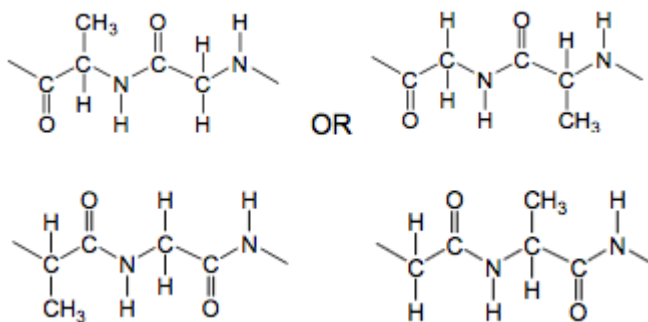


Mark schemes

1



(a) (i)

Only one molecule of each used

M1 for 2 amide links

M2 CH₂ and CH(CH₃)

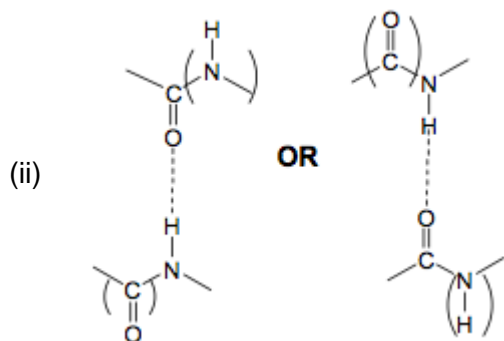
Allow 1 mark after one error

Dipeptide max 1

Treat both trailing bonds missing as one error

Ignore n

2



No need to show lp

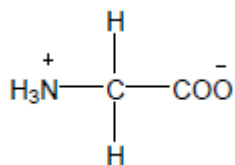
The covalent bond and the hydrogen bond either side of the H must be linear.

Allow



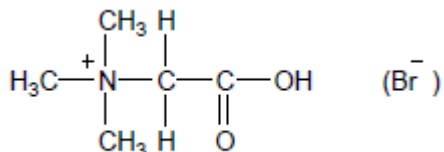
1

(b)

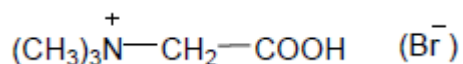


1

(c)



Allow

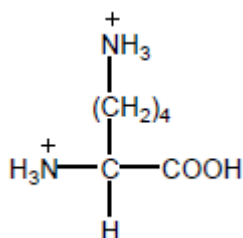


1

(d) 2-amino-3-hydroxybutanoic acid

1

(e)



1

[5]

4

(a) 2-deoxyribose

1

(b) Base A

If Base B stated, allow 1 mark only for response including hydrogen bonding

1

Top N-H forms hydrogen bonds to lone pair on O of guanine

1

The lone pair of electrons on N bonds to H-N of guanine

1

A lone pair of electrons on O bonds to lower H-N of guanine

Allow all 4 marks for a correct diagram showing the hydrogen bonding

Students could also answer this question using labels on the diagram

1

- (c) Allow either of the nitrogen atoms with a lone pair NOT involved in bonding to cytosine 1
- (d) Use in very small amounts / target the application to the tumour 1

[7]

5

- (a) **Wear plastic gloves:**
Essential – to prevent contamination from the hands to the plate 1

Add developing solvent to a depth of not more than 1 cm³:

Essential – if the solvent is too deep it will dissolve the mixture from the plate 1

Allow the solvent to rise up the plate to the top:

Not essential – the R_f value can be calculated if the solvent front does not reach the top of the plate 1

Allow the plate to dry in a fume cupboard:

Essential – the solvent is toxic
Allow hazardous 1

- (b) Spray with developing agent or use UV 1

Measure distances from initial pencil line to the spots (*x*) 1

Measure distance from initial pencil line to solvent front line (*y*) 1

R_f value = x / y 1

- (c) Amino acids have different polarities 1

Therefore, have different retention on the stationary phase or different solubility in the developing solvent 1

[10]

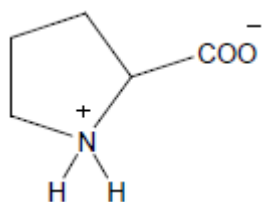
6

C

[1]

7

(a) (i)



Allow CO₂⁻ and NH₂⁺

1

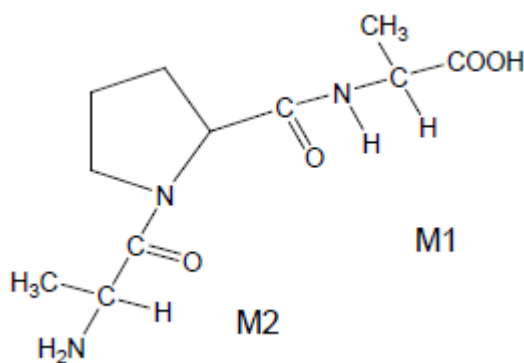
(ii) NOTE – **Two** marks for this clip

M1 for alanine section bonded through N

M2 for alanine section bonded through C

But penalise error in proline ring

1



Allow MAX 1 for correct tripeptide in polymer structure

1

(b) (i) 3-methylpent-2-ene

Ignore E-Z, commas, spaces or missing hyphens

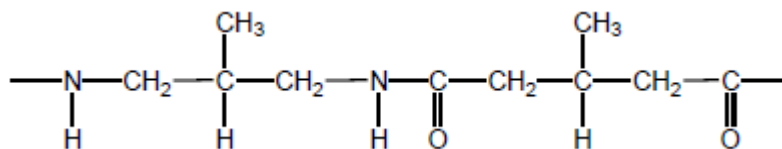
1

(ii) 4-amino-3-methylbutanoic acid

Ignore commas, spaces or missing hyphens

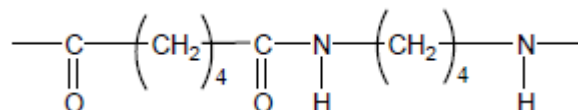
1

(iii)



or any polyamide section containing

8 carbons plus two C=O plus two N-H, such as



Trailing bonds are required

1

- (iv) Non polar OR no polar groups / bonds (for attack by water / acids / alkalis / nucleophiles or for hydrolysis)

C-C bonds are strong

1

[7]

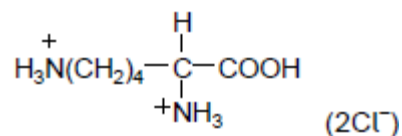
8

- (a) 2,6-diaminohexanoic acid

Ignore additional , or – or spaces.

1

- (b) (i)



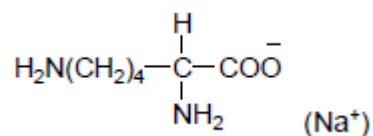
NB both N must be protonated.

Allow $-\text{NH}_3^+$ allow CO_2H Allow $-\text{H}_3\text{N}$.

Penalise $-\text{C}_4\text{H}_8-$ here.

1

- (ii)



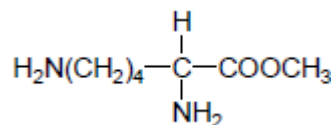
Allow CO_2^- .

Allow $-\text{H}_2\text{N}$.

Allow $-\text{COONa}$ but penalise $\text{O}-\text{Na}$ bond shown.

1

- (iii)

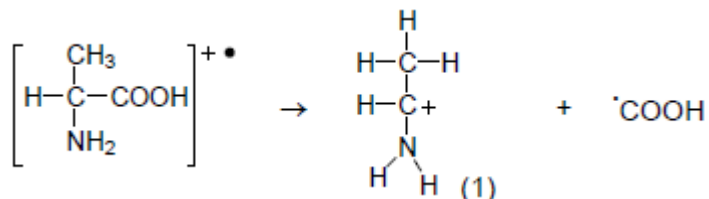


Allow CO_2CH_3 .

Allow $-\text{NH}_3^+$ or $-\text{H}_2\text{N}$.

1

(c)



1 for displayed formula of fragment ion.

1 for molecular ion of alanine AND radical.

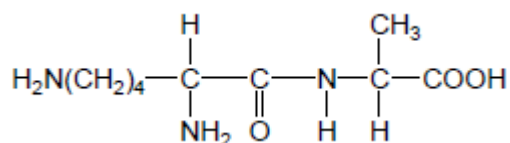
Allow molecular ion without brackets and fragment ion in brackets with outside +.

Allow dot anywhere on radical.

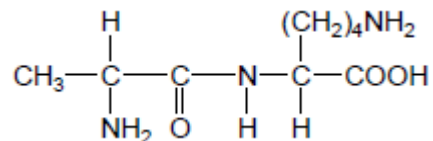
Allow $[\text{C}_3\text{H}_7\text{NO}_2]^+ \cdot$ for molecular ion.

2

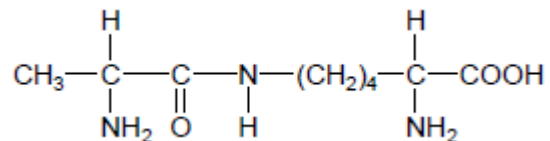
(d)



OR



OR



Dipeptide, not repeating unit /.

Allow CO_2H Allow $-\text{H}_2\text{N}$.

Allow $-\text{CONH}-$.

1

(e) M1 In acid lysine has double positive or more positive charge

1

M2 (Lysine ion) has greater affinity / greater attraction / adheres better / sticks better to polar / stationary phase

M2 only scores after a correct M1.

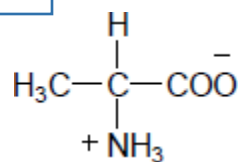
Ignore greater retention time.

1

[9]

9

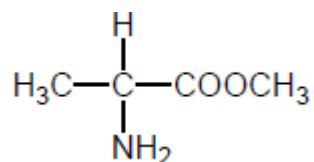
(a)



Allow $-\text{NH}_3^+$ and $+\text{NH}_3^-$

1

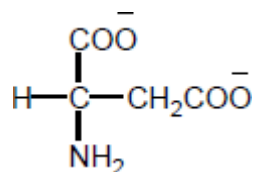
(b)



Allow protonated form, i.e. $-\text{NH}_3^+$ or $+\text{NH}_3^-$

1

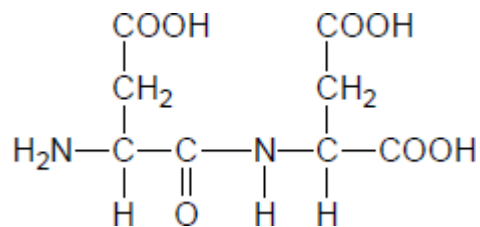
(c)



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

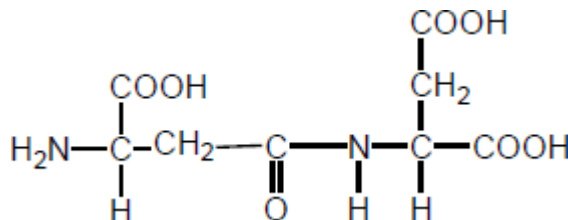
1

(d)



Allow zwitterion with any COO^-

Allow use of "wrong" COOH

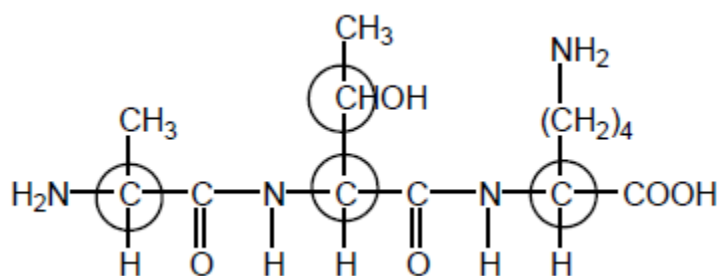


1

[4]

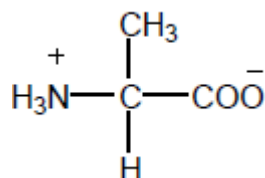
10

(a) (i)

*These four only*

1

(ii)

*Allow -NH₃⁺ and ⁺NH₃-*

1

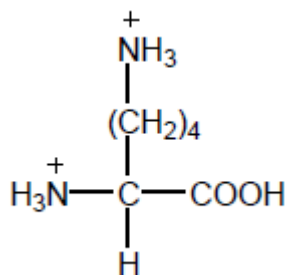
(iii) 2-amino-3-hydroxybutanoic acid*Ignore 1 in butan-1-oic acid*

Do not penalise commas or missing hyphens

Penalise other numbers

1

(iv)

*Allow -NH₃⁺ and ⁺NH₃-*

1

(b) (i) Condensation

Allow polyester

1

(ii) propane-1,3-diol*Must have e**Allow 1,3-propanediol*

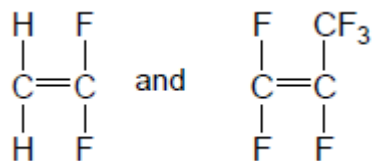
1

(c) (i) Addition

Not additional

1

(ii)

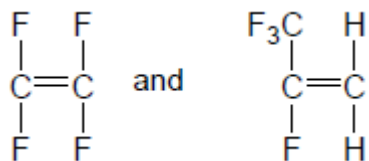


Allow monomers drawn either way round

Allow bond to F in CF₃

1

OR



1 for each structure within each pair

1

(d) c

If wrong, CE = 0

1

C-C or C-F bonds too strong

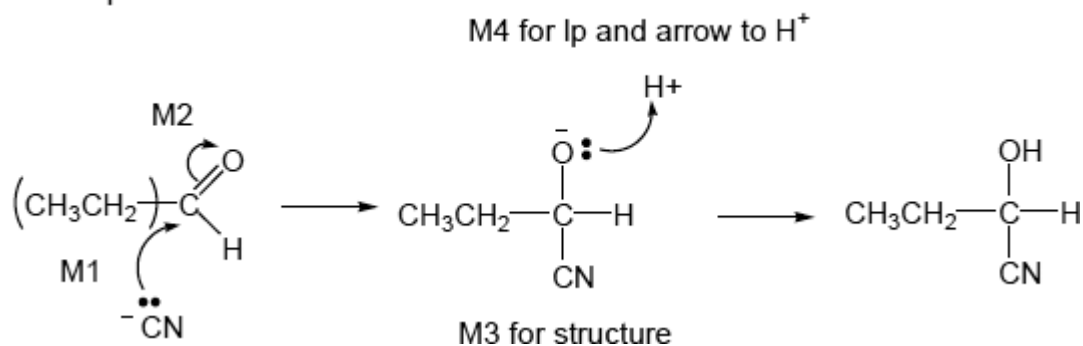
1

[11]

11

(a)

nucleophilic addition



- allow :CN⁻
- M2 not allowed independent of M1, but
- allow M1 for correct attack on C⁺
- + rather than δ+ on C=O loses M2
- M3 is for correct structure including minus sign but lone pair is part of M4
- Allow C₂H₅
- M1 and M4 for lp and curly arrow

1

(b) 2-bromobutanenitrile*Allow 2-bromobutane-1-nitrile*

1

(c) **M1** ammonia or NH₃*Ignore temp or pressure*

1

M2 excess (ammonia) excess tied to NH₃ and may score in M1 unless contradicted

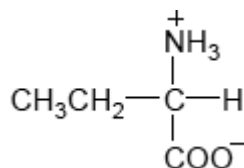
Ignore concentrated or sealed container, Acid loses conditions mark

1

M3 nucleophilic substitution*Allow close spelling*

1

(d) (i)

*Allow C₂H₅**Allow -CO₂⁻**Allow +NH₃⁻**Don't penalize position of + on NH₃*

1

- (ii) **M1** electrostatic forces between ions in **X** **QOL**
Allow ionic bonding.

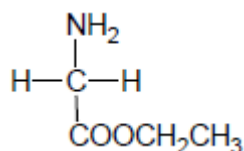
1

Marks independent

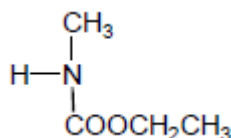
- M2** (stronger than) hydrogen bonding between $\text{CH}_3\text{CH}_2\text{CH}(\text{OH})\text{COOH}$
CE mention of molecules of **X** or inter molecular forces between **X**
 loses both marks

1

- (e) (i)



OR

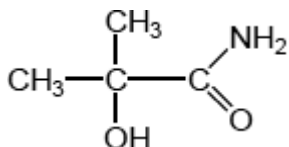


Isomer of $\text{C}_4\text{H}_9\text{NO}_2$

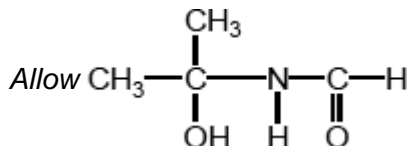
Allow NH_2-

1

- (ii)



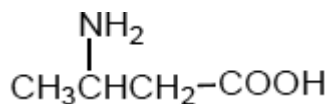
Isomer of $\text{C}_4\text{H}_9\text{NO}_2$ allow NH_2-



1

- (iii) $\text{H}_2\text{N}-\text{CH}_2\text{CH}_2\text{CH}_2-\text{COOH}$ or $\text{H}_2\text{N}-(\text{CH}_2)_3-\text{COOH}$
Isomer of $\text{C}_4\text{H}_9\text{NO}_2$ allow NH_2-

OR

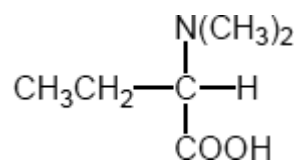


*Do **not** allow $-\text{C}_3\text{H}_6-$*

*Beware – do not credit **X** itself*

1

(f)



Answer has 6 carbons so **NOT** isomer of **X**

Allow C_2H_5

Must have bond from C to N not to methyl group

1

[16]