

**Chemistry B (Salters)**

Advanced GCE

Unit **F334**: Chemistry of Materials

**Mark Scheme for June 2013**

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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.

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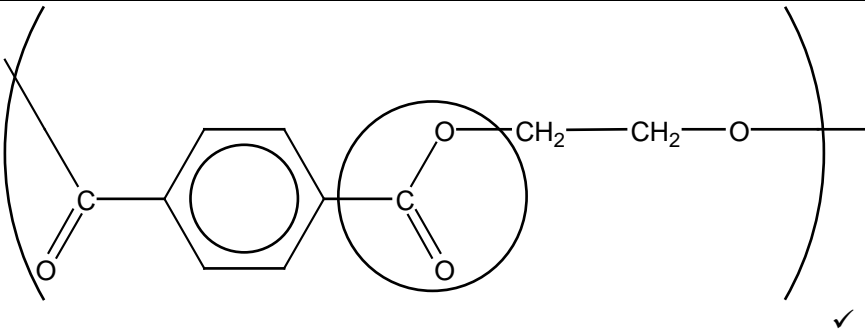
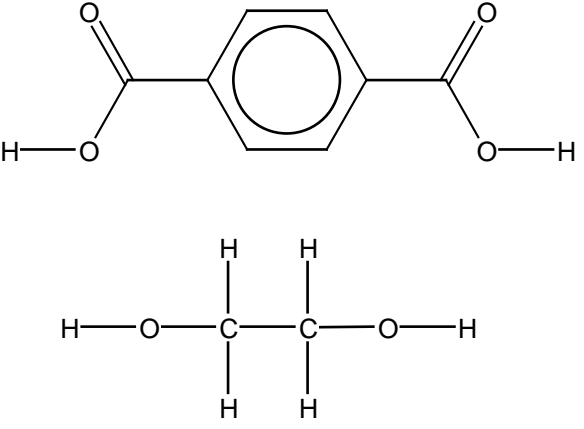
## 1. Annotations

Annotation	Meaning
	Benefit of doubt
	Contradiction
	Cross
	Error carried forward
	Ignore
	Not answered question
	Benefit of doubt not given
	Not good enough
	Rounding error
	Repeat
	Noted but no credit given
	Error in no. of significant figures
	Tick
	Omission mark

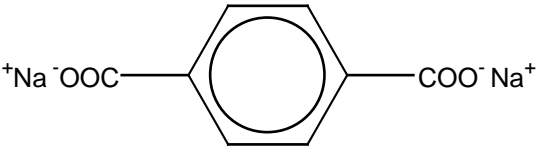
2. Abbreviations, annotations and conventions used in the detailed Mark Scheme (to include abbreviations and subject-specific conventions).

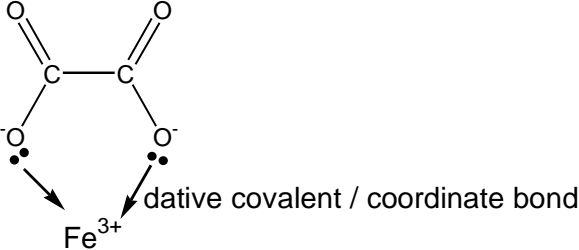
<b>Annotation</b>	<b>Meaning</b>
<b>DO NOT ALLOW</b>	Answers which are not worthy of credit
<b>IGNORE</b>	Statements which are irrelevant
<b>ALLOW</b>	Answers that can be accepted
( )	Words which are not essential to gain credit
—	Underlined words must be present in answer to score a mark
<b>ECF</b>	Error carried forward
<b>AW</b>	Alternative wording
<b>ORA</b>	Or reverse argument

All questions should be annotated with ticks to show where marks have been awarded in the body of the text.  
All questions where an ECF has been applied should also be annotated with the ECF annotation.

Question		Answer	Marks	Guidance
1	(a) (i)		1	<b>ALLOW</b> 'circle' to include the two adjacent C atoms
	(ii)	<p>permanent (dipole) – permanent dipole (bond/forces) ✓</p> <p>instantaneous (dipole) – induced dipole (bond/forces) ✓</p>	2	<p><b>NOT</b> just 'permanent dipole bond/forces'</p> <p><b>DO NOT ALLOW</b> pd etc</p> <p><b>ALLOW</b> van der waals</p> <p><b>IGNORE</b> permanent (dipole) – induced dipole (bond/forces)</p> <p><b>Each mention of any other type of bond in addition to both of these is a CON</b></p>
	(iii)	 <p>1 mark for each monomer correct ✓✓</p>	2	<p><b>ALLOW</b> -COC/ for -COOH</p> <p><b>ALLOW</b> -OH, HOCH<sub>2</sub>CH<sub>2</sub>OH</p>

Question		Answer	Marks	Guidance
1	(a) (iv)	condensation / esterification <b>AND</b> water / H <sub>2</sub> O ✓	1	If –COCl/ in (a) (iii) <b>MUST</b> have HCl/ <b>NOT</b> H <sub>2</sub> O
	(b) (i)	vapours are condensed / turned into liquid <b>AW</b> ✓  <b>mixture needed to be heated for a long time</b> (to break down polymers / for reaction to occur ) <b>OR</b> <b>no reactants or products / vapours are lost</b> <b>OR</b> <b>high temperature required for reaction</b> ✓	2	<b>IGNORE</b> any reference to 'fire' / toxic <b>NOT</b> just 'vapours' fall back down etc. i.e. 'vapours' need state change  <b>NOT</b> solution
	(ii)	<b>choice of solvent:</b> dissolves salt well at higher temp but very little / none at room temp <b>AW</b> ✓  <b>method:</b> 1. use <b>hot solvent</b> ✓ 2. dissolve in <b>minimum amount</b> of solvent ✓ 3. <b>leave to crystallise/cool</b> ✓ 4. <b>filter</b> off crystals, (soluble) <b>impurities are left in filtrate / solution</b> <b>AW</b> ✓ 5. wash with (cold) <u>solvent</u> and dry ✓  <b>MP4 is QWC – i.e. for linking removal of impurities to filtration</b>	6	<b>ALLOW</b> boiling point of solvent is lower than the melting point of the salt  <b>IGNORE</b> any reference to <b>INSOLUBLE</b> impurities
	(c) (i)	bonds (in a molecule) <u>absorb</u> ✓  specific/different/certain (IR) frequencies/wavelengths ✓  <b>alternative for 1<sup>st</sup> &amp; 2<sup>nd</sup> marking points:</b> <u>absorbing different frequencies</u> ✓ causes different bonds to vibrate ✓	2	<b>IGNORE</b> references to energy <b>NOT</b> 'electrons in bonds'

Question		Answer	Marks	Guidance
1	(c) (ii)	<p><b>structure of A</b></p>  <p style="text-align: right;">✓</p> <p>No OH bond since no <u>broad</u> absorption peak above about (2500-3200) / 3000 (cm<sup>-1</sup>) ✓</p> <p>C=O absorption peak at about 1720-1740 (cm<sup>-1</sup>) (so must be carboxylate <b>AW</b>) ✓</p> <p><b>structure of B</b></p> <p>HOCH<sub>2</sub>CH<sub>2</sub>OH ✓</p> <p>OH bond since (broad) absorption peak about 3200-3600 (cm<sup>-1</sup>) ✓</p>	5	<p><b>REMEMBER</b> marking points are independent <b>ALLOW</b> any correct structural formulae</p> <p><b>ALLOW</b> without Na<sup>+</sup></p> <p><b>IR data may be drawn on the spectra, please check</b></p> <p>reference to any functional group other than a carboxylic acid / carboxylate is a <b>CON</b> <b>ALLOW</b> frequency within stated range</p> <p><b>ALLOW</b> any correct structural formula</p>
	(d) (i)	temperature <u>below</u> which the polymer turns glassy/brittle ✓	1	'below' may be expressed by reducing temperature / cooling / shown in a diagram
	(ii)	it would soften / melt / turn into liquid/fluid ✓	1	
	(e) (i)	chains are further apart / less close together in PBT ✓ so has <u>weaker</u> intermolecular bonding/forces than PET ✓ so chains in PBT can move over one another more easily ✓	3	<b>ORA</b> <b>IGNORE</b> references to ordered chains etc. <b>IGNORE</b> fewer/less imb/fs
	(ii)	butane-1,4-diol  butane / butan / but <b>AND</b> diol = ✓ 1,4- = ✓	2	<b>IGNORE</b> commas & dashes  '1,4-' must be between 'butane' & 'diol'
<b>Total</b>			<b>28</b>	

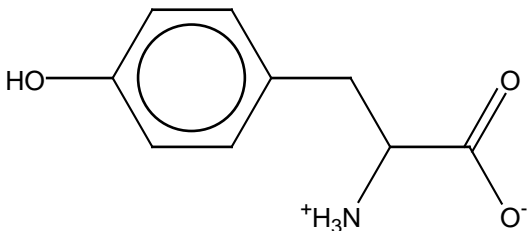
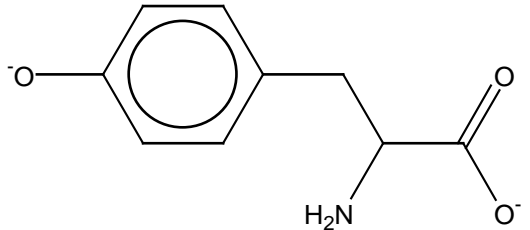
Question		Answer	Marks	Guidance
2	(a) (i)	 <p>dative covalent / coordinate bond</p> <p>lone pairs as shown ie must link to bond (any type of drawn line) ✓</p> <p><b>two</b> bonds shown as arrows from O<sup>-</sup> pointing to a single <b>Fe<sup>3+</sup></b> ✓</p> <p>dative (covalent)/coordinate bond labelled (anywhere on diagram) ✓</p>	3	<p><b>ECF</b> allow this marking point if the C=O are used instead of the O<sup>-</sup> (so max mark of 2 if incorrect Os used)</p> <p><b>CON</b> if any other bond is specifically labelled</p>
	(ii)	<p><math>[\text{Fe}(\text{C}_2\text{O}_4)_3]^{3-}</math> ✓</p> <p><u>octahedral</u> ✓</p>	2	<p><b>ALLOW</b> without square brackets</p> <p><b>IGNORE</b> separate correct charges for <b>both</b> Fe &amp; C<sub>2</sub>O<sub>4</sub> as long as overall charge 3- is shown</p> <p><b>ALLOW</b> structural formula</p>
	(b) (i)	<p>the <math>E^\ominus</math> of <b>CO<sub>2</sub>/(COOH)<sub>2</sub></b> half-cell is <b>more negative/less positive</b> than that of the <b>Fe<sup>3+</sup>/Fe<sup>2+</sup></b> half-cell</p> <p><b>OR</b></p> <p><math>E_{\text{cell}} = +1.26 \text{ V}</math>, so reaction is feasible ✓</p> <p><b>(COOH)<sub>2</sub></b> will release electrons / reduce Fe<sup>3+</sup></p> <p><b>OR</b> Fe<sup>3+</sup> will gain electrons / oxidise <b>(COOH)<sub>2</sub></b> ✓</p>	2	<p><b>ALLOW</b> (in this question only) <math>E^\ominus</math> of the <b>Fe<sup>3+</sup></b> half-cell etc. (there are only 2 half-cells given)</p>
	(ii)	<p><math>2\text{Fe}^{3+}(\text{aq}) + (\text{COOH})_2(\text{aq}) \rightarrow 2\text{Fe}^{2+}(\text{aq}) + 2\text{CO}_2(\text{g}) + 2\text{H}^+(\text{aq})</math></p> <p>correct formulae <b>AND</b> balanced ✓</p> <p>state symbols correct ✓</p>	2	<p><b>ACCEPT</b> CO<sub>2</sub>(aq)</p> <p>If <b>balanced with electrons</b> on either side max mark = 1</p> <p>State symbol mark may be awarded if species are correct even if equation is reversed</p>

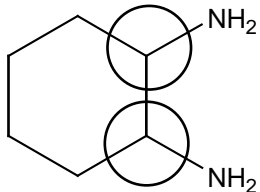
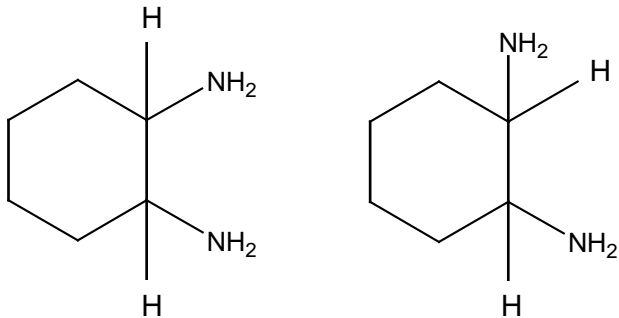


Question		Answer	Marks	Guidance
2	(b) (iii)	<p style="text-align: center;">3d                                      4s</p> <p>any 6 electrons for Fe<sup>2+</sup> and 5 for Fe<sup>3+</sup> = 1<sup>st</sup> mark ✓ correct arrangements for both = 2<sup>nd</sup> mark ✓</p>	2	
	(iv)	half-filled <u>d</u> shell (is more stable) AW ✓	1	<b>AW</b> eg only 1 electron in <b>each</b> of the d orbitals
	(c) (i)	<p>1. manganate(VII) solution in <u>burette</u> ✓</p> <p>2. <u>pipette</u> known/stated volume of ethanedioate solution <b>OR</b> graduated/volumetric <u>pipette</u> for ethanedioate solution ✓</p> <p>3. add acid ✓</p> <p>4. warm / heat solution / 60°C ✓</p> <p>5. titrate (AW) until <u>pink</u> colour persists/remains AW ✓</p> <p>6. <i>no indicator needed because only MnO<sub>4</sub><sup>-</sup>(aq) is coloured</i> <b>OR</b> a colour change takes place during the reaction AW ✓</p> <p><b>QWC</b> to gain the 1<sup>st</sup> mark the spelling of <b>burette</b> has to be correct at least <b>once in the answer</b></p>	6	<p>If pipette &amp; burette <b>used wrong way round</b> then 1 mark only for points 1 and 2 <b>AND</b> mark 5 is only available if purple changes to colourless <b>IGNORE</b> 'bulb'</p> <p><b>ALLOW</b> acid / 'acidified' in either solution</p> <p><b>NOT</b> purple alone here <b>MUST HAVE</b> pink</p> <p><b>ALLOW</b> purple–colourless <b>NOT</b> colourless–purple <b>INCORRECT</b> colour change is a <b>CON</b> eg orange–green, purple–pink etc.</p>

Question			Answer	Marks	Guidance
2	(c)	(ii)	<p>1. moles of <math>\text{KMnO}_4</math> used in titration = <b>18.40/1000</b> x <b>0.0500</b> ✓ = 0.0009200</p> <p>2. moles potassium ethanedioate used in titration = <b>5/2</b> x answer from 1 ✓ = 0.002300</p> <p>3. moles potassium ethanedioate in 100 <math>\text{cm}^3</math> = <b>4</b> x answer from 2 ✓ = 0.009200</p> <p>4. <math>M_r</math> of <math>\text{K}_2\text{C}_2\text{O}_4 \cdot \text{H}_2\text{O}</math> = <b>184 / 184.2</b> ✓</p> <p>5. mass potassium ethanedioate in 100 <math>\text{cm}^3</math> = <b>184.2</b> x (answer from 3) ✓ = 1.6946</p> <p>6. Answer = <b>1.695</b> or <b>1.69 g</b> ✓ to 4 or 3 sig figs</p>	6	<p>The marks are awarded for the working out given in bold</p> <p><b>ALLOW</b> ecf between each step</p> <p>3. may be done in <b>2 steps</b> via moles <math>\text{dm}^{-3}</math> and still scores only 1 mark</p> <p><b>Note:</b> 1 error means only 1 mark is lost eg incorrect <math>M_r</math> eg these are probably 5 marks but place ticks appropriately: <b>0.200</b> (/ <math>M_r</math> instead of x <math>M_r</math>) <b>3.39</b> (/2 missing) <b>5010</b> (in step 1: x(1000/18.40) rather than /)</p> <p>the following is probably 4 marks <b>0.42</b> (missing x4 and incorrect sf)</p>

Question		Answer	Marks	Guidance
2	(d) (i)	<p><b>X to Y:</b> increases ✓ rate speeds up as (catalyst) <math>\text{Mn}^{2+}</math> is formed ✓</p> <p><b>Y to Z:</b> decreases ✓ rate slows as reactants / <math>\text{C}_2\text{O}_4^{2-}</math> / <math>\text{MnO}_4^-</math> are/is used up / as concentrations of reactants fall ✓</p>	4	
	(ii)	colorimetry / use a colorimeter / visible spectroscopy / visible spectrophotometry ✓	1	<b>ALLOW</b> conductivity / gas volume <b>IGNORE</b> pH
	(iii)	<p><b>EITHER</b> (colourless) effervescence/fizzing/bubbling AW ✓</p> <p><b>OR</b> (purple/pink) colour fades AW ✓</p>	1	<p><b>IF MORE THAN ONE ANSWER</b> <b>MARK FIRST IN LIST ONLY</b></p> <p><b>NOT</b> gas forms <b>NOT</b> colour change <b>IGNORE</b> references to absorbance incorrect colour is a <b>CON</b></p>
<b>Total</b>			<b>30</b>	

Question		Answer	Marks	Guidance
3	(a) (i)	phenol / hydroxyl ✓ carboxyl / carboxylic acid ✓ amino / amine ✓	3	<b>NOT</b> hydroxide, alcohol <b>IGNORE</b> 'primary' but 'secondary' is a <b>CON</b>
	(ii)	(neutral) FeCl <sub>3</sub> / iron(III) chloride ✓ turns purple / violet (phenol present) ✓	2	<b>NOT</b> blue or pink If initial colour is given, it must be yellow, orange or colourless otherwise <b>CON</b>
	(b) (i)	contains a positive charge and a negative charge ✓  structure <b>ALL</b> correct ✓	2	<b>IGNORE</b> dipolar <b>MUST</b> indicate that there are only 1+ and 1- charge present this may be indicated by the structure drawn
	(ii)	acidic ✓ (because it has a) phenol group ✓	2	<b>ALLOW</b> structural formula for phenol <b>IGNORE</b> references to -COOH & -NH <sub>2</sub> groups
	(iii)	 1 mark for phenol group reacted correctly ✓ 1 mark for rest of the molecule correct ✓	2	<b>IGNORE</b> any Na <sup>+</sup> ions

Question		Answer	Marks	Guidance
3	(c)	(i)	2	<b>IGNORE</b> complementary <b>IGNORE</b> enzyme <b>NOT</b> 'react with'
		(ii)	1	<b>ALLOW</b> 'benefits outweigh side effects' <b>IGNORE</b> reference to 'disease'
	(d)	<p><b>optical isomers:</b></p>  <p>chiral Cs shown on the diagram ✓</p> <p>4 different groups around (each) C <b>OR</b> not superimposable on their mirror image ✓</p> <p><b>cis-trans isomers:</b> C-C between the chiral (AW) atoms is prevented from rotating by the ring structure ✓</p>  <p>2 correct diagrams for isomers ✓</p>	4	<p><b>NOT</b> 'functional groups'</p> <p><b>IGNORE</b> references to 'ring rotation'</p> <p>H's may not necessarily be shown as in MS <b>ACCEPT</b> if NH<sub>2</sub> groups only are shown with lines/wedges/dotted lines etc. <b>MUST CONVINC</b>e that we have cis &amp; trans isomers</p>
<b>Total</b>			<b>18</b>	

Question		Answer	Marks	Guidance
4	(a) (i)	order for $[\text{CH}_3\text{Cl}] = 1 \checkmark$ order for $[\text{H}_2\text{O}] = 2 \checkmark$	2	
	(ii)	rate = $k [\text{CH}_3\text{Cl}] [\text{H}_2\text{O}]^2 \checkmark$ overall order = 3 $\checkmark$	2	<b>ALLOW</b> with 'x's in rate equation <b>ECF</b> from (i) <b>ECF</b> from rate equation
	(iii)	slow step/rate determining step involves <b>one</b> $\text{CH}_3\text{Cl}$ (molecule) so it is 1 <sup>st</sup> order AW $\checkmark$ one $\text{OH}^-$ formed from the <b>two</b> $\text{H}_2\text{O}$ (molecules) so 2 <sup>nd</sup> order with respect to $\text{H}_2\text{O}$ AW $\checkmark$	2	<b>IGNORE</b> 'rds'
	(b)	hydrochloric acid $\checkmark$ methanoic acid $\checkmark$	2	<b>ALLOW</b> hydrogen chloride, formic acid <b>IGNORE</b> formulae
	(c)	<b>acidified</b> $\checkmark$ (potassium) <b>dichromate</b> / (sodium) dichromate / $\text{Cr}_2\text{O}_7^{2-}$ $\checkmark$  (add reagent to alcohol and) <b>distil</b> off aldehyde as it is formed $\checkmark$	3	any concentration of sulfuric acid / $\text{H}_2\text{SO}_4$ <b>DO NOT ALLOW</b> hydrochloric <b>OR</b> nitric acids  use of 'reflux' is a <b>CON</b>
	(d)	$1.56 \times 10^{-4} = k \times 1.82 \times 10^{-3} \checkmark$  $k = 0.0857 / 0.086 \checkmark$  $\text{s}^{-1} \checkmark$	3	<b>ALLOW</b> any correct rearrangement of equation <b>CORRECT ANSWER</b> gets both marks <b>ALLOW</b> two or more sig figs
<b>Total</b>			<b>14</b>	

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