

| Question Number | Correct Answer | Mark |
|-----------------|----------------|------|
| 1 (a) | A | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|------|
| 1 (b) | A | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|------|
| 1 (c) | D | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|------|
| 2 | B | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|------|
| 3 | D | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|------|
| 4 | C | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|------|
| 5 | D | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|------|
| 6 | D | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 7 | B | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 8 | C | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 9 | A | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 10 | C | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 11 | B | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 12 | C | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 13 | D | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 14 | C | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 15 | D | 1 |

| Question Number | Correct Answer | Mark |
|-----------------|----------------|----------|
| 16 | D | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--|------|
| 17 (a) | <p>Q: O-H</p> <p>ALLOW OH — O — H (1)</p> <p>R: C=O</p> <p>ALLOW — C = O — C = O (1)</p> <p>IGNORE names ACCEPT answers written on spectrum</p> | <p>Just 'alcohol' — OH</p> <p>Just 'carbonyl' — C O C-O</p> | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--------|------|
| (b) (i) | <p>Y = methanol / CH_3OH (1)</p> <p>Any two of the following: Molecular ion / M^+ / M_r / CH_3OH^+ / methanol = 32 CH_3^+ = 15 CH_3O^+ / CH_2OH^+ = 31 CHOH^+ / CH_2O^+ = 30 COH^+ = 29 CO^+ = 28 (1)</p> <p>Charges not required</p> <p>TE in second mark for two correct possible peaks from an incorrect compound.</p> | | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|---|------|
| 17 (b) (ii) | <p>Two (1) This mark may be scored if two shifts are given.</p> <p>Any two shifts correctly identified: -OH at 2.0-4.0 / any value in this range H-C-O at 3.0- 4.2 / any value in this range H in CH₃ OH at 3.39 (ppm) (1)</p> <p>Allow TE for ethanol with three peaks (1) and three correct shift values: -OH at 2.0-4.0 / any value in this range H-C-O at 3.0- 4.2 / any value in this range CH in an alkane at 0.1-1.9 (1)</p> | <p>CH in an alkane at 0.1-1.9</p> <p>Just CH₃ OH at 3.39</p> | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|------|
| (c) (i) | <p>Z contains two -OH/ one alcohol + one acid</p> <p>ALLOW two alcohol groups / is a diol</p> | | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|------|
| (c) (ii) | <p>Z is an acid / contains -COOH / contains -CO₂H/ contains a carboxylic acid group / contains H⁺</p> | | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---------------|------|
| (c) (iii) | <p>Z is a secondary alcohol/ a ketone is formed from Z /</p> <p>Z contains $\begin{array}{c} \\ \text{-C-OH} \\ \\ \text{H} \end{array}$ (1)</p> | Z is a ketone | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--------|------|
| (c) (iv) | <p>(Iodoform produced) so Z contains CH₃ CH(OH)-</p> <p>TE if Z is identified as a ketone in (iii): Z contains CH₃ C=O / Z is a methyl ketone</p> | | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|------|
| 17 (d) | <p>Answers will be based on several pieces of information (molecular formula, products of ester hydrolysis, answers to (c)) which may be contradictory if errors have been made.</p> <p>ALLOW TE marks for formulae which are chemically possible (ie no 5 bonded carbons etc) and based on most of the deductions but not necessarily all.</p> <p>Z is $\text{CH}_3\text{CH}(\text{OH})\text{CH}_2\text{COOH}$ (1) Stand alone mark</p> <p>ALLOW TE for an acid with OH in wrong position in Z if oxidation product identified as aldehyde</p> <p>TE for Z = $\text{CH}_3\text{COCH}_2\text{COOH}$ if identified as ketone in (iii)</p> <p>X is $\text{CH}_3\text{CH}(\text{OH})\text{CH}_2\text{COOCH}_3$ (1) Stand alone mark TE for a methyl ester of Z</p> | | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|----------------------------|------|
| 18 (a) (i) | Transesterification Ethanol transesterification | Substituted esterification | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|---|------|
| 18 (a) (ii) | To prevent hydrolysis/ to stop fatty acids forming / to stop breakdown of esters / water reacts with esters/ water is a better nucleophile than ethanol | To dilute ethanol Ethanol would react with water A reaction would occur (unspecified) | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---|------|
| 18 (b) | <p>(Vegetable) Fats/ oils are renewable (crude oil is not) /</p> <p>biodiesel comes from a renewable source /</p> <p>doesn't use up fossil fuel resources/</p> <p>carbon footprint is less /</p> <p>(closer to) carbon neutral /</p> <p>growing vegetables absorb CO₂</p> <p>If more than one answer is given, and one is incorrect, no mark</p> <p>Ignore comments on biodegradability</p> | <p>Just "made from plants"</p> <p>Just "crude oil is not sustainable"</p> <p>Less polluting produces less greenhouse gases / less CO₂</p> <p>Burns more cleanly</p> <p>Requires less energy for production</p> | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--|------|
| (c) | <p>Substances to be separated have different (forces of) attraction to / affinity for / solubilities in / adsorption to one or both of the mobile and stationary phases OWTTE (1)</p> <p>ALLOW absorption</p> <p>GC: mobile phase a (inert / unreactive) gas OR GC: mobile phase nitrogen / helium / argon / other named inert gas (1)</p> <p>GC: Stationary phase a liquid (on an (inert) solid) / a solid (1)</p> <p>HPLC: stationary phase a solid / silica (1)</p> <p>HPLC: mobile phase a liquid (1)</p> | <p>Different retention times without a reason why</p> <p>Different volatilities</p> <p>Different masses</p> <p>Different reactivity</p> <p>Different reactions</p> <p>Different interactions</p> | 5 |

TOTAL FOR SECTION C = 20 MARKS

| | | |
|--|---|--|
| <p>ALLOW gas given off that turns limewater cloudy</p> <p>OR</p> <p>Test : with blue litmus (1)</p> <p>Result: turns red (1)</p> <p>The test can be with any other indicator, including universal indicator, with the correct initial and final colour</p> <p>ALLOW pH meter (1)</p> <p>pH 4-6 (1)</p> <p>OR</p> <p>Test : add ethanol with conc H_2SO_4 (and warm) (1)</p> <p>Result: gives pleasant/fruity smell of ester (1)</p> <p>OR</p> <p>Test: add magnesium (1)</p> <p>Result: fizzing/bubbles etc (of hydrogen) (1)</p> <p>ALLOW gas given off that burns with a squeaky pop</p> | <p>Add sodium</p> <p>colourless gas evolved</p> | |
|--|---|--|

| Question Number | | Reject | Mark |
|------------------|--|---|----------|
| 19 (b)(i) | <p>Explanation of precedence/priority in terms of atomic numbers/masses of the attached groups</p> <p>OR</p> <p>Highest-precedent/priority groups on each carbon are on opposite sides of the molecule (1)</p> <p><i>E</i>-/entgegen (1)</p> <p>Mark independently</p> | Both CH ₃ /methyl groups on the same side so Z (0/2) | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|----------|
| (b)(ii) | <p>45 COOH⁺ /CO₂H⁺ (1)</p> <p>55 C₄H₇⁺</p> <p>OR</p> <p>C₃OH₃⁺ (1)</p> <p>ALLOW Structural/displayed formulae of ions</p> <p>Absence of + charge (1 max)</p> | | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|-----------------|----------|
| (b)(iii) | <p>If they say yes (0)</p> <p>(No) (Cleavage of the C—COOH bond in) both compounds gives fragment(s) of the same mass OR Both give the same peak(s)/fragment(s)</p> <p>Both give CO₂H⁺/ C₄H₇⁺ fragments</p> <p>The mark can be scored by referring to just one of the fragments/peaks/masses.</p> | 'No' on its own | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|--------------------------------|--|--------------------------------------|----------|
| *19(c)(i) QWC | <p>C is CH_3CHO (alone) (2)</p> <p>D is CH_3COCOOH (alone) (2)</p> <p>so tiglic acid must be B (1)</p> <p>tiglic acid mark can only be awarded if correct structures of either C or D are given.</p> <p>Any one of the following</p> <p>C must be an aldehyde (1)</p> <p>D is a ketone (1)</p> <p>Mention that CH_3CO present in either/both compounds (because of formation of iodoform) (1)</p> <p>If one or both of the structures are incorrect any of the last 3 marks can be awarded max 5</p> <p>If C and D are fully correct, but the wrong way round max 5</p> | CH_3COH 1 max | 6 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--|----------|
| (c)(ii) | <p>Doesn't distinguish <i>E</i>- isomer from <i>Z</i>- isomer/geometric isomers (so no)</p> <p>OR</p> <p>Doesn't distinguish which sides of $\text{C}=\text{C}$ functional groups are on</p> | Just isomers/ stereoisomers/ enantiomers | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---|------|
| (d)(i) | <p>CH₃CHO (1)</p> <p>ACCEPT displayed or skeletal</p> <p>Step 1</p> <p>(heat)using acidified potassium dichromate/or H⁺/Cr₂O₇²⁻ (1)</p> <p>distil (product as formed) conditional on dichromate (1)</p> <p>Step 2</p> <p>HCN with KCN</p> <p>OR</p> <p>KCN with H⁺/acid</p> <p>OR</p> <p>KCN with (cold) NaOH(aq)/alkali (1)</p> <p>ALLOW HCN with NaOH/alkali</p> <p>For step 2 Ignore conditions e.g. any references to heat</p> | <p>CH₃COH</p> <p>Manganate VII/KMnO₄</p> <p>Reflux</p> <p>HCN alone</p> | 4 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|---|------|
| (d)(ii) | <p>Nucleophilic addition</p> <p>Any recognisable spelling of 'philic' and addition, either order</p> <p>Both words needed</p> | <p>Nutrophilic addition</p> <p>Any other or additional words</p> | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|----------------------------------|--|---|----------|
| *19(d)(iii) QWC | <p>Ethanal is planar (at the reaction site)</p> <p>OR</p> <p>Ethanal is a planar molecule (1)</p> <p>Attack (from CN^- to give the cyanohydrin) is (equally likely) from either side/above or below/from both sides (of the molecule) (so a racemic mixture is formed) (1)</p> <p>Mark independently</p> | <p>Intermediate is planar Square planar</p> <p>Can attack carbocation from either side/any reference to SN1/SN2</p> | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--------|----------|
| (d)(iv) | <p>Receptors for the compound in the body are often stereospecific so only one stereoisomer is pharmacologically active</p> <p>OR</p> <p>Body recognises one (stereo)isomer</p> <p>ALLOW</p> <p>Only one (stereo)isomer is active</p> <p>OR</p> <p>One/the other isomer may be toxic/dangerous/harmful</p> <p>OR</p> <p>One isomer destroys body cells</p> <p>OR</p> <p>(Different) isomers have different biological/pharmacological/biochemical properties</p> | | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--------|----------|
| 20(a)(i) | <p>Formula showing -NH_3^+ and -COO^- / -CO_2^-</p> <p>Charges can be anywhere on functional group</p> <p>Rest of the molecule must be correct</p> <p>ALLOW displayed/part displayed formula</p> | | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---|----------|
| (a)(ii) | <p>Any two from</p> <p>High energy needed (to overcome) (1)</p> <p>strong ionic/electrostatic forces OR strong forces between oppositely charged ions/between positive and negative (1)</p> <p>between different (zwitter)ions</p> <p>OR</p> <p>between -NH_3^+ and -COO^-</p> <p>OR</p> <p>between one molecule and another</p> <p>OR</p> <p>Chains of zwitterions/molecules (1)</p> | <p>any reference to intermolecular forces eg (strongly) polar/bond polarity</p> <p>if they state the ionic bond is within the same molecule</p> | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|--|--------|----------|
| 20(a)(iii) | <div data-bbox="416 336 751 453" data-label="Chemical-Block"> </div> <p data-bbox="411 502 975 538">Correct peptide link (1)</p> <p data-bbox="411 572 975 644">Minimum two residues and extension to the rest of the molecule (1)</p> <p data-bbox="411 678 975 751">ALLOW -NHCH₂CONHCH₂CO- (2)</p> <p data-bbox="411 785 890 857">Drawn the other way round, i.e. starting with the carbonyl group</p> <p data-bbox="411 891 927 963">Brackets around outside with 'n' ie (.....)_n</p> <p data-bbox="411 998 858 1034">Second mark depends on first</p> | | 2 |

| Question Number | Acceptable Answer | Reject | Mark |
|-----------------------------|---|---|----------|
| *20(b) QWC | <p>Key Points</p> <p>KP1 Spot (of hydrolysate) on paper/tlc/thin layer chromatogram (1)</p> <p>KP2 Marker spots of known amino-acids/measure R_f (1)</p> <p>KP3 Run in (suitable) solvent/discussion of comparative solubilities in phases (1)</p> <p>KP4 (Spray with) ninhydrin (and heat) [Stand alone mark] (1)</p> <p>KP 5 Marker spots and the unknown spots correspond ALLOW Compare R_f values of marker spots with hydrolysate spots (1)</p> <p>OR</p> <p>If 2-d chromatography used (2 different solvents run in two directions at right angles):</p> <p>KP1 Spot (of hydrolysate) on paper/tlc/thin layer chromatogram (1)</p> <p>KP2 Run in (suitable) solvent in one direction (1)</p> <p>KP3 Develop in suitable/different solvent at right angles OR discussion of comparative solubilities in phases (1)</p> <p>KP4 Spray with ninhydrin (and heat) (1)</p> <p>KP5 Compare hydrolysate spots with same experiment for known amino acids (1)</p> <p>OR</p> | <p>Spot one amino acid/protein</p> <p>Water alone as solvent</p> <p>Spot one amino acid</p> | 5 |

| | | | |
|--|---|---------------------|--|
| | if column/GLC/GC used | | |
| | KP1 Put amino acid mixture (Hydrolysate) into column (1) | Spot one amino acid | |
| | KP2 Separately known amino-acids into column (1) | | |
| | KP3 Detect amino acids in effluent with Ninhydrin/mass spectrometry (1) | | |
| | KP4 Measure retention times/ discussion of comparative solubilities in phases (1) | | |
| | KP 5 Compare retention times (1) | | |

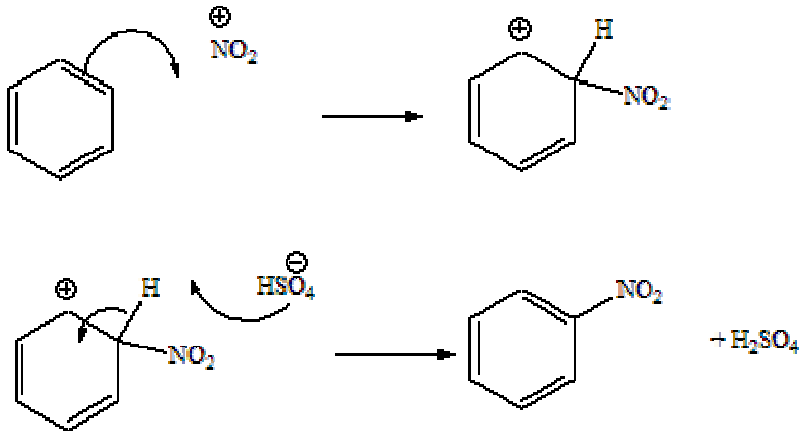
S

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|----------|
| 21(a)(i) | Not knowing the structure of the molecule (means that the reactions/reagents/reactants needed to make it are also unknown) ALLOW Structure not known | | 1 |

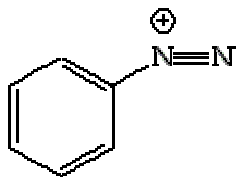
| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|----------|
| (a)(ii) | <p>Credit any reasonable arguments for example:</p> <p>First mark No longer any demand for madder/indigo</p> <p>OR</p> <p>Cheaper alternatives available (1)</p> <p>Second mark So the growing industries collapsed</p> <p>OR</p> <p>no market for crops</p> <p>OR</p> <p>farmers had to grow alternative crops</p> <p>OR</p> <p>decreased employment</p> <p>OR</p> <p>economic damage</p> <p>OR</p> <p>decreased GDP</p> <p>OR</p> <p>Loss of export (1)</p> | | 2 |

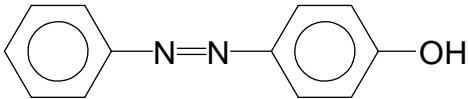
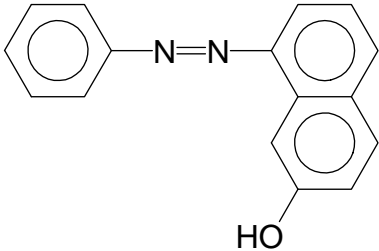
| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|----------|
| 21(b)(i) | <p>First mark Double bonds expected to react with bromine water turning it colourless</p> <p>OR</p> <p>Bromine water remained yellow/orange/red/brown (1)</p> <p>Second mark So benzene does not contain double bonds</p> <p>OR</p> <p>Double bonds not normal/not simply double bonds/any indication that double bonds are different</p> <p>OR</p> <p>His representation incorrect (1)</p> | | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|---------------------------------|----------|
| (b)(ii) | <p>The p/pi-/π/6 electrons (of carbon)</p> <p>OR</p> <p>π system (1)</p> <p>Electrons are delocalised around the ring (1)</p> <p>Which gives the molecule greater stability/need more energy to break the bonds in benzene (and hence a less exothermic hydrogenation enthalpy)</p> <p>Allow it is more stable (1)</p> | Harder to break/disrupt [alone] | 3 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|----------|
| 21(c) | <p> $2\text{H}_2\text{SO}_4 + \text{HNO}_3 \rightarrow \text{NO}_2^+ + \text{H}_3\text{O}^+ + 2\text{HSO}_4^-$ (1) OR $\text{H}_2\text{SO}_4 + \text{HNO}_3 \rightarrow \text{NO}_2^+ + \text{H}_2\text{O} + \text{HSO}_4^-$ OR $\text{H}_2\text{SO}_4 + \text{HNO}_3 \rightarrow \text{H}_2\text{NO}_3^+ + \text{HSO}_4^-$ and $\text{H}_2\text{NO}_3^+ \rightarrow \text{NO}_2^+ + \text{H}_2\text{O}$ </p> <p>Charges are needed for first mark</p>  <p>Attack on nitronium ion arrow must start on or in the benzene (1)</p> <p>Wheland intermediate Can be a part, but not complete circle, in correct place inside ring BUT part circle must cover minimum of 3 carbon atoms AND must not include where nitro group is attached and must positive charge somewhere (1)</p> <p>Either but only one of first two marks can be lost if bond is clearly to oxygen</p> <p>Arrow from H bond into the ring to produce either H^+ or H_2SO_4 and return to aromaticity (1)</p> | | 4 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---------------------|----------|
| 21(d)(i) | <p>< 0° C/temperature too low: reaction too slow/insufficient energy to overcome activation energy (1)</p> <p>> 10° C/temperature too high: diazonium ion decomposes/produces phenol (1)</p> | Will not take place | 2 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|----------------------------|----------|
| (d)(ii) |  <p>Positive charge can be on either N</p> <p>Cl⁻ may be given as well</p> <p>ALLOW circle in benzene ring and hydrogens/carbons displayed</p> <p>OR</p> <p>---N=N⁺ Is acceptable providing charge is on the end N</p> | Positive charge on wrong N | 1 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---|----------|
| (d)(iii) | <p>Adds phenol in sodium hydroxide/OH^-/alkali</p> <p>ALLOW 2-naphthol in sodium hydroxide/OH^- /alkali (1)</p> <p>Correct structure for the $-\text{N}=\text{N}-$ bond between 2 benzene rings (1)</p> <p>Remainder of molecule (1)</p> <p>which is either:</p> <div style="text-align: center;">  </div> <p>ALLOW anionic form of $-\text{OH}$</p> <p>OR if 2-naphthol is used it is:</p> <div style="text-align: center;">  </div> | Ignore position of $-\text{OH}$ group on the ring | 3 |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|-------------------------------------|----------|
| 21 (e) | <p>First mark</p> <p>—SO_3^- are solvated / hydrated</p> <p>Can be drawn with polar H of water</p> <p>OR</p> <p>Negative ion bonds with/attracted to water (1)</p> <p>Second mark</p> <p>Nitrogen/oxygen atoms hydrogen-bonded (to water)</p> <p>Can be drawn (1)</p> | Just sodium ions attracted to water | 2 |