is (more) stable (than $3d^6$ ) ✓	1	<b>ALLOW</b> 'paired electrons are less stable than if unpaired' ORA <b>ALLOW</b> in $3d^5$ electrons are in separate orbitals/d-subshells
2 (c) (i)	$\text{Fe}^{2+}(\text{aq}) + 2\text{OH}^-(\text{aq}) \rightarrow \text{Fe}(\text{OH})_2(\text{s})$ ✓✓	2	Formulae correct and balanced ✓ <b>ALLOW</b> $\text{Fe}^{2+}(\text{OH}^-)_2$ correct state symbols ✓ <i>if first mark not gained:</i> must be $(\text{aq}) + (\text{aq}) \rightarrow (\text{s})$ or $(\text{aq}) + (\text{aq}) \rightarrow (\text{s}) + (\text{aq})$
2 (c) (ii)	Red-brown ppt is iron(III) hydroxide / (hydrated) iron(III) oxide ✓ $\text{Fe}^{2+}$ / Fe(II) ions/ $\text{Fe}(\text{OH})_2$ are oxidised / lose electrons ✓  (by) oxygen ✓ THIS IS DEPENDENT ON Fe(II) ion/compound being oxidised	3	<b>ALLOW</b> correct formula, $\text{Fe}(\text{OH})_3$ / $\text{Fe}_2\text{O}_3$ / $\text{Fe}_2\text{O}_3 \cdot x\text{H}_2\text{O}$ <b>ALLOW</b> names May be shown by an equation: e.g. $\text{Fe}^{2+} \rightarrow \text{Fe}^{3+} + \text{e}^-$ $\text{Fe}(\text{OH})_2 + \text{O}_2 \rightarrow \text{Fe}_2\text{O}_3$ <b>IGNORE</b> 'by air'

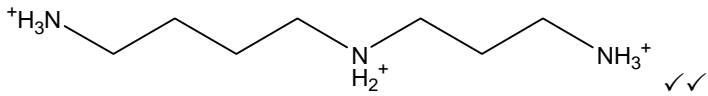
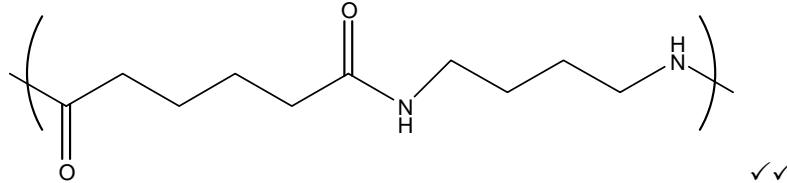
2 (d) (i)	Ce(SO <sub>4</sub> ) <sub>2</sub> ✓	1										
2 (d) (ii)	1. moles of Ce <sup>4+</sup> in titre = <b>0.100 x (18.5/1000)</b> (= 0.00185) ✓ 2. moles of Fe <sup>2+</sup> in 25.0 cm <sup>3</sup> = 0.00185 moles of Fe <sup>2+</sup> in 1000 cm <sup>3</sup> = 0.00185 x 1000 / 25.0 = <b>0.0740</b> ✓	2	The mark is for the working shown in bold  <b>ALLOW</b> answer to 2 sig figs i.e. 0.074 and ecf from 1									
2 (e) (i)	1. moles of A in 100 cm <sup>3</sup> = <b>0.1(0) / 213</b> (= 4.695 x 10 <sup>-4</sup> ) ✓ 2. moles of A in 1000 cm <sup>3</sup> = 10 x 0.10 / 213 = <b>4.7 x 10<sup>-3</sup></b> / 0.0047 ✓ 2sf	2	<i>Remember that in calculations correct answer gets full marks with or without working</i> <b>ALLOW</b> 4.70 x 10 <sup>-3</sup> (3 sf)									
2 (e) (ii)	<table><tr><th>wavenumber / cm<sup>-1</sup></th><th>bond</th><th>location</th></tr><tr><td>3150</td><td>O-H</td><td>carboxylic acid</td></tr><tr><td>1715</td><td>C=O</td><td>ketone AND/OR carboxylic acid</td></tr></table> <div>✓</div> <div>✓</div>	wavenumber / cm <sup>-1</sup>	bond	location	3150	O-H	carboxylic acid	1715	C=O	ketone AND/OR carboxylic acid	2	<b>BOTH bonds</b> correct ✓  <b>BOTH locations</b> correct ✓
wavenumber / cm <sup>-1</sup>	bond	location										
3150	O-H	carboxylic acid										
1715	C=O	ketone AND/OR carboxylic acid										
2 (e) (iii)	1. COOH / carboxylic (acid) / carboxyl group ✓ (reacts with alkali) to form:  2. ions in solution / a soluble salt / salt that dissolves / soluble carboxylate (allow formula) ✓  <b>OR</b>  2. carboxylate/anion of carboxylic acid (allow formula) forms bonds with water	2	<b>IGNORE</b> references to intermolecular bonding									

2 (f)	<p>Any <b>5</b> from the following <b>6</b> marking points but if <b>no QWC</b> <b>maximum mark is 4</b></p> <p>1. Use an appropriate/suitable filter <b>OR</b> a filter having the complementary colour (if named must be yellow/green) ✓</p> <p>2. (Put a sample of the reaction mixture into the colorimeter and) take <b>absorbance</b> readings at set (time/regular) intervals AW ✓</p> <p>3. convert <b>absorbance</b> readings to concentrations using the calibration curve ✓</p> <p>4. plot graph of concentration v time <b>OR</b> 1/time for reaction ✓</p> <p>5. determine/measure /find half-lives from graph ✓</p> <p>6. constant half-life = first order ✓</p>	5	<p><b>Please annotate with ticks to show marks awarded</b></p> <p>If complementary ignore any other colour</p> <p><b>QWC</b> 'absorbance' must be spelt correctly in either mark 2. or 3. for that mark to be allowed (this is NOT an extra mark)  <b>ALLOW</b> 'absorbency' but NOT 'absorbancy'  <b>NOTE</b> if no 'absorbance', max mark = 4</p> <p>3. <b>ALLOW</b> for only one absorbance reading</p> <p>4. <b>IGNORE</b> rate</p> <p><b>ALLOW</b> points 5. and 6. may be shown using labelled diagrams</p> <p>5. 'graph' must refer to concentration v time plot</p> <p>5. find rate of reaction by drawing tangents on graph ✓</p> <p>6. if concentration doubles and rate doubles = first order / plot of rate v [B] gives (diagonal) line (through the origin) / directly proportional ✓</p>
<b>Total</b>		<b>24</b>	

Question	Expected Answers	Marks	Additional Guidance
3 (a) (i)		1	<p><b>ALLOW</b> the linkage to proline ring (C,O and N atoms circled)</p> <p><b>DO NOT ALLOW</b> if only the bond in C-N is circled</p>
3 (a) (ii)		1	<b>ALLOW</b> H <sub>2</sub> N-CH <sub>2</sub> -COOH / full structure
3 (a) (iii)		2	<p><b>Structure must be a zwitterion to score</b></p> <p><b>ALLOW</b> COO<sup>-</sup></p> <p><b>ALLOW</b> + charge on H or N</p>
3 (b) (i)	<p>(at high temperatures / 50°C) intramolecular/hydrogen bonds break ✓</p> <p>and active site lost/altered/changed ✓</p>	2	<p><b>IGNORE</b> intermolecular / any other types of intramolecular force / changing hydrogen bonds</p> <p><b>ALLOW</b> bonds in tertiary structure</p> <p><b>IGNORE</b> references to denaturing</p> <p><b>ALLOW</b> 'active site is deformed/distorted' / no longer complementary/fits substrate</p> <p><b>ALLOW</b> 'tertiary structure' for 'active site'</p>

Question	Expected Answers	Marks	Additional Guidance
3 (b) (ii)	(change of pH) affects charges on polar/some/certain groups/ active site <b>OR</b> ionisable groups are altered ✓  prevents correct interactions/bonds between enzyme and substrate AW ✓	2	<b>ALLOW</b> a correct example (e.g. COOH, COO <sup>-</sup> , NH <sub>2</sub> , NH <sub>3</sub> <sup>+</sup> ) <b>IGNORE</b> references to denaturing  <b>ALLOW</b> ionic interactions/bonds are disrupted between enzyme and substrate <b>ALLOW</b> substrate does not fit/bind/react
3 (c) (i)	Rate = k x [P] x [enzyme] ✓  mol <sup>-1</sup> dm <sup>+3</sup> s <sup>-1</sup> ✓	2	<b>ALLOW</b> 'hydroxylase' or E or enz or complete name for enzyme <b>ALLOW</b> '(rate equation) = k x [P] x [enzyme]'; must have '='  <b>ALLOW</b> units in any order and dm <sup>3</sup> <b>ALLOW</b> '/' for '-1' e.g. dm <sup>+3</sup> / mol/ s and sec <sup>-1</sup> ecf for units
3 (c) (ii)	(When [P] is low) not all enzyme active sites are filled/saturated <b>OR</b> P can form a P-enzyme substrate (can be given as an equation) <b>OR</b> active sites available for substrates <b>OR</b> P can bind to active sites AW ✓  (as [P] increases) rate increases <b>in proportion</b> (so first order) AW ✓	2	<b>DO NOT ALLOW</b> 'there are an excess of enzymes' <b>ALLOW</b> the rate determining step is the formation of P- enzyme substrate / rds involves <u>one molecule</u> of P      <b>DO NOT ALLOW</b> 'as [P] increases rate increases' alone. There must be some indication of how the rate increases e.g. rate doubles as [P]/conc. of P/amount of P/number of molecules of P/P doubles
3 (c) (iii)	all the <u>active sites</u> are filled/saturated (any increase in [P] will not affect the reaction rate) <b>OR</b> no <u>active site</u> is available for P to bind to/react with ✓  (so) <b>order</b> becomes/is <b>zero</b> ✓	2	<b>ALLOW</b> rds involves the breakdown of the enzyme-substrate complex (which does not depend on the concentration of P)

Question	Expected Answers	Marks	Additional Guidance
3 (d)	<p>Any <b>two</b> from:</p> <p><b>speeds up</b> reaction rate</p> <p><b>reduces the number of steps</b> in a synthesis ✓</p> <p><b>improves the atom economy</b> AW ✓</p> <p><b>reduces the amount of energy/heat</b> required AW ✓</p> <p><b>easier separation</b> methods ✓</p> <p>enzymes can be <b>reused/recycled</b> ✓</p> <p>uses less toxic solvents/producing <b>less hazardous waste</b> no/fewer <b>organic solvents</b> used ✓</p> <p><b>reduces</b> use of more <b>toxic catalysts</b> ✓</p>	2	<p><b>ALLOW</b> it is a one step process</p> <p><b>ALLOW</b> lower temperature/pressure used/needed/required</p> <p><b>IGNORE</b> renewed</p>
	<b>Total</b>	<b>16</b>	

Question	Expected Answers	Marks	Additional Guidance
4 (a)	1,4-diaminobutane ✓✓	2	1,4-diamino ✓ <b>ALLOW</b> 1,4-diamine <b>DO NOT ALLOW</b> 1,4-butandiamine  butane ✓ <b>ALLOW</b> butan (often in middle of name) but <b>DO NOT ALLOW</b> buta  1,6-diaminohexane scores 1 mark  <b>IGNORE</b> gaps, commas and dashes
4 (b)	 ✓✓	2	extra H <sup>+</sup> on one amino group ✓  all correct ✓  <b>ALLOW</b> +ve charge on N or H <b>IGNORE</b> length of carbon chains / missing Hs on carbons <b>ALLOW</b> 1 mark if ALL 3 amino group Hs are correct but positive charge missing
4 (c) (i)	 ✓✓	2	amide group correct ✓ completely correct (including carbon chains) ✓  <b>ACCEPT</b> molecule the other way around. <b>ALLOW</b> structural formula or without brackets
4 (c) (ii)	(Secondary) amide ✓	1	<b>DO NOT ALLOW</b> peptide
4 (c) (iii)	Hydrogen chloride / HCl ✓  a small molecule/HCl has been eliminated/formed ✓	2	<b>DO NOT ALLOW</b> hydrochloric acid for the first mark  <b>IGNORE</b> 'water formed'

4 (d)	<p>Water (rather than HCl) is formed in the reaction ✓</p> <p>HCl is toxic/harmful/dangerous/polluting (to the environment) / HCl needs to be disposed of <b>ORA</b> ✓</p> <p><b>ALTERNATIVE ANSWER</b> <b>C</b> contains chlorine ✓ which requires extra energy/resources to make AW ✓</p>	2	<p><i>mark independently</i></p> <p><b>ALLOW</b> HCl causes acid rain / is corrosive</p> <p><i>second mark depends on first</i></p>
4 (e) (i)	<p><i>Stanyl</i>: hydrogen bond(ing) ✓</p> <p><i>poly(ethene)</i>: instantaneous (dipole)-induced dipole (bonds) ✓</p>	2	<p><b>ALLOW</b> id-id bonding / van der Waals forces</p>
4 (e) (ii)	<p>intermolecular bonds in polythene are <b>weaker</b> than those in <i>Stanyl</i> ✓ <b>ORA</b></p> <p><b>less energy</b> / lower temperatures needed to break the imb in poly(ethene) / separate chains / enable chains to slide <b>ORA</b> ✓</p> <p><b>chains can move / slide over</b> each other (and polymer softens) ✓</p>	3	<p><b>Please annotate with ticks to show where ALL marks are awarded</b></p> <p><b>ALLOW</b> intermolecular forces <b>ALLOW</b> for both marks named imb from 4 (e) (i), provided <i>Stanyl</i> bonds are stronger</p> <p><b>ALLOW</b> less heat</p>
<b>Total</b>		<b>16</b>	

Question	Expected Answers	Marks	Additional Guidance									
5 (a) (i)	<p>Since <math>E^\ominus</math> for Cu(/Cu<sup>2+</sup>) is more negative than Ag(/Ag<sup>+</sup>) <b>ORA</b> ✓</p> <p>electrons will flow/move (from Cu) to Ag / from Cu(/Cu<sup>2+</sup> to Ag/Ag<sup>+</sup> ✓</p>	2	<p><i>mark independently</i></p> <p><b>ALLOW</b> smaller/larger since both <math>E^\ominus</math> are positive <b>IGNORE</b> lower/higher <b>OR</b> references to redox processes <b>ALLOW</b> '<math>E^\ominus_{\text{cell}}</math>' for '<math>E^\ominus</math>', <math>E^\ominus</math> for Cu(/Cu<sup>2+</sup>) is less positive than Ag(/Ag<sup>+</sup>) <b>ALLOW</b> 'from copper to silver' or 'Cu to Ag<sup>+</sup>' <b>IGNORE</b> 'electrons flow through water' / attract / gain <b>DO NOT ALLOW</b> 'from Cu<sup>2+</sup>'</p>									
5 (a) (ii)	0.46 V ✓	1	<b>ALLOW</b> + or – 0.46									
5 (a) (iii)	Cu + 2Ag <sup>+</sup> → Cu <sup>2+</sup> + 2Ag ✓	1	<b>IGNORE</b> state symbols									
5 (b) (i)	<p>Oxidising agent = H<sup>+</sup> /H<sub>3</sub>O<sup>+</sup> ✓</p> <p><math>E^\ominus</math> values are measured with respect/compared to the (standard) H<sub>2</sub>/H<sup>+</sup> half-cell AW / <math>E^\ominus</math> H<sub>2</sub>/H<sup>+</sup> = 0 ✓</p> <p>metals with a negative electrode potential value will be oxidised by / will react with H<sup>+</sup> ions/HCl AW <b>ORA</b> ✓</p>	3	<p><b>ALLOW</b> hydrogen ions</p> <p><b>ALLOW one mark</b> for saying acids/H<sup>+</sup> can oxidise Zn but not Cu</p>									
5 (b) (ii)	<p>1. Moles of Cu<sup>2+</sup> in 250 cm<sup>3</sup> = <b>0.150 x (250/1000)</b> ✓</p> <p>2. Mass of copper in sample = 0.0375 x 63.5 = <b>2.381</b> ✓</p> <p>3. % of Cu in brass = 2.381 / 3.97 x 100 = <b>60</b> ✓</p>	3	<p>1. The mark is for the working shown in bold</p> <p><b>ALLOW</b> any number of sig. figs. <b>ALLOW</b> ecf from 1. and 2. <b>DO NOT ALLOW</b> 59</p>									
5 (b) (iii)	<table><tr><td></td><td>copper(II) complex formed with EDTA<sup>4-</sup></td><td rowspan="4">✓ ✓ ✓</td></tr><tr><td>formula</td><td>[Cu(EDTA)]<sup>2-</sup></td></tr><tr><td>shape</td><td>octahedral</td></tr><tr><td>coordination number</td><td>6</td></tr></table>		copper(II) complex formed with EDTA <sup>4-</sup>	✓ ✓ ✓	formula	[Cu(EDTA)] <sup>2-</sup>	shape	octahedral	coordination number	6	3	<p><i>Mark separately</i></p> <p>1 mark for each correct answer</p> <p><b>ALLOW</b> without square brackets</p>
	copper(II) complex formed with EDTA <sup>4-</sup>	✓ ✓ ✓										
formula	[Cu(EDTA)] <sup>2-</sup>											
shape	octahedral											
coordination number	6											
<b>Total</b>		<b>13</b>										