Oxford Cambridge and RSA

## GCE

## Chemistry B (Salters)

Unit H033/02: Chemistry in depth
Advanced Subsidiary GCE

Mark Scheme for June 2016

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This mark scheme is published as an aid to teachers and students, to indicate the requirements of the examination. It shows the basis on which marks were awarded by examiners. It does not indicate the details of the discussions which took place at an examiners' meeting before marking commenced.

All examiners are instructed that alternative correct answers and unexpected approaches in candidates' scripts must be given marks that fairly reflect the relevant knowledge and skills demonstrated.

Mark schemes should be read in conjunction with the published question papers and the report on the examination.

OCR will not enter into any discussion or correspondence in connection with this mark scheme.

Annotations

| Annotation | Meaning |
| :--- | :--- |
| A | Correct response |
| A | Incorrect response |
| BOD | Omission mark |
| CON | Benefit of doubt given |
| RE | Contradiction |
| SF | Rounding error |
| ECF | Error in number of significant figures |
| L1 | Level 1 |
| L2 | Level 2 |
| L3 | Level 3 |
| NBOD | Benefit of doubt not given |
| SEEN | Ignore |
| I |  |

Abbreviations, annotations and conventions

| Annotation | Meaning |
| :---: | :--- |
|  | alternative and acceptable answers for the same marking point |
| DO NOT ALLOW | Separates marking points |
| IGNORE | Statements which are irrelevant |
| ALLOW | Answers that can be accepted |
| ( ) | Words which are not essential to gain credit |
| ECF | Underlined words must be present in answer to score a mark |
| AW | Or reverse argument forward |
| ORA |  |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (a) |  | $\mathrm{C}_{4} \mathrm{H}_{10}(\mathrm{~g} / \mathrm{l})+61 / 2 \mathrm{O}_{2}(\mathrm{~g}) \rightarrow 4 \mathrm{CO}_{2}(\mathrm{~g})+5 \mathrm{H}_{2} \mathrm{O}(\mathrm{I})$ <br> $\checkmark$ for balanced equation <br> $\checkmark$ for state symbols | 2 | NOT multiples |
|  | (b) |  | FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer $=\Delta_{\mathrm{c}} \mathrm{H}=\mathbf{- 1 4 1 9}\left(\mathrm{kJ} \mathrm{mol}^{-1}\right)$ award 3 marks <br> Calculates energy transferred to water $\mathrm{q}=50.00 \times 4.18 \times(74-19)=11495(\mathrm{~J}) \checkmark$ Calculates number of moles of butane burned $=0.47 / 58.0$ or $0.008(1) \mathrm{mol}$ $\left(\Delta_{\mathrm{c}} \mathrm{H}=-[11495 / 0.008(1)]=-1419136 \mathrm{~J} \mathrm{~mol}^{-1}\right)$ $\Delta_{c} H=-1419\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ | 3 | Correct answer - $1419 \pm 1 \mathrm{~kJ} \mathrm{~mol}^{-1}$ scores 3 marks <br> ALLOW ECF between steps <br> ALLOW final answer to 2 or more sf (eg -1440 $\mathrm{kJ} \mathrm{mol}^{-1}$ if early rounding is evident) <br> Final MP must include negative sign |
|  | (c) | (i) | -2850--2950 (kJ mol$\left.{ }^{-1}\right) \checkmark$ | 1 | Must have negative sign |
|  |  | (ii) | ```Any two from: \(\checkmark \checkmark\) loss of fuel by evaporation / escape of unburned butane evaporation of water incomplete combustion / reaction non-standard conditions /states heat used to raise temp of calorimeter``` | 2 | Answers can be in any order <br> Ignore 'not fully reacted' as this makes it unclear whether the candidate is talking about the vol of butane or the combustion reaction Ignore measurement errors |
|  |  | (iii) | One from: <br> use a (draught) shield because this will reduce heat lost (to the surroundings); burn the butane in oxygen / because this will ensure that the combustion is more complete; use bomb calorimeter ensures complete combustion / reduces heat loss; use cover over lighter during weighing to prevent evaporation; insulate can to reduce heat loss | 1 | Must have method plus explanation to score <br> Ignore changes to vol of water / mass of fuel / length of time for combustion / move flame nearer to can / <br> ALLOW 'put lid on can' NOT 'use a polystyrene cup' |


| Ques | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: |
| (d) | Any one from: <br> reacting carbon and hydrogen doesn't (necessarily) make butane <br> Carbon and Hydrogen do not react together (under standard conditions) $\checkmark$ | 1 | ALLOW reference to formation of a mixture of products or alternative product(s) / side reactions |
| (e) | skeletal formula <br> systematic name: (2-)methylpropane $\checkmark$ | 1 | IGNORE dashes, commas and spaces in the name Needs BOTH skeletal formula AND name |
| (f) | ```FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer \(=(+) 486.6\) / (+)487 (kJ mol-1) award 3 marks energy absorbed in breaking bonds \(=3(413)+(358)+x+11 / 2(498)\) \(=2344+\mathrm{x}(\mathrm{kJ})\) AND energy evolved in making bonds \(=2(805)+4(x)\) \(=1610+4 \mathrm{x}(\mathrm{kJ}) \checkmark\) Overall energy change (Bonds broken - bonds made \(=\Delta \mathrm{H}\) ) \(=[2344+x]-[1610+4 x]=-726 \mathrm{~kJ} \mathrm{~mol}-1\) OR \(2344-1610+726=3 x\) \(1460=3 x \checkmark\) \(x=(+) 486.6 /(+) 487(\mathrm{~kJ} \mathrm{~mol}-1) ~ \checkmark\)``` | 3 | Correct answer $\mathbf{+ 4 8 6 . 6 / + 4 8 7 \mathrm { kJ } \mathrm { mol } ^ { - 1 }}$ scores 3 marks <br> ALLOW ECF between steps <br> 2344 and 1610 in calculation scores 1 mark if no other mark scored <br> ALLOW OH for x in calculation <br> ALLOW -sign if evaluation of their expression for x is correct |
|  | Total | 14 |  |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 2 | (a) |  | 8-11 | 1 | Accept any value between 7.1-14 |
|  | (b) |  | FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer $=\mathbf{0 . 3 4 ( m o l ) ( 2 ~ s f ) ~ a w a r d ~} \mathbf{3}$ marks mass of $\mathrm{Mg}(\mathrm{OH})_{2}$ in $250 \mathrm{~cm}^{3}=8 / 100 \times 250(=20) \mathrm{g} \checkmark$ $\mathrm{M}_{\mathrm{r}} \mathrm{Mg}(\mathrm{OH})_{2}=58.3 \mathrm{~g} \mathrm{~mol}^{-1}$ <br> Moles $\mathrm{Mg}(\mathrm{OH})_{2}=(20 / 58.3)=0.34(\mathrm{~mol})(2 \mathrm{sf}) \checkmark$ | 3 | Allow ecf throughout <br> Final answer MUST be to 2 sf |
|  | (c) | (i) | (it is the) oxidation state/number of the sulfur ${ }^{\checkmark}$ | 1 | incorrect number is CON ALLOW 6/+6/6+ <br> ALLOW oxidisation/ |
|  | (c) | (ii) | $\mathrm{Mg}(\mathrm{OH})_{2}$ is not (completely)soluble / forms a suspension (in water) $\sqrt{ }$ | 1 | ALLOW cloudiness of suspension obscures colour of indicator/ makes it difficult to identify end-point ALLOW medicine for $\mathrm{Mg}(\mathrm{OH})_{2}$ |
|  | (d) | (i) | (it is the) mean/average of the concordant titres / repeats 1 and 3 | 1 | ALLOW only used titres agreeing to within $0.1 \mathrm{~cm}^{3}$ / repeat 2 not included as it is an anomalous result(outlier) |
|  | (d) | (ii) | FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer $=0.0166$ ( mol ) award 2 marks $\begin{aligned} & \text { amount of } \mathrm{NaOH}(\text { in titre })=(16.65 / 1000 \times 1.99) \\ & =0.0331 \mathrm{~mol} \checkmark \end{aligned}$ <br> amount of $\mathrm{H}_{2} \mathrm{SO}_{4}$ in excess $=(0.5 \times 0.0331)$ $=0.0166(\mathrm{~mol}) \checkmark$ | 2 | Allow ecf from incorrect titre used in calculation <br> Do not accept 0.0165 , incorrect rounding of 0.01655 Final answer to 2sf or more. |


| Question |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| (d) | (iii) | FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer $=19.20-19.55(\mathrm{~g})$ award 3 marks amount of $\mathrm{H}_{2} \mathrm{SO}_{4}$ initially $=(25.0 / 1000 \times 2.00)=0.05 \mathrm{~mol}$ amount of $\mathrm{H}_{2} \mathrm{SO}_{4}$ used up = number of moles $\mathrm{Mg}(\mathrm{OH})_{2}$ $=(0.05-0.0166)=0.0334 \mathrm{~mol}$ $\begin{aligned} & \text { mass of } \mathrm{Mg}(\mathrm{OH})_{2} \text { in } 250 \mathrm{~cm}^{3}=(250 / 25 \times 0.0334 \times 58.3) \\ & =19.47 / 19.5(\mathrm{~g}) \downarrow \end{aligned}$ | 3 | ALLOW ecf from incorrect value in d(ii) <br> ALLOW ecf throughout <br> ALLOW 3 or more sf throughout, but rounding must be correct |
| (d) | (iv) | $(0.06 / 25.0 \times 100)=0.2(\%) \checkmark$ | 1 | Correct answer without working scores the mark |
| (e) | (i) | $\mathrm{Mg} \rightarrow \mathrm{Mg}^{2+}+2 \mathrm{e}^{-} / \mathrm{Mg}-2 \mathrm{e}^{-} \rightarrow \mathrm{Mg}^{2+} \checkmark$ | 1 | Ignore state symbols |
| (e) | (ii) | (magnesium) loses (two) electrons $\checkmark$ | 1 | ACCEPT the oxidation state (of the magnesium) increases (from 0 to +2 ) ignore species losing electrons unless incorrectly named |
| (e) | (iii) | $\mathrm{H}^{+} /$hydrogen (ion) $\checkmark$ | 1 | ACCEPT hydrochloric acid/ HCl NOT H/H2 |
| (f) | (i) | $\mathrm{Mg}^{2+}(\mathrm{aq})+2 \mathrm{OH}^{-}(\mathrm{aq}) \rightarrow \mathrm{Mg}(\mathrm{OH})_{2}(\mathrm{~s})$ <br> $\checkmark$ for balanced ionic equation <br> $\checkmark$ for state symbols | 2 | DO NOT ALLOW spectator ions <br> ALLOW state symbol mark if 'magnesium hydroxide' given as solid and all other species as aq |
|  | (ii) | $\mathrm{Ba}(\mathrm{OH})_{2}$ is more soluble (in water) $\checkmark$ | 1 | ORA <br> ALLOW $\mathrm{Ba}(\mathrm{OH})_{2}$ will not precipitate |
|  |  | Total | 19 |  |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (a) | (i) | $\mathrm{CH}_{3} \mathrm{COOH} \checkmark$ | 1 | ALLOW any unambiguous structure |
|  |  | (ii) | ester $\checkmark$ | 1 |  |
|  |  | (iii) | FIRST CHECK THE ANSWER ON THE ANSWER LINE <br> If answer = $\mathbf{7 5}$ (\%) award 2 marks <br> Mr of aspirin $\begin{aligned} & =[(12.0 \times 9)+(16.0 \times 4)+(1.0 \times 8)]=180.0 \\ & \% \text { atom economy } \\ & =180.0 /(138.0+102.0) \times 100 \\ & \text { OR } 180 /(180+60) \times 100 \\ & =75(\%) \quad \end{aligned}$ | 2 | ALLOW ecf from incorrect value for Mr of aspirin |
|  | (b) |  | 2-hydroxybenzoic acid $\checkmark$ | 1 | ALLOW salicylic acid / phenol group (present) IGNORE dashes, commas and spaces NOT phenol on its own |
|  | (c) |  | dissolve (crude) aspirin/solid in hot/warm ethanol/solvent use the minimum volume/amount of ethanol/solvent $\checkmark$ (allow to) cool/crystallise <br> filter, wash (with cold ethanol) and (allow to) dry $\checkmark$ | 4 | 'Dissolve in a minimum amount of hot ethanol' scores 2 marks (MP1 and MP2) <br> ALLOW crystals to form |
|  | (d) |  | the range (it) will be wider $\checkmark$ | 1 | ALLOW (it will be) lower ALLOW recrystallized product will be higher/narrower range |
|  | (e) |  | $\begin{aligned} & \text { FIRST CHECK THE ANSWER ON THE ANSWER LINE } \\ & \text { If answer }=45-46(\%) \text { award } 2 \text { marks } \\ & (138.0 \mathrm{~g} 2 \text {-hydroxybenzoic acid } \rightarrow 180.0 \mathrm{~g} \text { aspirin }) \\ & 1.15 \mathrm{~g} 2 \text {-hydroxybenzoic acid } \rightarrow(1.15 / 138.0 \times 180.0) \\ & =1.50 \mathrm{~g} \text { aspirin } \checkmark \\ & \% \text { yield }=(0.68 / 1.50 \times 100)=45(.3)(\%) \checkmark \end{aligned}$ | 2 | Calculates number of moles as <br> $1.15 / 138=0.00833 \mathrm{~mol}$ and $0.68 / 180=0.00378 \mathrm{~mol}(1)$ Allow ecf from incorrect value of Mr from 3a(iii) $\% \text { yield }=0.00378 / 0.00833 \times 100=45.3 \%$ <br> ALLOW 2 or more sf |



| Question |  | Answer | Marks | Guidance |
| :---: | ---: | :--- | :---: | :--- |
| (g) | No (observable) reaction with paracetamol $\checkmark$ <br> Effervescence/fizzing/bubbling/gas with aspirin $\checkmark$ <br> Aspirin contains a carboxyl/carboxylic acid functional <br> group (ORA) $\checkmark$ | $\mathbf{3}$ | ALLOW 'nothing happens' for 'no reaction' |  |
|  |  | Total | $\mathbf{2 1}$ | ALLOW 'dissolve' for aspirin and 'does not dissolve' for <br> paracetamol <br> ALLOW ' $\mathrm{CO}_{2}$ gas but any other named gas is CON |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (a) | (i) | Reagents ..... acidified (potassium) dichromate(VI) AND <br> Conditions ...... reflux $\checkmark$ | 1 | ALLOW (potassium) dichromate in (sulfuric) acid ( VI ) in 'dichromate( VI )' is not required but must be correct if included <br> ACCEPT $\mathrm{H}^{+} / \mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}$ |
|  | (a) | (ii) | $\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH}+2[\mathrm{O}] \rightarrow \mathrm{CH}_{3} \mathrm{COOH}+\mathrm{H}_{2} \mathrm{O} \checkmark$ | 1 | ALLOW $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{OH}$ for ethanol / 2[O] over the arrow DO NOT ALLOW $\mathrm{C}_{2} \mathrm{H}_{6} \mathrm{O}$ for ethanol or $\mathrm{C}_{2} \mathrm{H}_{4} \mathrm{O}_{2}$ for ethanoic acid as question asks for structural formulae Displayed or skeletal formulae are also both acceptable |
|  | (b) |  | $\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH}+2 \mathrm{O}_{2} \rightarrow 2 \mathrm{CO}+3 \mathrm{H}_{2} \mathrm{O} \checkmark$ | 1 | ALLOW either $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{OH}$ or $\mathrm{C}_{2} \mathrm{H}_{6} \mathrm{O}$ for ethanol ALLOW $\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH}+\mathrm{O}_{2} \rightarrow 2 \mathrm{C}+3 \mathrm{H}_{2} \mathrm{O}$ <br> $\mathrm{OR} \mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH}+2.5 \mathrm{O}_{2} \rightarrow \mathrm{CO}_{2}+\mathrm{CO}+3 \mathrm{H}_{2} \mathrm{O}$ (or doubled) OR any other balanced equation that includes C and/or CO as a product Ignore state symbols |
|  | (c) |  | ethanol and ethanoic acid - both hydrogen-bonds $\checkmark$ ethanoic acid has stronger/more H-bonds / id - id <br> H -bonds stronger than pd-pd / H-bonds are the strongest and stronger bonds take more energy to break (ORA) $\checkmark$ <br> Ethanal - permanent dipole - permanent dipole/pd - pd $\checkmark$ | 4 | ALLOW pd : pd / Van der Waal forces as an alternative to id : id <br> ALLOW <br> Permanent dipole - dipole / dipole - permanent dipole / permanent - permanent dipole |
|  | (d) | (i) | FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer $=\mathbf{7 5 0}\left(\mathbf{c m}^{3}\right)$ award $\mathbf{2}$ marks moles of $\mathrm{CH}_{3} \mathrm{CHO}=(0.55 / 44)=0.0125 \mathrm{~mol}$ $0.0125 \mathrm{~mol} \mathrm{CH}_{3} \mathrm{CHO}$ requires $(21 / 2 \times 0.0125)$ <br> $=0.03125 \mathrm{~mol} \mathrm{O}_{2}$ <br> volume of $\mathrm{O}_{2}=(0.03125 \times 24000)=750\left(\mathrm{~cm}^{3}\right) \checkmark$ | 2 | ALLOW 2 or more sf. Throughout <br> Allow ecf <br> ALLOW vol of $\mathrm{O}_{2}=300\left(\mathrm{~cm}^{3}\right)$ from correct calculation of Moles $\mathrm{CH}_{3} \mathrm{CHO} \times 24000$ ie $0.0125 \times 24000$ for 1 mark |
|  | (d) | (ii) | $\begin{aligned} & \mathrm{Mr} \mathrm{CO}_{2} /\left(\mathrm{Mr} \mathrm{CO}_{2}+\mathrm{Mr} \mathrm{H}_{2} \mathrm{O}\right) \times 100 \\ & 44 /(44+18) \times 100=71 \% \checkmark \end{aligned}$ | 1 | Correct answer $=71 \%$ without working scores Allow 2 or more sf |


| Quest | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: |
| (e) | Please refer to the marking instructions on page 4 of this mark scheme for guidance on how to mark this question <br> Level 3 <br> (5-6 marks) <br> Calculates both the empirical formula from the \% composition and the molecular formula using MS. <br> AND <br> Uses IR spectrum to identify $\mathrm{C}=\mathrm{O}$ bond and one other bond (or lack of bond) present in the structure. <br> AND <br> Draws correct detailed conclusion for ester formula/structure from above and MS fragment data <br> The conclusion relates to the evidence and is clear and logically structured. <br> Level 2 <br> (3-4 marks) <br> Concludes A is an ester supported by evidence from molecular formula/Mr and some IR data <br> OR <br> Concludes it is ethanal supported by empirical formula and appropriate evidence using IR / MS data <br> The conclusion relates to the limited evidence and is clear and logically structured. <br> Level 1 <br> (1-2 marks) <br> States empirical formula or Mr of Compound A using evidence from MS / \% composition <br> OR <br> suggests it is an ester/aldehyde from IR evidence alone <br> Pieces of evidence given are related in some way <br> Level 0 <br> Insufficient or irrelevant science | 6 | Indicative scientific points may include: <br> Formula/Mass Spec evidence <br> - empirical formula $\mathrm{C}_{2} \mathrm{H}_{4} \mathrm{O}$ with calculation from \% data, $\mathrm{C}=(54.5 / 12.0)=4.54, \mathrm{H}=(9.1 / 1.0)=9.1, \mathrm{O}$ $=(36.4 / 16.0)=2.275, \mathrm{C}=2(1.99), \mathrm{H}=4, \mathrm{O}=1)$ <br> - $\mathrm{Mr}=88$ identified from molecular ion peak in mass spectrum <br> - molecular formula $=$ empirical formula (mass) $\times 2$ $=\mathrm{C}_{4} \mathrm{H}_{8} \mathrm{O}_{2}$ <br> Infra-red evidence <br> - $\mathrm{C}=\mathrm{O}$ bond in ester (aldehyde/ester) present, absorption is 1740 OR in range 1720-1740 (cm-1) <br> - O-H bond in carboxylic acid not present, no (broad) absorption in range 2500-3300 (cm-1) <br> - C-O bond present as absorption in range 1250 -$1300\left(\mathrm{~cm}^{-1}\right)$ <br> Conclusion <br> - ester <br> - $\mathrm{CH}_{3} \mathrm{COOC}_{2} \mathrm{H}_{5}$ (structure or name, ethyl ethanoate) <br> - because of fragment(s) identified in mass spectrum |
|  | Total | 16 |  |

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