

## Mark schemes

1

- (a) Heating speeds up (hydrolysis / breaking of peptide bonds)

**OR** forms non-sweet (amino acids)

1

- (b) (2-)aminobutanedioic acid OR

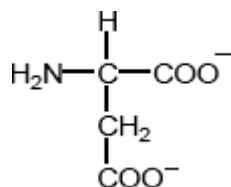
*2 not necessary but penalise other numbers at start*

(2-)aminobutane(-1,4-)dioic acid

*1,4 not necessary but penalise other numbers and 1,4 must be in correct place (QoL)*

1

- (c)

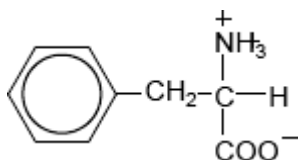


*allow  $-\text{CO}_2^-$*

*allow  $\text{NH}_2-$*

1

- (d)



*allow  $-\text{CO}_2^-$*

*allow  $+\text{NH}_3-$*

*don't penalize position of + on  $\text{NH}_3$*

1

- (e) (i) **M1** Compounds/molecules with same structural formula

*Not just structure*

1

**M2** But with bonds/atoms/groups arranged differently in space or in 3D

*Allow –with different spatial arrangement of atom/bond/group*

1

*Independent marks*

- (ii) (Plane) polarised light

1

Rotated in opposite directions

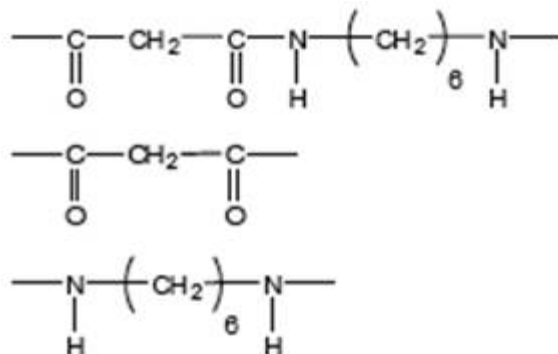
*Not bent or turned or twisted; not different directions (QoL)*

1

[8]

2

(a) (i)



Allow  $\text{---CONH---}$  or  $\text{---COHN---}$

*Mark two halves separately*

*lose 1 each for missing trailing bonds at one or both ends or error in peptide link or either or both of H or OH on ends*

1

Not allow  $\text{---(C}_6\text{H}_{12}\text{)---}$

*Ignore n*

1

(ii) **M1** in polyamides - H bonding

1

**M2** in polyalkenes - van der Waals forces

*Penalise forces between atoms or van der Waals bonds*

1

**M3** Stronger forces (of attraction) in polyamides

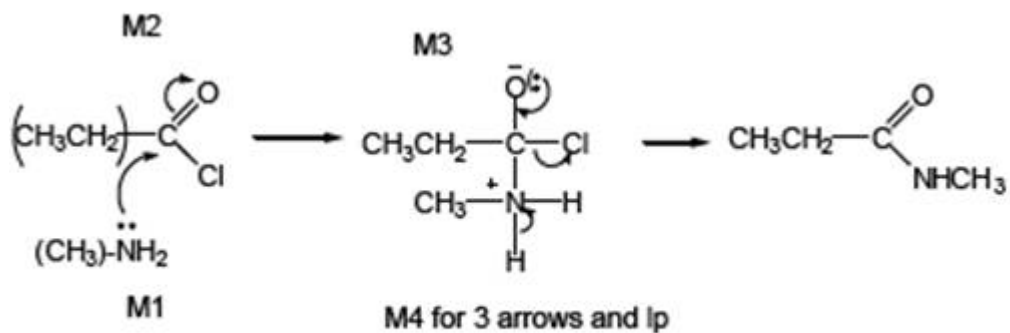
Or H bonding is stronger

(must be a comparison of correct forces to score M3)

*Do not award if refer to stronger bonds*

1

(b) (i) (nucleophilic) addition elimination



Not allow N-H<sub>2</sub>

*Minus sign on NH<sub>2</sub> loses M1*

1

*M2 not allowed independent of M1, but allow M1 for correct attack on C+*

*+ rather than δ+ on C=O loses M2*

*If Cl lost with C=O breaking, max 1 for M1*

*M3 for correct structure with charges but*

*lp on O is part of M4*

*only allow M4 after correct/ very close M3*

*For M4, ignore NH<sub>3</sub> removing H<sup>+</sup> but lose*

*M4 for Cl removing H<sup>+</sup> in mechanism,*

*but ignore HCl as a product*

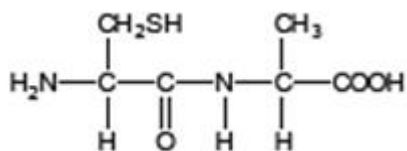
4

(ii) N-methylpropanamide

*Not N-methylpropaneamide*

1

(c)



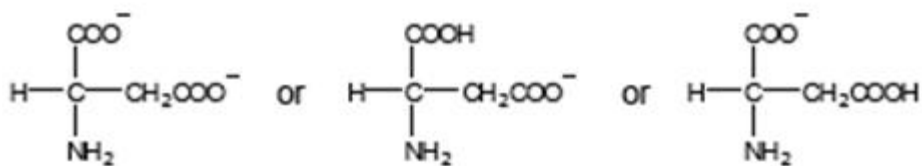
*Allow -CONH- or -COHN-*

1

(d) (i) 2-amino-3-hydroxypropanoic acid

1

(ii)



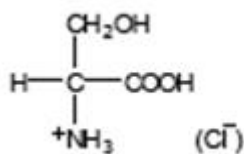
Must be salts of aspartic acid

*allow*  $-\text{CO}_2^-$

*allow*  $\text{NH}_2^-$

1

(iii) Penalise use of aspartic acid once in d(iii) and d(iv)



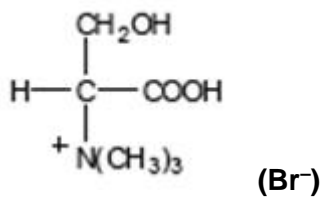
*allow*  $-\text{CO}_2\text{H}$

*allow*  $\text{}^+\text{NH}_3-$

*don't penalize position of + on  $\text{NH}_3$*

1

(iv) Penalise use of aspartic acid once in d(iii) and d(iv)



*allow*  $-\text{CO}_2^-$

*must show C-N bond*

*don't penalize position of + on  $\text{N}(\text{CH}_3)_3$*

1

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3

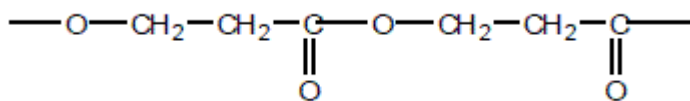
(a) 3-hydroxypropanoic acid

*allow* 3-hydroxypropionic acid

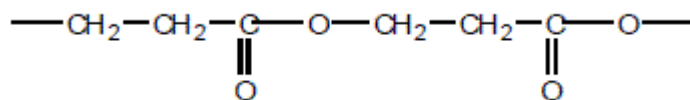
*must be correct spelling*

1

- (b) (i) must show trailing bonds



or can start at any point in the sequence, e.g.



*not allow dimer*

*allow  $\text{---O---CH}_2\text{CH}_2\text{COOCH}_2\text{CH}_2\text{CO---}$*

*or  $\text{---CH}_2\text{CH}_2\text{COOCH}_2\text{CH}_2\text{COO---}$*

*ignore ( ) or n*

*NB answer has a total of 6 carbons and 4 oxygens*

1

- (ii) condensation (polymerisation)

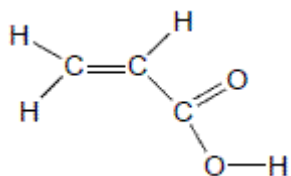
*Allow close spelling*

1

- (c) (i) C=C or carbon-carbon double bond

1

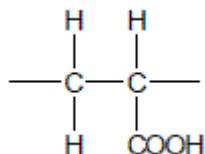
- (ii)



*must show **ALL** bonds including O-H*

1

- (iii) must show trailing bonds

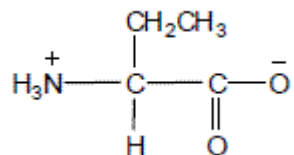


*allow polyalkene conseq on their c(ii)*

*ignore n*

1

(d)

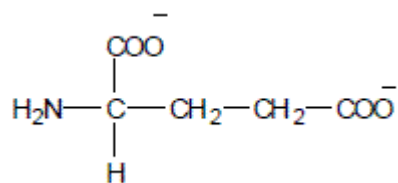


allow  $\text{NH}_3^+$  —

allow  $\text{COO}^-$

1

(e) (i)



In (e), do not penalise a slip in the number of carbons in the  $-\text{CH}_2\text{CH}_2-$  chain, but all must be bonded correctly

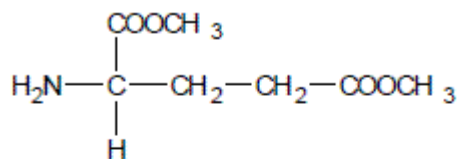
NB two carboxylate groups

Allow  $\text{COONa}$  or  $\text{COO}^- \text{Na}^+$  but not covalent bond to Na

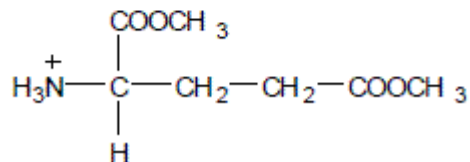
allow  $\text{NH}_2-$

1

(ii)



OR



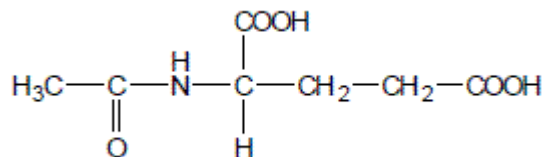
In (e), do not penalise a slip in the number of carbons in the  $-\text{CH}_2\text{CH}_2-$  chain, but all must be bonded correctly

NB two ester groups

allow  $\text{NH}_2-$  or  $^+\text{NH}_3-$

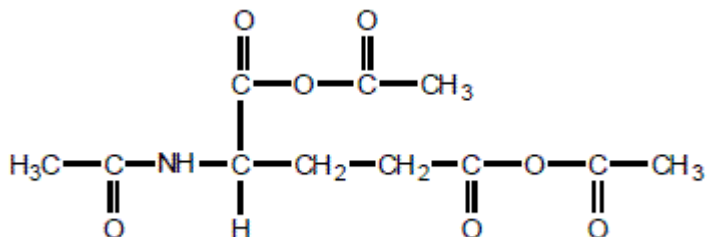
1

(iii)



In 4(e), do not penalise a slip in the number of carbons in the  $-\text{CH}_2\text{CH}_2-$  chain, but all must be bonded correctly

allow anhydride formation on either or both COOH groups (see below) with or without amide group formation



1

(f) **M1** phase or eluent or solvent (or named solvent) is moving or mobile

1

**M2** stationary phase or solid or alumina/silica/resin

1

**M3** separation depends on balance between solubility or affinity (of compounds) in each phase

**OR**

different adsorption or retention

**OR**

(amino acids have) different  $R_f$  values

**OR**

(amino acids) travel at different speeds or take different times

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4

(a) (i) hydrolysis

*not hydration*

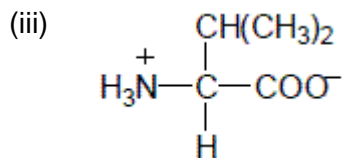
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(ii) 2-aminopropanoic acid

*ignore alanine*

*QoL*

1

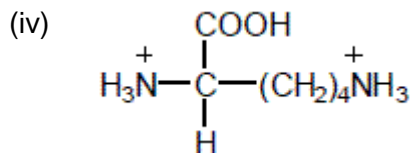


allow  $-\text{CO}_2^-$

allow  $^+\text{NH}_3-$

don't penalize position of + on  $\text{NH}_3$

1

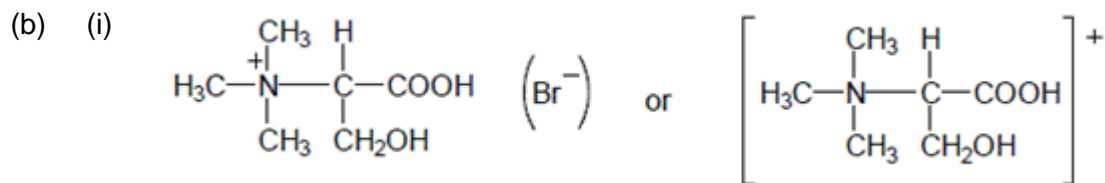


allow  $-\text{CO}_2^-$

allow  $^+\text{NH}_3-$

don't penalize position of + on  $\text{NH}_3$

1

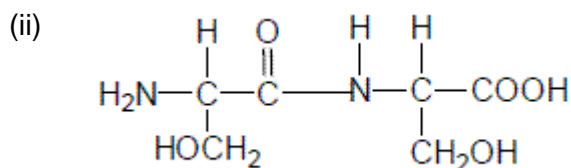


allow  $-\text{CO}_2\text{H}$

allow limit as  $\begin{array}{c} | \\ -\text{C}- \\ | \\ \text{CH}_2\text{OH} \end{array}$

+ on N or outside [ ]

1



allow  $-\text{CO}_2\text{H}$  allow  $-\text{CONH}-$  or  $-\text{COHN}-$

allow  $\text{NH}_2-$

allow limit as  $\begin{array}{c} | \\ -\text{C}- \\ | \\ \text{CH}_2\text{OH} \end{array}$

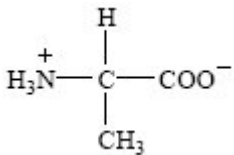
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[6]



5

(a) (i)



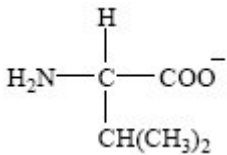
allow  $-\text{CO}_2^-$

allow  $^+\text{NH}_3-$

don't penalize position of + on  $\text{NH}_3$

1

(ii)



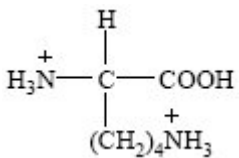
allow  $-\text{CO}_2^-$

allow  $\text{NH}_2-$

allow  $\text{C}_3\text{H}_7$

1

(iii)



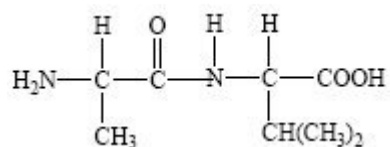
allow  $-\text{CO}_2\text{H}$

allow  $^+\text{NH}_3-$

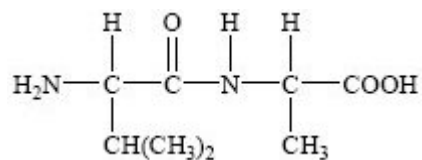
don't penalize position of + on  $\text{NH}_3$

1

(b)



1



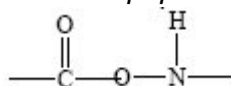
allow  $-\text{CO}_2\text{H}$

allow  $\text{NH}_2-$

allow  $\text{C}_3\text{H}_7$

allow as zwitterions

if error in peptide link e.g.



if twice, penalise both times

not polymers

if wrong amino acid in both can score Max 1

1

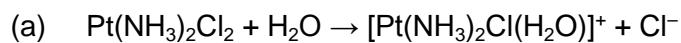
(c) chromatography or electrophoresis

ignore qualification to chromatography

1

[6]

6



Correct product

1

Balanced equation

1

(b) (i) Hydrogen bond

1

Oxygen (or nitrogen)

Only score this mark if type of bond is correct

1

(ii) Co-ordinate

1

Nitrogen (or oxygen)

Bond type must be correct to score this mark but allow M2 if bond is covalent

1

- (c) Killing them or causing damage (medical side effects)  
 Allow any correct side effect (e.g. hair loss)  
 Allow kills healthy (or normal) cells

1

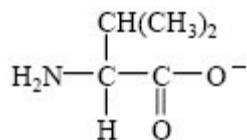
May attach to DNA in normal cells

1

[8]

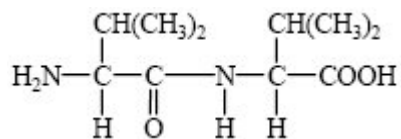
7

- (a) (i)



1

- (ii)

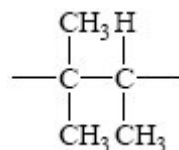


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- (iii) hydrogen bonding (do not allow H-bonding) QWC  
 do not penalise any error twice.

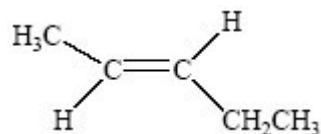
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- (b) (i)



1

- (ii)

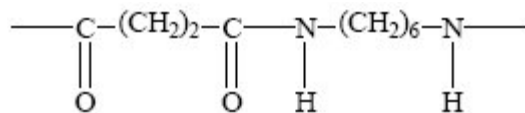


1

- (iii) Isomer must be saturated or must not contain a double bond

1

(c)



2

(d) (i) heat/reflux with aqu NaOH

1

poly(alkene) is inert/ no reaction

1

polyamide is hydrolysed (or undergoes hydrolysis)  
to form acid salt and alcohol QWC

1

(ii) e.g combustion

1

heat energy produced

1

toxic gases produced

1

[14]