# wеес cbac 

## GCE A LEVEL MARKING SCHEME

SUMMER 2018

A LEVEL (NEW) CHEMISTRY - UNIT 4 1410U40-1

## INTRODUCTION

This marking scheme was used by WJEC for the 2018 examination. It was finalised after detailed discussion at examiners' conferences by all the examiners involved in the assessment. The conference was held shortly after the paper was taken so that reference could be made to the full range of candidates' responses, with photocopied scripts forming the basis of discussion. The aim of the conference was to ensure that the marking scheme was interpreted and applied in the same way by all examiners.

It is hoped that this information will be of assistance to centres but it is recognised at the same time that, without the benefit of participation in the examiners' conference, teachers may have different views on certain matters of detail or interpretation.

WJEC regrets that it cannot enter into any discussion or correspondence about this marking scheme.

## UNIT 4: ORGANIC CHEMISTRY AND ANALYSIS

## MARK SCHEME

## GENERAL INSTRUCTIONS

## Recording of marks

Examiners must mark in red ink.
One tick must equate to one mark, apart from extended response questions where a level of response mark scheme is applied.
Question totals should be written in the box at the end of the question.
Question totals should be entered onto the grid on the front cover and these should be added to give the script total for each candidate.
Extended response questions
A level of response mark scheme is applied. The complete response should be read in order to establish the most appropriate band. Award the higher mark if there is a good match with content and communication criteria. Award the lower mark if either content or communication barely meets the criteria.

## Marking rules

All work should be seen to have been marked.
Marking schemes will indicate when explicit working is deemed to be a necessary part of a correct answer.
Crossed out responses not replaced should be marked.

## Marking abbreviations

The following may be used in marking schemes or in the marking of scripts to indicate reasons for the marks awarded.
cao = correct answer only
ecf = error carried forward
bod $=$ benefit of doubt
Credit should be awarded for correct and relevant alternative responses which are not recorded in the mark scheme.

## Section A

| Question |  |  | Marking details | Marks available |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  | AO1 | AO2 | AO3 | Total | Maths | Prac |
| 1. | (a) |  |  |  |  | 1 |  | 1 |  |  |
|  | (b) |  | ${ }^{-} \mathrm{CN}$ accept $\mathrm{CN}^{-}$ | 1 |  |  | 1 |  |  |
| 2. |  |  | butanoic acid accept methylpropanoic acid / pentanedioic acid |  | 1 |  | 1 |  |  |
| 3. |  |  | $2 \mathrm{~mol} \mathrm{NH}_{3}$ from 1 mol amide $\mathbf{T}$ 0.060 mol from 0.030 mol amide $\mathbf{T}$ $M_{\mathrm{r}}$ of the amide 3.90/0.030 $=130$ $M_{r}{ }^{\prime} R^{\prime}=130-88=42$ <br> therefore $\mathrm{C}_{3} \mathrm{H}_{6} / \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2}$ |  | 1 | 1 | 2 | 1 |  |


| Question |  |  | Marking details | Marks available |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  | AO1 | AO2 | AO3 | Total | Maths | Prac |
| 4. | (a) |  |  |  |  | 1 |  | 1 |  |  |
|  | (b) |  | 2-aminopropanoic acid exists as zwitterions / $\mathrm{CH}_{3} \mathrm{CHNH}_{3}{ }^{+} \mathrm{COO}^{-}$ <br> it has strong ionic character therefore much stronger forces between molecules than the other acids (1) |  | 2 |  | 2 |  |  |
| 5. |  |  | $275.4 \mathrm{dm}^{3}$ gaseous material from 222 g RDX therefore $1 \mathrm{~m}^{3}$ from $222 \times 1000 / 275.4=806$ answer must be given to 3 sig figs (1) |  | 1 | 1 | 2 | 2 |  |
|  |  |  | Section A total | 1 | 7 | 2 | 10 | 3 | 0 |

Section B

| Question |  |  | Marking details | Marks available |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  | A01 | AO2 | AO3 | Total | Maths | Prac |
|  | (a) | (i) |  | $\mathrm{C}_{8} \mathrm{H}_{16} \mathrm{O}$ | 1 |  |  | 1 |  |  |
|  |  | (ii) | $\begin{aligned} & \mathrm{NaOH}(\mathrm{aq}) /{ }^{-} \mathrm{OH}(\mathrm{aq}) \\ & \text { nucleophilic substitution (1) } \end{aligned}$ | 2 |  |  | 2 |  | 1 |
|  |  | (iii) | accept any of following acidified dichromate(VI) acidified potassium dichromate(VI) $\mathrm{H}^{+}, \mathrm{Cr}_{2} \mathrm{O}_{7}^{2-}$ acidified manganate(VII) acidified potassium manganate(VII) acidified permanganate $\mathrm{H}^{+}, \mathrm{MnO}_{4}^{-}$ | 1 |  |  | 1 |  | 1 |
|  | (b) | (i) | $\mathrm{CHI}_{3}$ | 1 |  |  | 1 |  | 1 |
|  |  | (ii) | possible formulae <br> accept an unsaturated methyl secondary alcohol $\begin{equation*} \text { e.g. } \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}=\mathrm{CHCH}_{2} \mathrm{CH}(\mathrm{OH}) \mathrm{CH}_{3} \tag{1} \end{equation*}$ <br> methylketone / $\mathrm{CH}_{3} \mathrm{C}=\mathrm{O}$ (1) <br> or $\mathrm{CH}(\mathrm{OH}) \mathrm{CH}_{3}$ (that can be oxidised to $\mathrm{CH}_{3} \mathrm{C}=\mathrm{O}$ ) |  | 1 | 1 | 2 |  |  |



| Question |  |  | Marking details | Marks available |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  | A01 | AO2 | AO3 | Total | Maths | Prac |
| 7. | (a) | (i) |  | both curly arrows needed (1) correct structure of intermediate | 1 | 1 |  | 2 |  |  |
|  |  | (ii) | electrophilic substitution | 1 |  |  | 1 |  |  |
|  | (b) |  | Ione pair of electrons on the nitrogen atom(s) | 1 |  |  | 1 |  |  |
|  | (c) | (i) | orange / red solid | 1 |  |  | 1 |  |  |
|  |  | (ii) | the derivatives formed using 2,4-DNPH have suitable melting temperatures for identification / precise melting temperatures |  |  | 1 | 1 |  | 1 |
|  | (d) |  | dissolve the (impure) oxime in a minimum volume (1) hot methanol / warm methanol (1) <br> water bath / electrical heater (1) <br> allow to cool (1) <br> filter off oxime and dry (1) | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | 1 | 5 |  | 5 |


| Question |  | Marking details | Marks available |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  | A01 | AO2 | AO3 | Total | Maths | Prac |
| (e) |  |  | (broad) peak at $3200-3550 \mathrm{~cm}^{-1}$ due to $\mathrm{O}-\mathrm{H}$ bond present in the oxime and not in the amide (1) <br> peak at $1650-1750 \mathrm{~cm}^{-1}$ due $\mathrm{C}=\mathrm{O}$ bond present in the amide and not in the oxime (1) <br> peak at $3300-3500 \mathrm{~cm}^{-1}$ due to $\mathrm{N}-\mathrm{H}$ bond present in the amide and not in the oxime (1) |  | 3 |  | 3 |  |  |
|  |  | Question 7 total | 6 | 6 | 2 | 14 | 0 | 6 |


| Question |  |  | Marking details | Marks available |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  | A01 | AO2 | AO3 | Total | Maths | Prac |
| 8. | (a) |  |  | $\mathrm{F}_{3} \mathrm{CCl} \rightarrow \mathrm{F}_{3} \mathrm{C} \cdot+\mathrm{Cl} \cdot$ |  | 1 |  | 1 |  |  |
|  | (b) | (i) | the higher the ratio of $\mathrm{HF}: \mathrm{CH}_{2} \mathrm{Cl}_{2}$ the higher the yield of $\mathrm{CH}_{2} \mathrm{~F}_{2}$ |  |  | 1 | 1 |  |  |
|  |  | (ii) | $\begin{equation*} \mathrm{n}\left(\mathrm{CH}_{2} \mathrm{~F}_{2}\right)=0.0356 \tag{1} \end{equation*}$ $\begin{equation*} \text { mass of } \mathrm{CH}_{2} \mathrm{~F}_{2}=0.0356 \times 52.02=1.85 \tag{1} \end{equation*}$ <br> ecf possible from incorrect $M_{r}$ |  | 2 |  | 2 | 1 |  |
|  | (c) | (i) | high pressure because there are more (gaseous) moles on the left than on the right | 1 |  |  | 1 |  |  |
|  |  | (ii) | high temperature as the endothermic reaction needs heat for the position of equilibrium to move to the right | 1 |  |  | 1 |  |  |
|  |  | (iii) | separation can be achieved by (fractional) distillation (1) <br> reduce the temperature of the mixture to below $-30^{\circ} \mathrm{C}$ and let the mixture warm up slowly ( $1,1,1,2$-tetrafluoroethane will distil off first and can be condensed in a cold trap) (1) |  |  | 2 | 2 |  | 2 |


| Question |  |  | Marking details | Marks available |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  | A01 | AO2 | AO3 | Total | Maths | Prac |
| (d) | (i) |  |  | each end of the $\mathrm{C}=\mathrm{C}$ double bond has two different groups bonded to it | 1 |  |  | 1 |  |  |
|  | (ii) | 1 | electrophilic addition | 1 |  |  | 1 |  |  |
|  |  | II | (intermediate) carbocations / carbonium ions have similar stabilities / similar reactivities / similar activation energies / are formed at similar rates |  |  | 1 | 1 |  |  |
|  | (iii) |  |   <br> correct formulae (1) <br> chiral centres correctly shown on both compounds <br> award (1) for one correct formula and chiral centre |  | 2 |  | 2 |  |  |
|  | (iv) |  |   |  | 1 |  | 1 |  |  |
|  |  |  | Question 8 total | 4 | 6 | 4 | 14 | 1 | 2 |


| Question |  |  |  | Marking details | Marks available |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  |  |  | A01 | AO2 | AO3 | Total | Maths | Prac |
| 9. | (a) |  |  | Indicative content <br> - percentage of chlorine in this compound is $32.1 \%$ which is consistent with that of 2,4-D <br> - colourless gas with $\mathrm{NaHCO}_{3}$ indicates compound is a carboxylic acid <br> - $\mathrm{n}(\mathrm{NaOH})$ is $1.85 \times 10^{-3} \mathrm{~mol}$ <br> - mole ratio is $1: 1$ therefore only one COOH group <br> - number of moles of compound in $25.0 \mathrm{~cm}^{3}$ is also $1.85 \times 10^{-3} \mathrm{~mol}$ therefore $1.85 \times 10^{-2}$ in $250 \mathrm{~cm}^{3}$ <br> - $M_{r}$ of compound is 221 which is consistent with that of 2,4-D <br> test results consistent with structure of 2,4-D <br> - no chloride ions on reflux with NaOH therefore chlorine atoms are bonded directly to the ring, not in alkyl side-chains <br> - no white precipitate with aqueous bromine therefore not a phenol <br> - aqueous bromine not decolourised therefore no $\mathrm{C}=\mathrm{C}$ double bonds <br> spectral data consistent with that of 2,4-D <br> - ${ }^{1} \mathrm{H}$ NMR spectrum of this compound would show three peaks; aromatic (area of 3), $-\mathrm{CH}_{2}-$ (area 2), O-H (area 1) <br> - ${ }^{13} \mathrm{C}$ NMR of this compound would show eight peaks as it has 8 different carbon environments <br> - take its melting temperature and compare to a book value <br> - mix a sample with some actual 2,4-D, and take its melting temperature - see if the value is unchanged | 2 | 2 | 2 | 6 | 2 | 4 |






UNIT 4: ORGANIC CHEMISTRY AND ANALYSIS
SUMMARY OF MARKS ALLOCATED TO ASSESSMENT OBJECTIVES

| Question | AO1 | AO2 | AO3 | Total | Maths | Prac |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| Section A | 1 | 7 | 2 | 10 | 3 | 0 |
| 6. | 6 | 4 | 4 | 14 | 0 | 6 |
| 7. | 6 | 6 | 2 | 14 | 0 | 6 |
| 8. | 4 | 6 | 4 | 14 | 1 | 2 |
| 9. | 3 | 6 | 6 | 14 | 4 | 1 |
| 10. | 22 | 37 |  |  | 80 | 10 |
| Totals |  |  |  |  |  |  |

