



Mark Scheme (Results)

January 2018

Pearson Edexcel International Advanced
Level In Chemistry (WCH05) Paper 01
General Principles of Chemistry II –
Transition Metals and Organic Nitrogen
Chemistry

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General Marking Guidance

- All candidates must receive the same treatment. Examiners must mark the first candidate in exactly the same way as they mark the last.
- Mark schemes should be applied positively. Candidates must be rewarded for what they have shown they can do rather than penalised for omissions.
- Examiners should mark according to the mark scheme not according to their perception of where the grade boundaries may lie.
- There is no ceiling on achievement. All marks on the mark scheme should be used appropriately.
- All the marks on the mark scheme are designed to be awarded. Examiners should always award full marks if deserved, i.e. if the answer matches the mark scheme. Examiners should also be prepared to award zero marks if the candidate's response is not worthy of credit according to the mark scheme.
- Where some judgement is required, mark schemes will provide the principles by which marks will be awarded and exemplification may be limited.
- When examiners are in doubt regarding the application of the mark scheme to a candidate's response, the team leader must be consulted.
- Crossed out work should be marked UNLESS the candidate has replaced it with an alternative response.

Section A (multiple choice)

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 1 | <p>The only correct answer is D</p> <p><i>A is not correct because the electrode is wrong and Mn^{2+} ions are missing from the solution</i></p> <p><i>B is not correct because the electrode is wrong</i></p> <p><i>C is not correct because Mn^{2+} ions are missing from the solution</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 2 | <p>The only correct answer is C</p> <p><i>A is not correct because the oxidation number of Cr is + 3 and Mn is + 2</i></p> <p><i>B is not correct because the oxidation number of Cr is + 6 and Ti is + 3</i></p> <p><i>D is not correct because the oxidation number of Cr is + 6 and Mn is + 7</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 3 | <p>The only correct answer is A</p> <p><i>B is not correct because the oxidation number should be $+3 - 2 = +1$</i></p> <p><i>C is not correct because the oxidation number should be $+3 - 2 = +1$</i></p> <p><i>D is not correct because the oxidation number should be $+3 - 2 = +1$</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 4(a) | <p>The only correct answer is A</p> <p><i>B is not correct because the complex is linear</i></p> <p><i>C is not correct because the complex is square planar</i></p> <p><i>D is not correct because the complex is octahedral</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 4(b) | <p>The only correct answer is B</p> <p>A is not correct because the oxidation number of the metal is +3</p> <p>C is not correct because the oxidation number of the metal is +2</p> <p>D is not correct because the oxidation number of the metal is +4</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 5(a) | <p>The only correct answer is C</p> <p>A is not correct because it is not $-4/2$</p> <p>B is not correct because it is not $-4/2$</p> <p>D is not correct because it is not $-4/2$</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 5(b) | <p>The only correct answer is C</p> <p>A is not correct because incorrect number of Z ligands</p> <p>B is not correct because incorrect number of Z ligands and incorrect charge</p> <p>D is not correct because incorrect charge</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 6(a) | <p>The only correct answer is A</p> <p>B is not correct because although copper(I) oxide is reddish brown, it is an incorrect product</p> <p>C is not correct because copper(II) oxide is black not brown and it is an incorrect product</p> <p>D is not correct because zinc sulfate is white not brown</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 6(b) | <p>The only correct answer is B</p> <p><i>A is not correct because it is not a white solid</i></p> <p><i>C is not correct because it is not a white solid</i></p> <p><i>D is not correct because it is soluble</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 7 | <p>The only correct answer is B</p> <p><i>A is not correct because it is not an addition reaction</i></p> <p><i>C is not correct because not it is not a nucleophilic or addition reaction</i></p> <p><i>D is not correct because it is not a nucleophilic reaction</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 8(a) | <p>The only correct answer is D</p> <p><i>A is not correct because this is the wrong product</i></p> <p><i>B is not correct because this is the wrong product</i></p> <p><i>C is not correct because this is the wrong product</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 8(b) | <p>The only correct answer is C</p> <p><i>A is not correct because it is electrophilic not nucleophilic</i></p> <p><i>B is not correct because electromeric effect outweighs inductive effect</i></p> <p><i>D is not correct because it is a nucleophile not an electrophile</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 9(a) | <p>The only correct answer is A</p> <p><i>B is not correct because the locants are incorrect</i></p> <p><i>C is not correct because there are no locants for the amine groups</i></p> <p><i>D is not correct because there are no locants for the amine groups</i></p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 9(b) | <p>The only correct answer is D</p> <p>A is not correct because there is no carbonyl group</p> <p>B is not correct because there is no carbonyl group</p> <p>C is not correct because the amide groups are incorrect</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 10(a) | <p>The only correct answer is C</p> <p>A is not correct because it is not an acid</p> <p>B is not correct because it is not an acid</p> <p>D is not correct because it is optically active</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 10(b) | <p>The only correct answer is A</p> <p>B is not correct because Q forms phenol</p> <p>C is not correct because R does not form sodium benzoate</p> <p>D is not correct because S does not form sodium benzoate</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 11(a) | <p>The only correct answer is B</p> <p>A is not correct because it is a primary alcohol and should be a secondary alcohol</p> <p>C is not correct because it is a primary alcohol and should be a secondary alcohol</p> <p>D is not correct because it is a tertiary alcohol and should be a secondary alcohol</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|---|------------|
| 11(b) | <p>The only correct answer is B</p> <p>A is not correct because there would be no loss of carbon atom</p> <p>C is not correct because there would be no loss of carbon atom</p> <p>D is not correct because there would be no loss of carbon atom</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 11(c) | <p>The only correct answer is D</p> <p>A is not correct because this does not react with propanoic acid to give propanoyl chloride</p> <p>B is not correct because this does not react with propanoic acid to give propanoyl chloride</p> <p>C is not correct because this does not react with propanoic acid to give propanoyl chloride</p> | (1) |

| Question Number | Correct Answer | Mark |
|-----------------|--|------------|
| 11(d) | <p>The only correct answer is B</p> <p>A is not correct because it is not branched</p> <p>C is not correct because it is not an amine</p> <p>D is not correct because it is not an amine</p> | (1) |

Section B

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--------|------------|
| 12(a)(i) | $(\text{Al}^{3+} 1s^2)2s^22p^6$ (1) $(\text{Fe}^{3+} 1s^2)2s^22p^63s^23p^63d^5$ (1) ALLOW $2p_x^22p_y^22p_z^2$ / $3p_x^23p_y^23p_z^2$ ALLOW $4s^0$ included in Fe^{3+} / any other orbitals with 0 electrons IGNORE $1s^2$ written again | | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|-------------------------------------|------------|
| 12(a)(ii) | The increase in ionisation energy is balanced by an increase in hydration/lattice enthalpy ALLOW There is (only) a gradual/ small increase in (successive) ionisation energies for iron OR Iron has (several removable) d electrons of similar energies OR The 4s / 3d electrons / orbitals have similar energies (for iron) OR The energy difference / gap between 4s and 3d is small (for iron) OR The ionisation energies are similar (for iron) IGNORE References to stability of half-full <i>d</i> -subshell / References to 3p electrons | Just 'iron is a transition element' | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|---|------------|
| *12(b) | <p>First mark - splitting (3)d orbitals / (3)d subshell are/is split (in energy by the ligands)</p> <p>ALLOW "d orbital splitting" (1)</p> <p>IGNORE Just 'there is an energy difference between the (3) d orbitals'</p> <p>Second mark - absorption Electrons / photons absorb energy</p> <p>ALLOW Electrons absorb (visible) light Frequencies / wavelengths (of visible light) are absorbed (1)</p> <p>Third mark - promotion Electrons are promoted (from lower to higher energy d orbital(s) / levels) OR Electrons move (from lower) to higher energy (d) orbital(s) / levels</p> <p>ALLOW d-d transitions occur / Electrons are excited to higher energy (d) orbital(s) / levels (1)</p> <p>Fourth mark - colour Reflected / transmitted / remaining light is coloured / is in the visible region</p> <p>ALLOW Complementary colour seen (The frequency of) reflected / transmitted / remaining light is seen (1)</p> <p>IGNORE Reference to electrons relaxing / dropping to the ground state / any reference to aluminium</p> | <p>d orbital 'singular' is split</p> <p>d shell is split</p> <p>Penalise omission of (3)d only in First mark</p> <p>Light emitted</p> | (4) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|-----------------------------|------------|
| 12(c)(i) | <div data-bbox="327 264 646 488" data-label="Chemical-Block"> </div> <p>3D octahedral shape Recognisable 3D octahedral shape</p> <p>ALLOW Wedges/dots instead of dashed lines going into the page or other recognisable representations (1)</p> <p>Note The word 'octahedral' does not rescue a poor diagram</p> <p>Oxygen atoms 6 oxygen atoms joined to Fe (with or without lone pairs)</p> <p>ALLOW O₂H Oxygens to Fe joined by single bonds / arrows for this mark (1)</p> <p>IGNORE Omission of brackets and charge Incorrect charge Name of shape, even if incorrect</p> | <p>Negative charge on O</p> | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|---|------------|
| 12(c)(ii) | <p>Any 2 points from 3</p> <p>First point Fe^{3+} is (small and) highly charged/ Fe^{3+} has a high charge density (1)</p> <p>Second point Fe^{3+} polarises / weakens / distorts the O–H bond OR Fe^{3+} attracts the electrons / electron cloud / electron density in the OH bond</p> <p>ALLOW Fe^{3+} polarises / distorts the water molecule (1)</p> <p>Third point (Solvent) water acts as a base OR (Solvent) water removes/ accepts a proton OR Water ligands donate a proton to (solvent) water</p> <p>ALLOW $[\text{Fe}(\text{H}_2\text{O})_6]^{3+}$ / the complex (ion) / water ligand acts as an acid (1)</p> <p>IGNORE Just 'deprotonation' / just 'acid-base reaction' / alkali instead of base/ the complex is a proton donor</p> | <p>Water ligands act as a base</p> <p>Disproportionation</p> | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---|------------|
| 12(d) | <p>First mark – electron deficient (The aluminium atom in) AlCl_3 has an empty (p) orbital (in the outer shell) / is electron deficient / has 6 electrons (in the outer shell)</p> <p>ALLOW 6 electrons shown around Al on a diagram (1)</p> <p>IGNORE Just 'AlCl_3 has an incomplete (p) subshell'</p> <p>Second mark - bond It can accept a pair of / two electrons (from chlorine) OR Chlorine donates a pair of / 2 electrons (to AlCl_3) OR It form a dative (covalent) bond (with chlorine)</p> <p>ALLOW Mention of chloride ion (1)</p> <p>IGNORE diagram</p> | <p>Reference to 3d / 3s orbitals</p> <p>Reference to aluminium ion</p> <p>Reference to 'an electron'</p> <p>Chlorine molecule / Cl_2</p> | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--|------------|
| 12(e)(i) | <p>Ligand exchange (reaction) OR Ligand substitution (reaction) OR Ligand replacement (reaction)</p> <p>Both words needed for the mark</p> | <p>Ligand change Ligand reaction Oxidation Reduction Redox</p> | (1) |

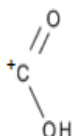
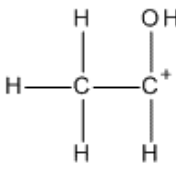
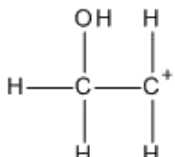
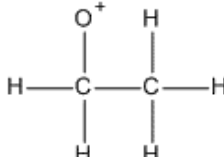
| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--------|------------|
| 12(e)(ii) | <p>Dot-and-cross diagram</p> <p>OR</p> <p>ALLOW Any symbols for electrons Bonds shown with 2 electrons on each e.g. \times (1)</p> <p>IGNORE Missing brackets and charge</p> <p>Structure and charge</p> <p>$^-S-C\equiv N$</p> <p>OR</p> <p>$S=C=N^-$</p> <p>Bonds correct and negative charge shown or stated on correct atom</p> <p>ALLOW This mark if bonds also shown in dot-and-cross diagram and negative charge on correct atom (1)</p> | | (2) |

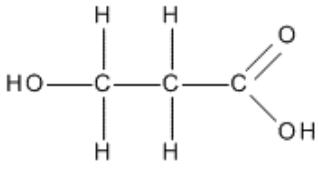
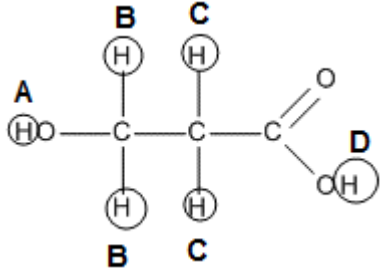
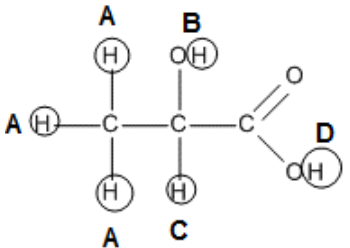
| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|--|--------|------------|
| 12(e)(iii) | <p>Dative covalent bond / bond from lone pair (of electrons) on sulfur/S and from nitrogen/N (to an empty orbital in Fe^{3+})</p> <p>IGNORE Ionic bond</p> | | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|------------|
| 12(f) | <p>Ionic equation with hydrochloric acid $\text{Al(OH)}_3 + 3\text{H}^+ \rightarrow \text{Al}^{3+} + 3\text{H}_2\text{O}$</p> <p>ALLOW $\text{Al(OH)}_3(\text{H}_2\text{O})_3 + 3\text{H}^+ \rightarrow \text{Al}^{3+} + 6\text{H}_2\text{O}$ $\text{Al(OH)}_3(\text{H}_2\text{O})_3 + 3\text{H}^+ \rightarrow [\text{Al}(\text{H}_2\text{O})_6]^{3+}$ $\text{Al(OH)}_3 + 3\text{H}^+ + 3\text{H}_2\text{O} \rightarrow [\text{Al}(\text{H}_2\text{O})_6]^{3+}$ (1)</p> <p>IGNORE $\text{H}^+ + \text{OH}^- \rightarrow \text{H}_2\text{O}$</p> <p>Ionic equation with sodium hydroxide $\text{Al(OH)}_3 + \text{OH}^- \rightarrow \text{AlO}_2^- + 2\text{H}_2\text{O}$ OR $\text{Al(OH)}_3 + \text{OH}^- \rightarrow [\text{Al(OH)}_4]^-$</p> <p>ALLOW $\text{Al(OH)}_3(\text{H}_2\text{O})_3 + \text{OH}^- \rightarrow [\text{Al(OH)}_4]^- + 3\text{H}_2\text{O}$ $\text{Al(OH)}_3(\text{H}_2\text{O})_3 + \text{OH}^- \rightarrow [\text{Al(OH)}_4(\text{H}_2\text{O})_2]^- + \text{H}_2\text{O}$ $\text{Al(OH)}_3 + 2\text{OH}^- \rightarrow [\text{Al(OH)}_5]^{2-}$ $\text{Al(OH)}_3(\text{H}_2\text{O})_3 + 2\text{OH}^- \rightarrow [\text{Al(OH)}_5]^{2-} + 3\text{H}_2\text{O}$ $\text{Al(OH)}_3 + 3\text{OH}^- \rightarrow [\text{Al(OH)}_6]^{3-}$ $\text{Al(OH)}_3(\text{H}_2\text{O})_3 + 3\text{OH}^- \rightarrow [\text{Al(OH)}_6]^{3-} + 3\text{H}_2\text{O}$ (1)</p> <p>If no other mark awarded, ALLOW 1 mark for two non-ionic / partially ionic equations e.g. $\text{Al(OH)}_3 + 3\text{HCl} \rightarrow \text{AlCl}_3 + 3\text{H}_2\text{O}$ and $\text{Al(OH)}_3 + \text{NaOH} \rightarrow \text{NaAl(OH)}_4$</p> <p>IGNORE State symbols, even if incorrect / missing square brackets</p> | | (2) |

(Total for Question 12 = 19 marks)

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--------|------------|
| 13(a)(i)* | <p>The methods below illustrate the allocation of marks. However, the marks may be scored by any correct method</p> <p>Correct molecular formulae with some working involving C/CO₂, H/H₂O and either O or use of 90 and subtraction of A_rs of C and H / COOH scores full marks</p> <p>Correct molecular formula with no working scores (1)</p> <p>Method 1</p> <p>mol CO₂ = 3.30/44 = 0.075 (= mol C) (1)</p> <p>mol H₂O = 1.35/18 = 0.075 (1)</p> <p>and</p> <p>mol H = 2 x mol H₂O = 0.150 or ratio C : H = 1 : 2 (1)</p> <p>mass O = 2.25 – ((12 x 0.075) + (1 x 0.150)) = 1.2 (g) (1)</p> <p>mol O = 1.2/16 = 0.075 (1)</p> <p>Method 2</p> <p>mass C = 3.30 x 12/44 = 0.90 (g) and mol C = 0.90/12 = 0.075 (1)</p> <p>mass H = 1.35 x 2/18 = 0.15 (g) and mol H = 0.15/1 = 0.15 (1)</p> <p>mass O = 2.25 – (0.90 + 0.15) = 1.2 (g) (1)</p> <p>mol O = 1.2/16 = 0.075 (1)</p> <p>Empirical and molecular formulae from Methods 1 and 2</p> <p>Empirical formula = CH₂O (1)</p> <p>Relative empirical formula mass CH₂O = 12 + (2 x 1) + 16 = 30</p> <p>So, molecular formula is C₃H₆O₃ (1)</p> <p>TE on incorrect moles but the ratio must be a whole number</p> <p>Method 3</p> <p>mol T = 2.25/90 = 0.025 (1)</p> <p>mol CO₂ = 3.30/44 = 0.075 (= mol C) (1)</p> <p>mol ratio T : C / CO₂ = 1 : 3 (1)</p> <p>mol H₂O = 1.35/18 = 0.075</p> <p>and</p> <p>mol H = 2 x mol H₂O = 0.150 or mol ratio C : H = 1 : 2 / 3 : 6 (1)</p> <p>90 - (36 + 6) = 48 and mol O = 48/16 = 3 (1)</p> <p>So, molecular formula is C₃H₆O₃ (1)</p> | | (6) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|---|------------|
| 13(a)(ii) | <p>Any two of (1 mark for each structure)</p> <div style="display: flex; justify-content: space-around; align-items: center;">   </div> <div style="display: flex; justify-content: space-around; align-items: center;">   </div> <p>ALLOW Structural formula i.e COOH^+, CH_3CHOH^+, $\text{CH}_2\text{OHCH}_2^+$, $\text{CH}_3\text{CH}_2\text{O}^+$</p> <p>ALLOW Structures in brackets with positive charge outside bracket</p> <p>IGNORE Position of positive charge $\text{C}_2\text{H}_5\text{O}^+$ Connectivity of the OH group</p> | <p>Extra bonds once only e.g. $-\text{COOH}^+$</p> <p>Negative charge or omission of charge once only</p> <p>HCO_2^+</p> | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|--|--------|----------|
| 13(a)(iii) | <p>Structure of compound T:</p>  <p style="text-align: right;">(1)</p> <p>IGNORE Connectivity of the OH group</p> <p>Explanation: Use of peak ratio Use of peak ratio e.g. protons A and D are ratio 1 : 1 (as they are single protons) and protons B and C are ratio 2 : 2 (as there are two protons in each environment)</p> <p>ALLOW Ratio of protons /hydrogens is 1 : 2 : 2 : 1 (1)</p> <p>Proton environments identified 4 proton environments clearly identified by symbols or words e.g.</p>  <p style="text-align: right;">(1)</p> <p>Note If compound T is identified as lactic acid, (1) mark awarded for identification of four proton environments only e.g.</p>  <p>No TE on any other structure</p> | | 3 |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|--|------------|
| 13(b)(i)* | <p>Reaction 1 Reagents: CH_3COCl / ethanoyl chloride and AlCl_3 / aluminium chloride / FeCl_3 / iron(III) chloride (1)</p> <p>Intermediate: stand alone or TE on acyl chloride used</p> <div data-bbox="341 465 657 613" data-label="Chemical-Block"> </div> <p>(1)</p> <p>Reaction 2 Reagents: conditional on a carbonyl compound HCN / hydrogen cyanide and KCN / potassium cyanide OR KCN / potassium cyanide and acid / H_2SO_4 / sulfuric acid / H^+ ions / hydrogen ions OR KCN / potassium cyanide and pH 8-10 / alkali</p> <p>ALLOW HCN / hydrogen cyanide and alkali / NaOH / sodium hydroxide / OH^- / hydroxide ions</p> <p>ALLOW Sodium for potassium and vice versa (1)</p> <p>Intermediate: stand alone</p> <div data-bbox="341 1352 657 1509" data-label="Chemical-Block"> </div> <p>(1)</p> <p>Reaction 3 Reagents: conditional on reaction with a CN group H^+ / hydrogen ions / (dilute) acid / name or formula of a strong acid</p> <p>ALLOW OH^- / hydroxide ions / alkali / name or formula of an alkali and followed by / then acidification / H^+ / hydrogen ions / (dilute) acid / name or formula of a strong acid (1)</p> <p>IGNORE Concentration of acid or alkali</p> | <p>Just 'CN^- in alkali'</p> <p>Additional reagent(s) e.g. KMnO_4</p> <p>Acid and alkali added together</p> | (5) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|---|------------|
| 13(b)(ii) | <div data-bbox="331 264 758 537"> <p>Diagram showing a segment of a polymer chain: $\text{---O---C(CH}_3\text{)(C}_6\text{H}_5\text{)---C(=O)---O---C(CH}_3\text{)(C}_6\text{H}_5\text{)---C(=O)---}$. The first ester linkage, ---C(=O)---O---, is circled.</p> </div> <p>OR</p> <div data-bbox="331 604 774 878"> <p>Diagram showing a segment of a polymer chain: $\text{---C(CH}_3\text{)(C}_6\text{H}_5\text{)---C(=O)---O---C(CH}_3\text{)(C}_6\text{H}_5\text{)---C(=O)---O---}$. The second ester linkage, ---C(=O)---O---, is circled.</p> </div> <p>First mark One correct ester linkage (as circled above) (1)</p> <p>Second mark Conditional on one (or more) ester linkage Rest of structure correct with 2 repeat units and extension bonds</p> <p>ALLOW C_6H_5 for benzene ring (1)</p> <p>IGNORE Brackets and n / bond lengths and bond angles</p> | <p>Os at both ends or no O at either end loses second mark only</p> | (2) |

(Total for Question 13 = 18 marks)

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--------|------------|
| 14(a)(i) | $\text{V}^{2+} + 2\text{H}_2\text{O} \rightarrow \text{VO}_2^+ + 4\text{H}^+ + 3\text{e}^{(-)}$ <p>ALLOW Multiples Reversible arrow, provided equation written in the direction shown $\text{V}^{2+} + 2\text{H}_2\text{O} - 3\text{e}^{(-)} \rightarrow \text{VO}_2^+ + 4\text{H}^+$</p> <p>IGNORE State symbols even if incorrect</p> | | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--|------------|
| 14(a)(ii) | <p>First mark – reducing agent Reducing agent: $\text{Fe}^{2+}(\text{aq})$ / iron(II) (ions) This can be shown in an equation (1)</p> <p>Justification: Second mark – V(V) to V(IV) $\text{Fe}^{2+}/\text{Fe}^{3+}$ electrode potential / SEP / E^\ominus value is less positive / lower than the $\text{VO}_2^+/\text{VO}^{2+}$ value / $(+)0.77 < (+)1.00$ (so V(V) is reduced to V(IV))</p> <p>OR $\text{VO}_2^+/\text{VO}^{2+}$ electrode potential / SEP / E^\ominus value is more positive / greater / higher than the $\text{Fe}^{2+}/\text{Fe}^{3+}$ value / $(+)1.00 > (+)0.77$ (so V(V) is reduced to V(IV))</p> <p>OR E^\ominus_{cell} for the reaction between VO_2^+ and Fe^{2+} is positive / $(+)0.23 \text{ V}$ / >0 (so V(V) is reduced to V(IV))</p> <p>ALLOW Any of the above 3 explanations if SO_2, Zn, V^{3+} or V^{2+} chosen as reducing agent e.g. E^\ominus for $\text{SO}_2/\text{SO}_4^{2-} = (+)0.83$ or E^\ominus for $\text{Zn}/\text{Zn}^{2+} = (+)1.76$ or E^\ominus for $\text{V}^{3+}/\text{VO}^{2+} = (+)0.66$ or E^\ominus for $\text{V}^{2+}/\text{V}^{3+} = (+)1.26$ (so V(V) is reduced to V(IV)) (1)</p> <p>Third mark – V(IV) to V(III) $\text{Fe}^{2+}/\text{Fe}^{3+}$ electrode potential / SEP / E^\ominus value is more positive / greater / higher than the $\text{VO}^{2+}/\text{V}^{3+}$ value / $(+)0.77 > (+)0.34$ (so V(IV) is not reduced to V(III))</p> <p>OR $\text{VO}^{2+}/\text{V}^{3+}$ electrode potential / SEP / E^\ominus value is less positive / lower than the $\text{Fe}^{2+}/\text{Fe}^{3+}$ value / $(+)0.34 < (+)0.77$ (so V(IV) is not reduced to V(III))</p> <p>OR E^\ominus_{cell} for the reaction between VO^{2+} and Fe^{2+} is negative / -0.43 V / <0 (so V(IV) is not reduced to V(III)) (1)</p> | <p>Incorrect value</p> <p>Incorrect value</p> <p>Incorrect value</p> | (3) |

| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|--|-------------------------------------|------------|
| 14(a)(iii) | <p>First mark - equation $2V^{3+} + H_2O \rightarrow V^{2+} + VO^{2+} + 2H^+$</p> <p>ALLOW Multiples Reversible arrow, provided reaction is written in the direction shown (1)</p> <p>IGNORE State symbols even if incorrect Cancelled / crossed out electrons</p> <p>Second mark - E^\ominus_{cell} value $E^\ominus_{\text{cell}} = -0.26 - 0.34 = -0.6(0) \text{ (V)}$</p> <p>NO TE on equation written in reverse (1)</p> <p>Third mark - feasibility E^\ominus_{cell} is negative / <0 and so the disproportionation is not feasible / V^{2+} and VO^{2+} will react to form V^{3+}</p> <p>ALLOW this mark even if an incorrect negative value is calculated for E^\ominus_{cell}</p> <p>TE on a positive value for E^\ominus_{cell} e.g. E^\ominus_{cell} is positive / $>$ and so the disproportionation is feasible (1)</p> | Equation with uncancelled electrons | (3) |

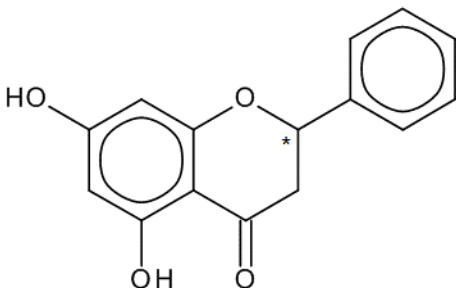
| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|------------|
| 14(b)(i) | <p>Correct answer, with or without working scores both marks</p> <p>First mark – mol I₂ Mol S₂O₃²⁻ used = $24.20 \times 0.100/1000$ = $0.00242 / 2.42 \times 10^{-3}$</p> <p>and Mol I₂ = $0.00242/2 = 0.00121 / 1.21 \times 10^{-3}$ (1)</p> <p>Second mark – conc Br₂ (Mol Br₂ = mol I₂ = 0.00121) Conc Br₂ = $0.00121 \times 1000/25.0 = 0.0484$ (mol dm⁻³) (1)</p> <p>TE on mol S₂O₃²⁻ and mol I₂</p> <p>IGNORE SF except 1SF</p> | | (2) |

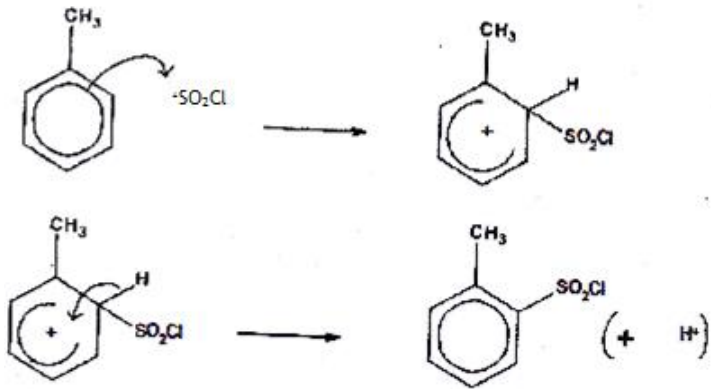
| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|--------|------------|
| 14(b)(ii) | <p>Allow correct expressions if intermediate values not evaluated</p> <p>First mark – original mol Br₂ Original mol Br₂ = $100.0 \times 0.0484/1000$ = $0.00484 / 4.84 \times 10^{-3}$ (1) TE on conc Br₂ in (i)</p> <p>Second mark – mol I₂ Mol S₂O₃²⁻ used = $16.80 \times 0.100/1000$ = $0.00168 / 1.68 \times 10^{-3}$</p> <p>and Mol I₂ = $0.00168/2 = 0.00084 / 8.4 \times 10^{-4}$ (1)</p> <p>Third mark – mol Br₂ reacted (mol Br₂ in excess = mol I₂ = 0.00084) Mol Br₂ reacted with S₂O₃²⁻ = $0.00484 - 0.00084$ = $0.00400 / 4.00 \times 10^{-3}$ (1) TE on original mol Br₂ and mol I₂/ Br₂ in excess</p> <p>Fourth mark – mol ratio Mole ratio S₂O₃²⁻ : Br₂ = $0.00100 : 0.00400$ = $1 : 4$ (1) TE on mol S₂O₃²⁻ and mol Br₂ reacted</p> <p>Fifth mark – equation – stand alone $S_2O_3^{2-} + 4Br_2 + 5H_2O \rightarrow 2SO_4^{2-} + 10H^+ + 8Br^-$</p> <p>ALLOW 8HBr + 2Br⁻ on RHS (1) No TE on incorrect mol ratio</p> <p>IGNORE State symbols even if incorrect</p> | | (5) |

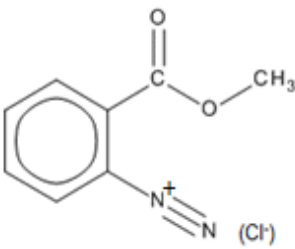
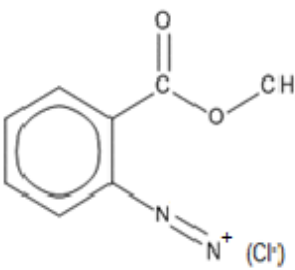
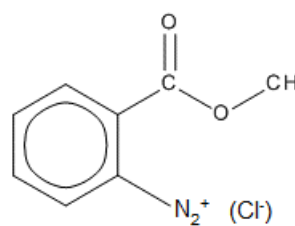
(Total for Question 14 = 14 marks)

Section C

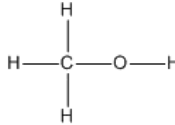
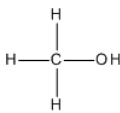
| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--|------------|
| 15(a)(i) | $C_{15}H_{12}O_4$ ALLOW symbols in any order, i.e. $C_{15}O_4H_{12}$, $H_{12}O_4C_{15}$, $H_{12}C_{15}O_4$, $O_4H_{12}C_{15}$, $O_4C_{15}H_{12}$ IGNORE Any other formulae as working | Numbers written as superscripts e.g. $C^{15}H^{12}O^4$ | (1) |

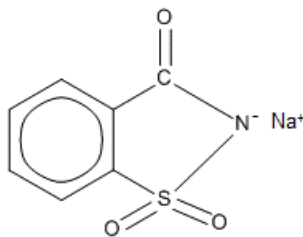
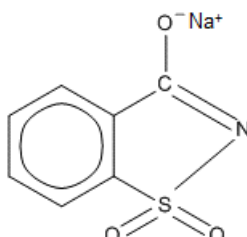
| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|---------------------------------------|------------|
| 15(a)(ii) |  ALLOW Any way of identifying the chiral carbon, including a circle, provided that it does not include any other carbon atoms | Any additional carbon atoms indicated | (1) |

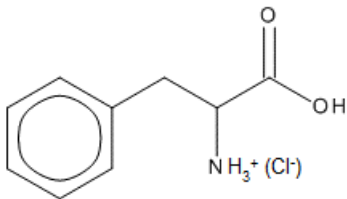
| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--|------|
| 15(b) |  <p>First mark – first curly arrow Curly arrow from on or within the circle to the S of $^+\text{SO}_2\text{Cl}$</p> <p>ALLOW Curly arrow from anywhere within the hexagon</p> <p>ALLOW Curly arrow to any part of the $^+\text{SO}_2\text{Cl}$ ion, including the + charge (1)</p> <p>Second mark - intermediate Intermediate structure including charge with horseshoe covering at least 3 carbon atoms and facing the tetrahedral carbon atom and some part of the positive charge must be within the horseshoe (1)</p> <p>Note Do not award this mark If benzene used instead of methylbenzene or if final product is not the 2-isomer</p> <p>Third mark – second curly arrow Curly arrow from C-H bond to anywhere in the hexagon, reforming the delocalised structure (1)</p> <p>IGNORE Missing H^+ Involvement of any other ion / molecule in removal of H^+</p> <p>Correct Kekule structures score full marks</p> | <p>Curly arrow on or outside the hexagon / incorrect electrophile / missing +</p> <p>Dotted bonds to H and SO_2Cl unless part of a 3D structure</p> | (3) |

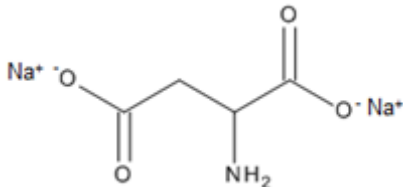
| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--|------------|
| 15(c)(i) |  <p>OR</p>  <p>OR</p>  <p>IGNORE Missing Cl⁻ on the structures shown above</p> | <p>+ on wrong nitrogen atom</p> <p>Covalent bond to Cl</p> | (1) |

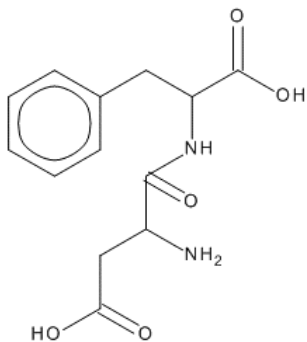
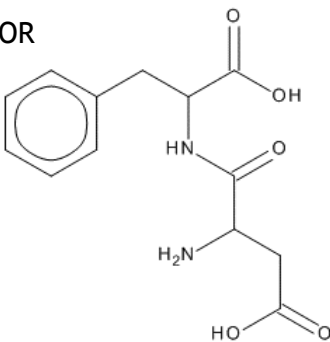
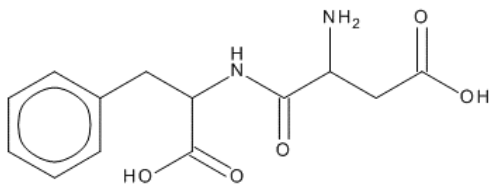
| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|--------|------------|
| 15(c)(ii) | <p>(alcoholic / ethanolic) ammonia / NH₃</p> <p>ALLOW Aqueous ammonia / NH₃(aq)</p> <p>IGNORE Concentration / heat</p> | | (1) |

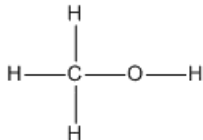
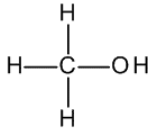
| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|--|---|------------|
| 15(c)(iii) | CH_3OH OR  OR  ALLOW CH_4O | CH_3OH^+ CH_4O^+ | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|--|------------|
| 15(c)(iv) |  OR  Both charges needed ALLOW Correct structure in brackets with charge outside and Na^+ | Bond between N and Na or between O and Na Partial charges | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|-----------------------|------------|
| 15(d)(i) |  <p>ALLOW displayed, structural or skeletal formula or any combination of these</p> <p>ALLOW -NH₃Cl as side group</p> <p>IGNORE Missing Cl⁻ ions</p> | Bond between N and Cl | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|---|------------|
| 15(d)(ii) |  <p>ALLOW No charges, provided there is no bond between O and Na</p> <p>ALLOW Displayed, structural or skeletal formula or any combination of these e.g. COONaCH₂CH(NH₂)COONa</p> <p>IGNORE Missing Na⁺ ions</p> | Partial charges δ^+ / δ^- | (1) |

| Question Number | Acceptable Answers | Mark |
|-------------------|--|------------|
| 15(d)(iii) | <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">  </div> <div style="text-align: center;"> <p>OR</p>  </div> </div> <p>OR</p> <div style="text-align: center;">  </div> <p>Any one amide/peptide link shown e.g. $\begin{array}{c} \text{O} \\ \parallel \\ \text{---C---N---} \end{array}$ /-CONH-</p> <p>Do not award this mark if other functional groups are joined directly to the amide (1)</p> <p>Rest of structure correct</p> <p>ALLOW Displayed or structural formula or any combination of these (1)</p> <p>IGNORE Bond angles and bond lengths</p> | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|-------------------|------------|
| 15(d)(iv) | <p>Methanol OR Any unambiguous formula e.g. CH₃OH /</p> <div style="text-align: center;">  <p>/</p>  </div> <p>If name and formula are given, both must be correct</p> | CH ₄ O | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--|------------|
| 15(e)(i) | <p>LiAlH_4 / lithium tetrahydridoaluminate / lithium aluminium hydride (in dry ether)</p> <p>OR</p> <p>NaBH_4 / sodium tetrahydridoborate / sodium borohydride (in aqueous / alcohol solution)</p> <p>IGNORE</p> <p>Lithal / heat</p> | <p>LiAlH_4 in water / aqueous solution</p> <p>Just '[H] / H⁻'</p> | (1) |

| Question Number | Acceptable Answers | | | Reject | Mark |
|------------------|---|---|----------------------------------|--|------------|
| 15(e)(ii) | Reagent | Glucose | Sorbitol | Observations not linked to a reagent Potassium dichromate(VI) | (3) |
| | Fehling's / Benedict's (solution and heat/boil) | Red / red-brown /brown / orange and precipitate | Stays blue / no change / no ppt | | |
| | Tollens' (reagent) / ammoniacal silver nitrate (and warm) | Silver mirror or grey /black /silver and precipitate | No change / no ppt | | |
| | Brady's (reagent) / 2,4-dinitro-phenyl-hydrazine / 2,4-DNP(H) | Orange / yellow / red and precipitate | Stays orange /no change / no ppt | | |
| | Reagent (1) | | | | |
| | Matching observation for glucose (1) | | | | |
| | Matching observation for sorbitol (1) | | | | |
| | ALLOW Correct formulae | | | | |
| | ALLOW No reaction / no observation / nothing happens for sorbitol | | | | |
| | IGNORE Sodium hydroxide in Tollens' reagent No TE on incorrect reagents | | | | |

| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|--|--------|------------|
| 15(e)(iii) | <div data-bbox="379 275 703 499"> </div> <p data-bbox="368 555 416 589">OR</p> <div data-bbox="379 600 703 824"> </div> <p data-bbox="368 880 416 913">OR</p> <div data-bbox="379 925 703 1070"> </div> <p data-bbox="368 1093 954 1193">ALLOW Structural or displayed formulae or any combination of these</p> <p data-bbox="368 1238 818 1339">IGNORE Bond angles and bond lengths OH connectivity</p> | | (1) |

(Total for Section C = 19 marks)
TOTAL FOR PAPER = 90 MARKS