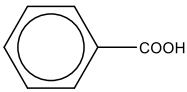
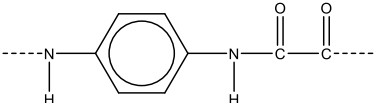
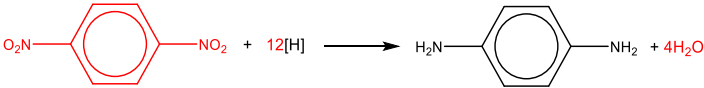
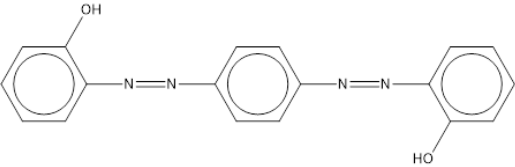
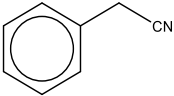
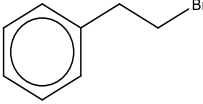
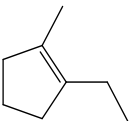
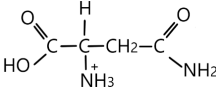
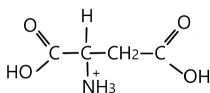


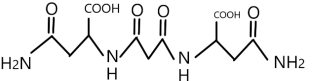
|          |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               |
|----------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| 7(a)     | <p><b>M1:</b> <math>\text{ClCH}_2\text{COOH}</math> <math>\text{BrCH}_2\text{COOH}</math> <math>\text{CH}_3\text{COOH}</math> <math>\text{CH}_3\text{CH}_2\text{OH}</math><br/>most acidic least acidic</p> <p><b>M2:</b> electron withdrawing groups / electronegative groups / negative inductive effect<br/><b>AND</b> weakens O–H bond / more stable anion</p> <p><b>OR</b><br/>electron donating groups / positive inductive effect <b>AND</b> strengthens O–H bond / less stable anion</p> <p><b>M3 M4:</b></p> <ul style="list-style-type: none"> <li>chlorine is more electron-withdrawing / electronegative than bromine (related to <math>\text{XCH}_2\text{COOH}</math>)</li> <li>electron withdrawing of <math>\text{C=O}</math> / negative inductive effect of <math>\text{C=O}</math> (related to <math>\text{COOH}</math>)</li> <li>positive inductive effect / electron donating of alkyl / ethyl / R group (related to <math>\text{ROH}</math>)</li> </ul> <p>any two [1], all three [2]</p> |
| 7(b)     | <p><b>M1:</b></p>  <p><b>M2:</b> <math>\text{CO}_2</math></p>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                |
| 7(c)(i)  |  <p><b>M1:</b> correct displayed amide bond with <math>\text{C=O}</math>, <math>(\text{C}_6\text{H}_5)\text{-N}</math> and <math>(\text{C-})\text{C=O}</math></p> <p><b>M2:</b> rest of the structure correct with continuation bonds</p>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| 7(c)(ii) |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| 7(d)(i)  | <p><b>M1:</b> <math>\text{Cl}^- \text{N} \equiv \text{N}^+ - \text{C}_6\text{H}_4 - \text{N}^+ \equiv \text{N} \text{Cl}^-</math></p> <p><b>M2:</b></p>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     |

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| 7(d)(ii) | <p><math>\text{HNO}_2</math> (and <math>\text{HCl}</math>) <b>AND</b> <math>\leq 10^\circ\text{C}</math></p> <p><b>OR</b></p> <p><math>\text{NaNO}_2</math> <b>AND</b> <math>\text{HCl}</math> <b>AND</b> <math>\leq 10^\circ\text{C}</math></p>                              |
| 9(a)     | <p><b>M1:</b> (basicity linked to) ability of lone pair / p-orbital to <b>AND</b> accept / coordinate with a proton / <math>\text{H}^+</math></p> <p><b>M2:</b> (lone pair of) electrons on N is delocalised into <math>\text{C=O}</math> group (and make amides neutral)</p> |

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| 9(b)(i)  | <p><b>M1: M</b></p>  <p><b>M2: N</b></p>                                                                                                                                             |
| 9(b)(ii) | $\text{LiAlH}_4$                                                                                                                                                                                                                                                                                                                                           |
| 9(c)     | ketone / carbonyl <b>AND</b> secondary / $2^\circ$                                                                                                                                                                                                                                                                                                         |
| 9(d)(i)  |  <p>curly arrow 1: from O lone pair to C atom (in <math>\text{C=C}</math>)<br/>curly arrow 2: from <math>\text{C=C}</math> bond to the right-hand O atom in <math>\text{O=O}</math><br/>curly arrow 3: from <math>\text{O=O}</math> bond to <math>\text{O}^+</math></p> |

|          |                                                                                     |
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| 9(d)(ii) |  |
|----------|-------------------------------------------------------------------------------------|

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|------|---------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| 8(a) | the pH at which an amino acid exists as a zwitterion                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
| 8(b) | <p><b>M1:</b> <math>[\text{HOOCCH}(\text{NH}_3)\text{CH}_2\text{CONH}_2]^+</math></p> <p><b>M2:</b> <math>[\text{HOOCCH}(\text{NH}_3)\text{CH}_2\text{COOH}]^+</math></p> <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">  </div> <div style="text-align: center;">  </div> </div> |

|          |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                       |
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| 8(c)     | <p><b>M1:</b> HOCH<sub>2</sub>CH(NH<sub>2</sub>)CH<sub>2</sub>CH<sub>2</sub>NH<sub>2</sub><br/> <b>M2:</b> HOCH<sub>2</sub>CH(NH<sub>2</sub>)CH<sub>2</sub>CH<sub>2</sub>OH</p> <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;"> <math display="block">\begin{array}{c} \text{H} \\   \\ \text{HOCH}_2 - \text{C} - \text{CH}_2 - \text{CH}_2\text{NH}_2 \\   \\ \text{NH}_2 \end{array}</math> </div> <div style="text-align: center;"> <math display="block">\begin{array}{c} \text{H} \\   \\ \text{HOCH}_2 - \text{C} - \text{CH}_2 - \text{CH}_2\text{OH} \\   \\ \text{NH}_2 \end{array}</math> </div> </div> |
| 8(d)(i)  | HOOCCH <sub>2</sub> COOH + 2SOCl <sub>2</sub> → ClOCCH <sub>2</sub> COCl + 2SO <sub>2</sub> + 2HCl [1]                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                |
| 8(d)(ii) | <p>HOOCCH(CH<sub>2</sub>CONH<sub>2</sub>)NHCOCH<sub>2</sub>CONHCH(CH<sub>2</sub>CONH<sub>2</sub>)COOH</p> <p><b>M1:</b> three molecules joined, with four correct amide groups only<br/> <b>M2:</b> whole structure correct as above, displayed formula accepted, skeletal formula as below accepted</p>                                                                                                                                                                                                                                                                                             |
| 8(e)     | <p><b>M1:</b> hydrolysis of amide to carboxylate ion<br/> <b>M2:</b> hydrolysis of –COOH to carboxylate ion and rest of molecule</p> <p>–OOCCH(NH<sub>2</sub>)CH<sub>2</sub>COO<sup>–</sup></p>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                       |

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| 8(f)     | <ul style="list-style-type: none"> <li>three monomer residues</li> <li>all linkages correct</li> <li>correct trailing bonds</li> <li>repeat unit marked contains correct atoms</li> </ul> <p>Any two [1], any three [2], all four [3]</p> <div style="text-align: center;"> <math display="block">\begin{array}{c} \text{CH}_2\text{CONH}_2 \qquad \qquad \text{CH}_2\text{CONH}_2 \\   \qquad \qquad \qquad   \\ -\text{COCHNH}-\left[ \text{COCHNH} \right]_n-\text{COCHNH}- \\   \\ \text{CH}_2\text{CONH}_2 \end{array}</math> </div> |
| 8(g)(i)  | <ul style="list-style-type: none"> <li>rotate the plane of the plane polarised light</li> <li>by the same angle</li> <li>in opposite directions</li> </ul> <p>Any two [1], all three [2]</p>                                                                                                                                                                                                                                                                                                                                              |
| 8(g)(ii) | racemic mixture                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |