

GCE MARKING SCHEME

CHEMISTRY AS/Advanced

SUMMER 2012

(a) Any valid ester structure with formula $C_{10}H_{12}O_2$

Examples:

(b) (i) Compound X [1]

(iii) Rotate the plane of polarised light in opposite directions [1]

(c)

Reagent(s)	Observation if the test is positive	Compound(s) that would give a positive result
I ₂ / NaOH (aq)	Yellow solid	X
Na ₂ CO ₃ (aq)	Bubbles of colourless gas / effervescence	W
FeCl ₃ (aq)	Dark purple/blue/green - do not accept 'precipitate'	X, Z

(1 mark for each box) [6]

[1]

(d) (i) Heat / Alkaline / Potassium manganate(VII) / then acidify
(1 mark for Potassium manganate + 1 other point; 2 marks for all)
[2]

(ii) I. Addition polymer – One large molecule formed only / Condensation polymer – one large molecule with small molecules (e.g. water) lost. (1)

Addition polymer – one starting material / Condensation – two starting materials

OR Addition polymer – one functional group in each molecule/ Condensation polymer –
two functional groups in each molecule

(1) [2]

II.

$$-\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{CH}_{2}-\mathsf{CH}_{2}}{\overset{\mathsf{O}}{\longleftarrow}}\underset{\mathsf{C}}{\overset{\mathsf{O}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}}\underset{\mathsf{C}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}{\overset{\mathsf{O}}\underset{\mathsf{C}}}$$

(e) (i) NaBH₄ / LiAlH₄ or name(1) Reduction (1) [2]
- ignore conditions unless LiACH₄ - do not accept 'redox'
in water
(ii)

$$O = C \\ O =$$

Accept structures with only one –OH group reacted. [1]

[19 marks]

(a) (i) Alanine forms a zwitterion (1)
 Forces between alanine molecules are ionic bonding (1)
 Ionic bonding much stronger than hydrogen bonding / van der Waals (1)

Max 2 marks [2]

(ii) 1 mark for each correct structure

[2]

(iii) 1 mark for correct identification of peptide link

[1]

(b) Enzymes / Structural proteins / Hormones or specific example

[1]

(c) 1 mark for arrows in first stage; 1 mark for correct intermediate; 1 mark for arrow giving gain of proton in second stage (from HCN or from H⁺).

$$R \xrightarrow{C} C \xrightarrow{D} R \xrightarrow{D} R \xrightarrow{D} C \xrightarrow{D} R \xrightarrow{D} R \xrightarrow{D} R \xrightarrow{D} R \xrightarrow{D} C \xrightarrow{D} R \xrightarrow{D}$$

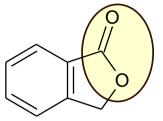
[3]

(d) Soda lime

[1]

[10 marks]

(a) (i) [1]



Phthalide

(ii) [1]

- (b) Distillation / Chromatography [1]
- (c) Hydrogenation of 3-butyl phthalide removes a benzene ring (1)

 Benzene ring is more stable than alkene/ Reference to delocalisation energy (1) [2]
- (d) 62.1% [1]
- (e) (i) Greater variety of different phthalides that can be produced [1]
 - (ii) Higher atom economy / less waste / carbon monoxide is toxic [1]
 - do not accept references to yield
- (f) Silver nitrate and ammonia / Tollen's reagent (1); Q = Silver mirror (1); R = No reaction (1)

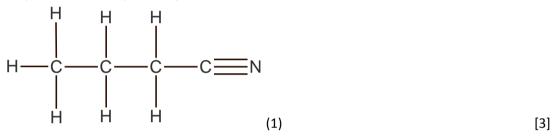
OR 2,4,-DNP (1); Orange precipitate with Q (1); No reaction with R (1)

OR Fehling's solution (1); Orange solid with Q (1); No reaction with R (1) [3]

[11 marks]

- (a) (i) Nucleophilic substitution / Hydrolysis [1]
 - (ii) Dissolved in alcohol (1) Propene or unambiguous structure (1) [2]
 - (iii) Potassium manganate(VII) / Potassium dichromate(VI) must be **name** (1)
 Oxidation (1) [2]
 - (iv) (Add Potassium dichromate(VI)) and distil off the propanal from the reaction mixture [1]
- (b) (i) Step 1: Potassium cyanide in ethanol / Heat (1)

Step 2: Heat with aqueous hydrochloric acid (or other acid) (1)



- (ii) Two points from different bullet points 1 mark each.
- Atom economy / Amount of waste / Whether waste material was recyclable / Whether waste was toxic.
- Amount of energy required / temperature required / pressure required / conditions used
- Rate of production / time
- Availability of catalyst
- Cost of reactants / Availability of reactants / toxicity of reactants.
- Two step processes usually have lower yields than one step processes / percentage
 yield

[2]

- Purification method / separation
- (c) (i) Butanoic acid is $C_4H_8O_2$ so $M_r = 88$ (1)

Percentage carbon = 48/88 x 100 = 54.5%; percentage hydrogen = 8/88 = 9.1%;

Percentage oxygen = 32/88 = 36.4% (At least two of these for 1)

OR empirical formula for butanoic acid = C_2H_4O (1) and

calculate empirical formula from percentage masses = C_2H_4O (1)

- (ii) Structure 1 mark + 4 marks for explanations.
 - Product is ethyl ethanoate. (1)
 - Two points from the following required for each mark– MAX 4 marks
 - Sweet-smelling = ester
 - Peak at 1.0ppm implies CH₃
 - Peak area 3 = CH₃
 - Peak area 2 = CH₂
 - o Triplet shows CH₃ is next to a CH₂ group.
 - o Singlet shows CH₃ no hydrogen atoms bonded to adjacent carbon.
 - Peak at 2.1 ppm suggests this is next to C=O.
 - Quartet shows CH₂ is adjacent to a CH₃ group.
 - o Peak at 4.0 ppm shows it is -O-CH₂-
 - o IR Peak at 1752 cm⁻¹ = C=O
 - IR Peak at 2981 cm⁻¹ = C-H or O-H
 - o Cannot be –OH as we know there is no –OH in NMR spectrum

[5]

QWC: selection of a form and style of writing appropriate to purpose and to complexity of subject matter. (1)

QWC: organisation of information clearly and coherently; use of specialist vocabulary where appropriate. (1) [2]

[20 marks]

(a) (i) (Concentrated) nitric acid / (concentrated) sulfuric acid / Temperature of 40-80°C (Any 2 = 1 mark; All 3 = 2 marks)

Electrophilic substitution (1)

[3]

(ii) I. Peak area is proportional to amount of substance (1)

Percentage =
$$(30 / 38) \times 100 = 79\%$$
 (1

(Can obtain both marks from correct percentage)

[2]

II. $45 = COOH^{+}$, $46 = NO_{2}^{+}$, $122 = C_{6}H_{4}NO_{2}^{+}$ and $167 = C_{7}H_{5}NO_{4}^{+}$.

$$(Any 2 = 1 mark; All 4 = 2 marks)$$

[2]

(iii) I. Lower melting point / melts over a range [1]

II. 1 mark for each point.

- Dissolve in the minimum volume
- Of hot water
- Filter hot
- Allow to cool
- Filter
- Dry residue under suction / in oven below 142°C

Max 4 marks [4]

QWC: legibility of text, accuracy of spelling, punctuation and grammar, clarity of meaning.[1]

(b) (i) Tin and concentrated hydrochloric acid

[1]

(ii) Below 10°C (1)

(1) [2]

(iii) N=N double bond is chromophore (1)

Compound absorbs blue /green / complementary colours to red / all colours but red (1)

Remaining frequencies are transmitted, giving the red colour seen. (1)

Any 2 out of 3 [2]

(c) Nitrogen has a lone pair (1) which can accept a proton (1)

[2]

[20 marks]