



# Chemistry

Advanced GCE A2 7882

Advanced Subsidiary GCE AS 3882

# **Mark Schemes for the Units**

# January 2009

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# **2811 Foundation Chemistry**

Q	Question		Expected Answers	Marks	Additional Guidance
1	(a)	i	20√	1	
		ii	2√	1	
		iii	5 √	1	
	(b)	i	69.8 ✓	1	
		ii	96.0 ✓	1	Allow 96
		iii	moles of NaNO <sub>3</sub> = 0.05 $\checkmark$	2	4.8 g worth 1 (wrong $M_r$ )
			mass = 0.05 x 85 = 4.25 (g) ✓		Accept 4.3 but not 4.2
					(ecf for calculated moles x 85)
		iv	$2.51 \times 10^{21} $	1	<b>Allow</b> $2.5 \times 10^{21}$
					Calc: 2.508333333 x 10 <sup>21</sup>
					<b>Allow</b> calc value and any degree of correct rounding down to $2.5 \times 10^{21}$
	(c)		(+)7 ✓	1	Sign not required but do not credit '–7' Accept VII
			Total	9	

Questic	en Expected Answers	Marks	Additional Guidance
2 (a)	Energy change when each atom in 1 mole $\checkmark$ of gaseous atoms $\checkmark$ loses an electron $\checkmark$ (to form 1 mole of gaseous 1+ ions)	3	Not 'element' alone Compensate missed marks from correct equation
(b)	Si ✓ Sharp rise in successive ionisation energy between 4th and 5th IE ✓ marking a change to a new shell / there are 4 electrons in the outer shell ✓	3	Not consequential Not sub-shell
(c)	<ul> <li>atomic radii increases/ there are more shells/atoms get bigger ✓</li> <li>there is more shielding/ more screening ✓</li> <li>ionisation energy decreases because the Increased shielding and distance outweigh the increased nuclear charge / the nuclear attraction decreases ✓</li> </ul>	3	<ul> <li>USE annotations with ticks, crosses, con, ecf, etc for this part.</li> <li>'down the group' not required</li> <li>'more' is essential allow 'more electron repulsion from inner shells'</li> <li>Allow 'nuclear pull' ignore any reference to 'effective nuclear charge'</li> </ul>
	Total	9	

Qı	uestio	n	Expected Answers	Marks	Additional guidance
3	(a)		Ca <sup>2+</sup> : 20 protons; 18 electrons ✓ Cl <sup>-</sup> : 17 protons; 18 electrons ✓	2	
	(b)		cation shown with either 8 or 0 electrons AND anion shown with 8 electrons AND correct number of crosses and dots $\checkmark$ Correct charges on both ions $\checkmark$ $\begin{bmatrix} c_a \end{bmatrix}^{2^+} \begin{bmatrix} c_1 \\ c_1 \end{bmatrix}^-$ $\begin{bmatrix} c_1 \\ c_1 \end{bmatrix}^-$	2	For 1st mark, if 8 electrons shown around cation then 'extra' electron(s) around anion must match symbol chosen for electrons in cation. <i>Circles not required</i> Ignore inner shell electrons For charges, Allow: 2[CI <sup>-</sup> ] 2[CI <sup>-</sup> [CI <sup>-</sup> ] <sub>2</sub> (brackets not required except for last one) Do not allow: for CaCl <sub>2</sub> , [Cl <sub>2</sub> ] <sup>2-</sup> [Cl <sub>2</sub> ] <sup>-</sup> [2CI] <sup>2-</sup> [CI] Max 1 if only one CI <sup>-</sup>
	(c)		solid: ions are fixed (AW) $\checkmark$ aqueous: ions are free (to move) (AW) $\checkmark$	2	If charge carriers are wrong but comparison is given, then award one mark, <i>e.g.</i> solid: <i>electrons</i> are fixed in lattice <i>AND</i> aqueous: <i>electrons</i> are free to move $\checkmark$ (1 mark
	(d)	i	molar mass CaCO <sub>3</sub> : 100.1 (g mol <sup>-1</sup> ) ✓ 4.85/100.1 = 0.0485 mol ✓	2	Not 100 for molar mass calc. 0.048451548 <b>Allow</b> rounding of calculator value back to 2 sig figs allow 0.048-0.049 <b>ECF</b> If working shown for an incorrect molar mass, then the 2nd mark can be awarded as 4.85/calculated molar mass
		ii	5.38 or 5.39 g or 5.4 g ✓	1	For information: 0.0485 x 111.1 = 5.39 0.048451548 x 111.1 = 5.38

Q	Question		Expected Answers	Marks	Additional guidance
					<b>ECF</b> : moles from (i) x 111.1 or 111
		iii	0.0970 or 0.097 or 0.0969 ✓	2	For information: $2 \times 0.0485 = 0.0970 \text{ mol}$ $2 \times 0.048451548 = 0.0969$ ECF moles from (i) x 2
			volume = 64.7 or 64.6 cm <sup>3</sup> or 65 $\checkmark$		For information (0.0970/1.50) x 1000 = 64.7 cm <sup>3</sup> (0.0969/1.50) x 1000 = 64.6 cm <sup>3</sup> ECF (moles above/1.50) x 1000
	(e)		$\begin{array}{rcl} \text{Ca/CaO/Ca(OH)}_2 \checkmark \\ \text{Ca} &+ & 2\text{HCI} &\longrightarrow & \text{CaCI}_2 &+ \text{H}_2 \ / \\ \text{CaO} &+ & 2\text{HCI} &\longrightarrow & \text{CaCI}_2 &+ & \text{H}_2\text{O} \ / \\ \text{Ca(OH)}_2 &+ & 2\text{HCI} &\longrightarrow & \text{CaCI}_2 &+ & 2\text{H}_2\text{O} \ \checkmark \end{array}$	2	Ignore state symbols Allow any other suitable alternatives
	(f)	i	Ca : H : S : O = $19.82/40.1$ : $0.99/1$ : $31.74/32.1$ : $47.45/16$ or 1 : 2 : 2 : 6 $\checkmark$ empirical formula = CaH <sub>2</sub> S <sub>2</sub> O <sub>6</sub> $\checkmark$	2	Using atomic numbers gives CaHS <sub>2</sub> O <sub>6</sub> worth 1 Allow Ca(HSO <sub>3</sub> ) <sub>2</sub> !
		ii	$Ca(OH)_2 + 2SO_2 \longrightarrow CaH_2S_2O_6 \checkmark$	1	If you see it, allow Ca(HSO <sub>3</sub> ) <sub>2</sub> !
			Total	16	

Q	uestic	on	Expected Answers	Marks	Additional Guidance
4	(a)		down group/from CI to I/, number of electrons/shells increases ✓	3	Answers involving ionisation energies score 0
			more/ stronger/ van der Waals'/ intermolecular forces/ induced dipoles/ instantaneous dipoles ✓		
			greater forces to break/more energy has to be put in to break forces $\checkmark$		
	(b)		$I_2 \rightarrow IO^-$ , 0 to +1 $\checkmark$ : oxidised	3	Sign not required but do not credit '-1'
			$I_2 \rightarrow I^-$ , 0 to $-1 \checkmark$ : reduced		' ' Sign required here
			<b>correct</b> 'oxidised' <b>and</b> 'reduced' above/I is both oxidised and reduced / disproportionation ✓		
	(c)	i	goes orange/red/yellow ✓	2	Ignore brown
			$Cl_2 + 2Br^- \longrightarrow Br_2 + 2Cl^- \checkmark$		Ignore spectator ions
		ii	Ag <sup>+</sup> (aq) + Cl <sup>-</sup> (aq) $\longrightarrow$ AgCl(s) correct equation $\checkmark$ correct state symbols $\checkmark$	2	Allow state symbols for (slightly) incorrect equations
	(d)	i	attraction of an atom for electrons $\checkmark$ in a (covalent) bond/ bonding pair $\checkmark$	2	
		ii	correct 3-D tetrahedral shape shown showing one outward	2	For bond into paper, accept:
			wedge and 1 inward wedge; 3 bonds below horizontal $\checkmark$ correct dipoles: $\delta$ + on C and $\delta$ – on each Cl $\checkmark$		
			$CI^{\delta-}$		Allow correct shape with no atom labels:
					Only need to show one dipole
		iii	the polarities/ dipoles cancel out / the molecule is symmetrical	1	
			Total	15	

Question	Expected Answers	Marks	Additional Guidance
5	Magnesium structure/bonding: giant ✓ metallic ✓ conducts by delocalised/free/mobile electrons ✓ melting point high because of the electrostatic attraction / attraction between (positive) ions and electrons ✓ Diamond does not conduct/poor conductor: no mobile charge carriers/electrons/ions ✓ structure/bonding: giant (✓) covalent ✓ melting point: high because strong/ lots of (covalent) bonds are broken ✓	10	USE annotations with ticks, crosses, con, ecf, etc for this part. Credit information if given in annotated diagrams Watch out for contradictions, especially of bonding type Allow: positive ions with a sea of electrons for both structure and bonding marks if labelled, one if not.
	<ul> <li>Ice         <ul> <li>does not conduct:                 no mobile charge carriers/electrons/ions ✓                 structure/bonding:                 H-bonds/intermolecular forces/ simple molecular ✓                 melting point:                 Low because H bonds/intermolecular/ weak forces                 between molecules (are broken)/ higher than expected                 because H-bonds broken/ H-bonds stronger than other                 (named) intermolecular forces ✓</li> </ul> </li> <li>QWC – At least two sentences that show legible text with         <ul> <li>accurate spelling, punctuation and grammar so that the                 meaning is clear. ✓</li> </ul> </li> </ul>	1	must refer to bonds being broken once QWC mark <b>must</b> be indicated with a tick or cross through the Quality of Written Communication prompt at the bottom of page 9.
			Then scroll up to start of (b), counting ticks.

# 2812 Chains and Rings

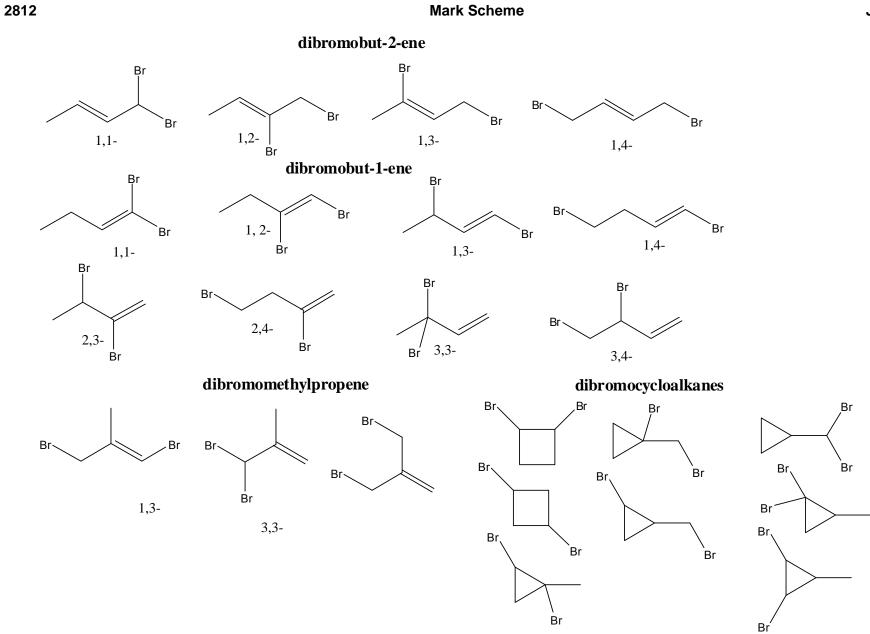
C	uest	ion	Expected Answers	Marks	Additional Guidance
1	(a)		compound/molecule that contains carbon & hydrogen only ✓	1	allow hydrocarbons contain carbon & hydrogen <u>only</u> allow molecules that contain carbon & hydrogen <u>only</u>
	(b)		$C_{14}H_{30} \longrightarrow C_8H_{18} + C_6H_{12} \checkmark$	1	allow $CH_3(CH_2)_{12}CH_3 \longrightarrow CH_3(CH_2)_6CH_3 + C_6H_{12}$ allow any isomer of $C_6H_{12}$ or any combination of alkenes that add up to $C_6H_{12}$ .
	(c)	(i)	$\begin{array}{c c} & & & \\ &$	1	allow different orientations as long as the two methyl groups are on adjacent Cs
		(ii)	hydrogen/H₂✓	1	no other correct response
	(d)	(i)	$\begin{array}{c c} & & & & CH_3 \\ & & & & \\ & & & \\ H_3C \end{array} \begin{array}{c} CH_2 \\ CH_3 \\ CH_3 \end{array} \begin{array}{c} CH_3 \\ CH_3 \end{array} \begin{array}{c} CH_3 CH_2CH_2(CH_3)C(CH_3)_3 \\ CH_3 CH_3 CH_3 CH_3CH_2(CH_3)C(CH_3)C(CH_3)_3 \end{array}$	1	<b>allow</b> any unambiguous form of 2,2,3-trimethylpentane

Q	uesti	ion	Expected Answers	Marks	Additional Guidance
		(ii)	$C_8H_{18} + 12\frac{1}{2}O_2 \longrightarrow 8CO_2 + 9H_2O \checkmark \checkmark$ 1 mark if all formulae are correct both marks if correctly balanced	2	allow $2C_8H_{18} + 25O_2 \longrightarrow 16CO_2 + 18H_2O$ allow structural, displayed or skeletal formula of $C_8H_{18}$ .
	(e)	(i)	(feedstock is obtained) from plants ✓ which can be re-grown ✓	2	allow made from sugar cane/beet/biomass for 1 mark not allow just sugar allow made from sugar because it can be re-grown for 2 marks not allow just fermentation allow fermentation from/of plants for first marking point
		(ii)	CO₂ used in photosynthesis is balanced by CO₂ released in combustion/ it is carbon neutral ✓	1	not allow does not produce greenhouse gasesallow doesn't emit any oxides of nitrogen/sulphurnot allow doesn't produce toxic gases/acid rainIf two statements are made and one is incorrect the markis lost e.g. is carbon neutral and does not producegreenhouse gases this gets * con
			Total	10	

G	Quest	ion	Expected Answers	Marks	Additional Guidance
2	(a)	i	1,1-dibromomethylpropene√	1	allow 1,1-dibromo-2-methylpropene allow 2-methyl-1,1-dibromopropene allow methyl-1,1-dibromopropene also allow any of the above with prop-1-ene
		ii	<i>M</i> <sub>r</sub> = 213.8 ✓	2	<b>not allow</b> $M_r = 214$ for first mark
			% = (159.8/213.8) x 100 = 74.7 ✓		<b>allow</b> any of: % = 75/74.74 or any correct rounding up to and including the calculator value of 74.74275023
					allow ecf for correct rounding of 74.76635514 if used <i>M</i> <sub>r</sub> 214 ecf for correctly calculating percentage from incorrect M <sub>r</sub> 37.4% scores 1 mark
		(iii)	any dibromobut-1-ene any dibromobut-2-ene (except 2,3-dibromobut-2-ene) any dibromomethylpropene (except 1,1-dibromomethylpropene) any dibromocyclobutane any dibromomethylcyclopropane	1	see page 10 at end of question for skeletal formulae of acceptable isomers Most common <b>incorrect</b> response is <i>trans</i> -2,3-dibromobut-2-ene Br C C C C C C C C
	(b)	i	decolourised ✓	1	not allow goes clear / discoloured allow turns colourless/orange colour disappears ignore "clear" if "decolourises and goes clear" i.e. not 'CON'
		ii	electrophilic addition ✓	1	
		iii	molecular formula = $C_4H_6Br_4 \checkmark$ empirical formula = $C_2H_3Br_2 \checkmark$	2	<b>allow</b> ecf from molecular formula $C_xH_yBr_z$

Quest	tion	Expected Answers	Marks	Additional Guidance
(c)		$H \xrightarrow{CH_3 Br} H \xrightarrow{CH_3 Br} H \xrightarrow{CH_3 Br} H \xrightarrow{H} H$	2	Ignore bond linkage
(d)	i	B is symmetrical ✓	1	allow A isn't symmetrical ignore A is asymmetric
	ii	$\begin{array}{c c} & CH_3 & Br \\ & &   &   \\ H - C - C - Br \\ &   &   \\ CH_3 & Br \\ & CH_3 & Br \\ &   &   \\ Br - C - C - H \\ &   &   \\ CH_3 & Br \\ & & \checkmark \end{array}$	2	Ignore bond linkage
e)	i	Br OH Br ✓	1	<b>Do not allow</b> bond linkage to H in the OH, bond must <b>clearly</b> go to the O
	ii	reagent: steam/H <sub>2</sub> O <sub>(g)</sub> ✓ conditions: phosphoric acid ✓	2	allow H <sub>2</sub> O but only if temp is quoted above 100°C allow sulphuric acid not allow acid catalyst allow reagent: phosphoric acid ✓ allow conditions: steam ✓ mention of alkali <i>× con</i> acid mark

Question	Expected Answers	Marks	Additional Guidance
(f)	$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	2	$\begin{array}{c ccccccccccccccccccccccccccccccccccc$
			ignore CH <sub>3</sub> bond linkage
	Total	18	



### Q3a should be marked as a complete question NOT by item response

Question	Expected Answers	Marks	Additional Guidance
3 (a) i	100 🗸	1	If incorrect ecf can be awarded for 3a(ii)
ii	0.05 ✓	1	Check for ecf from 3a(i) if incorrect check response to part (iii) which can score all 3 marks as <b>ecf</b> to incorrect answer in (ii)
	moles of cyclohexene = 1.8/82 / 0.02195 / 0.022 ✓ % yield = (0.022/0.05) x 100 = 43.9% ✓ % yield to 2 sig figs = 44% ✓	3	44% scores all 3 marks <b>allow</b> alternative method theoretical mass of cyclohexene = $0.05 \times 82 = 4.1(g)$ % yield = $(1.8/4.1) \times 100 = 43.9\%$ % yield to 2 sig figs = $44\%$ <b>ecf</b> if $M_r$ of cyclohexene is incorrect, the remaining two marks can be awarded e.c.f <b>ecf</b> % yield = $(0.022/incorrect answer to (a)(ii)) \times 100$ for <i>max 3 marks</i> <b>do not allow</b> moles of cyclohexene rounded to 0.02 which will then lead to 40% yield. <b>allow</b> 40% will score 2 out of the 3 available marks <b>allow</b> 36% for <i>max 1 mark</i>

Question	Exp	ected Ansv	wers		Marks	Additional Guidance
Question (b) i ii	Expo (peak between) 3230–3550 OH ✓ Na bubbles/fizzes/effervesces not allow hydrogen gas/ gas evolved			Sence of PCI <sub>5</sub> / SOCI <sub>2</sub> white fumes	Marks 2 1 1	do not allow $2500-3500 \text{ (cm}^{-1})$ For OH allow peak within stated range Ignore any reference to C–O peak allow RCOCI with observation of white fumes and product same as carboxylic acid If manganate(VII) used as oxidising agent then allow marks for observation (purple to colourless/green/brown) and product of cyclohexanone only <b>not allow</b> $C_6H_{11}ONa / C_6H_{11}OOCR / C_6H_{11}CI$ product mark must be related to correct reagent. If no
	charges not essential but do not allow		Ŭ,			product mark must be related to correct reagent. If no reagent then no product mark is possible allow one mark for bromocyclohexane as product if HBr used as reagent but no marks for reagent or observations not allow
		l		Total	10	

281	2
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Q	uesti	ion	Expected Answers	Marks	Additional Guidance
4	(a)		bond angle 109° 28' $\checkmark$	2	<b>allow</b> 109·5/ 109–110
	(b)	i	electron pair donor 🗸	1	allow lone pair (of electrons) donor
		II	$H \rightarrow G \rightarrow $	2	not allow any incorrect charges on reagents <i>*con 1</i> mark
		111	$CH_3CI + 2NH_3 \longrightarrow CH_3NH_2 + NH_4CI \checkmark$	1	allow $CH_3CI + 2NH_3 \longrightarrow CH_3NH_2 + NH_4^+ + CI^-$ not allow $CH_3CI + NH_3 \longrightarrow CH_3NH_2 + HCI$ not allow $CH_3CI + 2NH_3 \longrightarrow CH_3NH_2 + HCI + NH_3$
		iv	methylamine/aminomethane ✓	1	allow even if equation in (b)(iii) is incorrect.

Q	uestion	Expected Answers	Marks	Additional Guidance
	(c)	reaction would be faster ✓ C—I bond is weaker/has lower bond enthalpy (than C—CI bond) ✓	2	second mark is dependent on first mark <i>e.g. reaction is slower because C—I bond is weaker</i> scores no marks. <b>not allow</b> iodomethane / CH <sub>3</sub> I has lower/weaker bond energy/enthalpy <b>not allow</b> C—I bond is longer <b>allow</b> C—I bond is longer, therefore weaker <b>not allow</b> iodine bond is weaker
		Total	9	

Qu	estion	Expected Answers	Marks	Additional Guidance
5 (	(a)	alkanes are non-polar ✓ nucleophiles/electrophiles are <b>attracted</b> to polar substances ✓ C–H bonds are strong ✓ allow max of 2 from 3	2	allow C–H bonds have little/no polarity/no dipoles allow no regions of high or low electron density allow nucleophiles/electrophiles/reagents are <b>not</b> attracted to non-polar substances <b>not allow</b> attacks/reacts as an alternative to attracts allow bonds in alkanes are strong
(	(b)	Free radical substitution $\checkmark$ balanced equation $C_5H_{12}$ + $Br_2 \rightarrow C_5H_{11}Br$ + $HBr \checkmark$ mechanism $Br_2 \longrightarrow 2Br \bullet \checkmark$		if a different alkane is used <b>do not allow</b> mark for either propagation step but the rest can be marked ecf
		Br• + C <sub>5</sub> H <sub>12</sub> → HBr + •C <sub>5</sub> H <sub>11</sub> ✓ •C <sub>5</sub> H <sub>11</sub> + Br <sub>2</sub> → C <sub>5</sub> H <sub>11</sub> Br + Br• ✓ any two free radicals to show termination step ✓		If error in first propagation step ecf can be awarded for second propagation step allow any one of: $2Br \bullet \longrightarrow Br_2$ $Br \bullet + \bullet C_5H_{11} \longrightarrow C_5H_{11}Br$ $\bullet C_5H_{11} + \bullet C_5H_{11} \longrightarrow C_{10}H_{22}$ If H• formed in propagation allow ecf for a termination equation using the H•
		<ul> <li>conditions: uv ✓</li> <li>bond fission: homolytic fission ✓</li> <li>mixed products due to: <ul> <li>multiple substitution of H (in C<sub>5</sub>H<sub>12</sub>)</li> <li>several isomers of C<sub>5</sub>H<sub>11</sub>Br</li> <li>different products could be formed in termination step* any two from three ✓√</li> </ul> </li> </ul>	10	<ul> <li>allow sunlight/high temperature</li> <li>allow homolysis/homolytic cleavage</li> <li>do not allow free radicals are very reactive/difficult to control</li> <li>* must be stated not just assumed if they write more than one termination step.</li> </ul>

(	Question	Expected Answers	Marks	Additional Guidance
	QWC	Well structured answer and uses all three of initiation, propagation and termination correctly $\checkmark$	1	
		Tota	I 13	

## 2813/01 How Far? How Fast?/Experimental Skills 1 Written Paper

Q	Question		on Expected Answers		Additional Guidance
1	(a)	i	reaction slows ✓	1	Not becomes constant
		ii	because there are less particles per unit volume (as the reaction proceeds)/particles further apart/ the concentration decreases ✓		marks are stand alone
				2	
			(rate) of collision decreases ✓		allow successful collisions
	(b)		sketch to show		
			graph starting more steeply $\checkmark$		
			finishing at same level ✓	2	
			Total	5	

Q	uesti	on	Expected Answers	Marks	Additional Guidance
2	(a)	i	energy/enthalpy change to break 1 mole of a (covalent)		do <b>not</b> allow first mark:
			bond 🗸		if energy released
					if break and make
					if ionic
					if heat
			in the gaseous state $\checkmark$		2nd mark is stand alone
				2	ignore 'under standard conditions'
		ii	energy is put in to break the bond/ energy is needed to		not attraction between oppositely charged ions
			overcome the attraction (between electrons and nuclei) in		endothermic alone does not score.
			the bond ✓	1	
	(b)		bonds broken = $1(H-H) + 1(F-F) = 436 + (F-F)$ and bonds		ecf possible
			made = 2(H–F) = 1136 ✓		
			436 + (F–F) – 1136 = –542 ✓		
			b and anthermore $450$ (i.e. $mat^{-1}$ )	0	
	(-)		bond enthalpy = 158 (kJ mol <sup>-1</sup> ) $\checkmark$	3	
	(c)		enthalpy change is $\Delta H$ for $\frac{1}{2}H_2 + \frac{1}{2}CI_2 \rightarrow HCI \checkmark$		
			bonds broken = $218 + 121 = 339$ and bonds made		
				2	
	(4)	:	$= -432 \Delta H = -93 \text{ (kJ mol^{-1})} \checkmark$	2	allow partially insided
	(d)	1	a weak acid is partially dissociated $\checkmark$		allow partially ionised
			to form H⁺ ions/ protons ✓	2	can be shown in an equation
		ii	the H–F bond is stronger ora ✓	1	accept harder to break the bond.
		iii	$Na_2CO_3 + 2HF \rightarrow 2NaF + CO_2 + H_2O \checkmark$	1	ignore state symbols; allow H <sub>2</sub> CO <sub>3</sub>
		iv	$\text{CO}_3^{2-} + 2\text{H}^+ \rightarrow \text{. CO}_2 + \text{H}_2\text{O}/$		ignore state symbols
			$Na_2CO_3 + 2H^+ \rightarrow 2Na^+ + CO_2 + H_2O \checkmark$	1	do not allow the inclusion of spectator ions
		v	hydrochloric acid bubbles faster / solid disappears		
			faster ✓		
			concentrations (of acids) must be the same $\checkmark$	2	allow conditions must be the same
			Total	15	

Qı	uestio	n	Expected Answers	Marks	Additional guidance
3	(a)		high temperature needed (to send equilibrium to RHS) $\checkmark$		all stand alone marks
			since (forward) reaction is endothermic $\checkmark$		
			low pressure needed (to send equilibrium to RHS) $\checkmark$		
			since fewer moles of gas on LHS $\checkmark$	4	
	(b)		high temperature (gives a fast rate of reaction) $\checkmark$		all stand alone marks
			because a higher proportion of collisions exceed $Ea \checkmark$		accept because a 'higher proportion of collisions have enough energy to cause a reaction' do not accept 'more collisions' or 'more energetic collisions'
			high pressure (gives a fast rate of reaction) $\checkmark$		
			because molecules are closer together <b>and</b> collide at a faster rate/ because particles are more concentrated <b>and</b> collide at a faster rate ✓	4	not answer based on <i>E</i> <sub>a</sub>
	(C)		would use high temperature – for rate <b>and</b> yield $\checkmark$		
			but compromise on pressure $\checkmark$		compromise must be on pressure not temperature
			because high pressure gives a fast rate but poor yield $\checkmark$	3	both rate and yield needed
	(d)	i	the (total) enthalpy change for a reaction is the same whichever route is taken ✓ owtte	1	
		ii	cycle/ $\Sigma \Delta H$ (products) - $\Sigma \Delta H$ (reactants) $\checkmark$		cycle need not be drawn correctly/ drawn at all -107 scores 3
			$-75 - 242 + 210 = \Delta H_{\rm f} \checkmark$		
			∆ <i>H</i> <sub>f</sub> = - 107 (kJmol <sup>-1</sup> ) ✓	3	
			Total	15	

Q	uestic	on Expected Answers	Marks	Additional Guidance		
4	(a)	enthalpy profile diagram				
		y axis labelled energy/enthalpy and one curve drawn to include either horizontal lines or reactants and products labelled $\checkmark$		1 max for enthalpy profile diagram if diagrams are reversed		
		second curve included to start and finish at same energy level and one curve labelled (catalysed or uncatalysed) ✓				
		Boltzmann distribution				
		axes labelled x as energy <b>and</b> y as number/ fraction or % of particles/ molecules/atoms $\checkmark$		for x axis allow kinetic energy/ KE/ velocity/ enthalpy do not allow activation energy/ $E_a$		
		shape of curve ✓		must start at 0,0 and at higher energies must not touch or cross the x axis		
		catalysed and uncatalysed $E_{\rm a}$ shown as vertical lines touching or crossing the curve and labelled $\checkmark$		if 2 graphs are drawn , this mark does not score		
		explanation				
		to increase the rate of reaction more collisions/particles/molecules have to exceed $E_a$ /have enough energy to react $\checkmark$				
		a catalyst acts by lowering $E_a \checkmark$		if candidate says catalyst gives molecule more energy, this mark		
		by allowing the reaction to proceed via a different route $\checkmark$	8	does not score		
	(b)	equilibrium position unchanged ✓				
		rate of forward and reverse reaction increased by same amount $\checkmark$	2			
		Total	10			

## 2813/03 How Far? How Fast?/Experimental Skills 1 Practical Examination

Skill P: 16 marks (out of 19 available)

#### Analysis of the acids – 8 marks

#### Four different chemical tests must be described, one for each acid.

- If the same reagent is used twice, **award zero** on the second occasion.
- If the test described is ambiguous, the first two correct bullets are not credited.
- For each test, two or three bullets correct = 1 mark: four bullets correct = 2 marks

#### 1. Test for sulphuric acid (marks A1, A2)

Use of CuO (giving blue solution) scores 1 mark (max) as test for sulphuric acid

- Add aqueous barium nitrate/chloride Aq or solution is needed, but do not penalise twice in the Plan
- White precipitate indicates that the solution tested is sulphuric acid
- The precipitate is barium sulphate (can be credited from 's' state symbol in the equation)
- Equation or ionic equation for reaction Ba(NO3)2 + H2SO4  $\rightarrow$  BaSO4 + 2HNO3

#### 2. Use of AgNO3 (marks A3, A4)

- Add aqueous silver nitrate
- Cream/off-white precipitate obtained indicates HBr or white/milky precipitate indicates HCI
- Precipitate is silver chloride/bromide (can be scored from 's' state symbol in an equation)
- Equation or ionic equation for reaction (ignore state symbols)

#### 3. Test for the other hydrohalic acid (marks A5, A6)

#### Option (a) For HBr

- Add chlorine *or* <u>concentrated</u> sulphuric acid
- Goes orange (or appropriate observation)
   Do not allow brown as the colour of aq bromine
   Accept any reasonable colour if an organic solvent is used to show Br2 colour
- Chlorine is more reactive than bromine *or* chlorine displaces bromine (*owtte*)
- Equation *or* ionic equation for reaction

[2]

[2]

[2]

#### Option (b) For either HCl or HBr

- [Add silver nitrate] then add ammonia to the precipitate formed
- Use excess aqueous ammonia and shake the tube
- All three points required (excess, aqueous/solution, invert/shake/stir)
- For HCI, the precipitate dissolves/ goes clear
- For HBr, the precipitate is insoluble/ partially soluble/ does not dissolve or precipitate is soluble in concentrated ammonia

#### 4. Test for ethanoic acid (marks A7, A8)

## [2]

#### Option (a)

- Add magnesium, a named metal carbonate or NaHCO<sub>3</sub>
- Fizzing, bubbling *or* effervescence observed
- Test for gas outlined or reference to the slow reaction [compared to the other acids]
- Equation *or* ionic equation for reaction

#### Option (b)

- Add any named alcohol
- Heat with concentrated sulphuric acid
- [Sweet/fruity] smell of ester obtained
- Equation for chosen reaction Accept molecular formulae, but the -COO- ester linkage must be shown

#### Option (c) - 1 mark available, only, for these tests

	<ul><li>Add a named indicator</li><li>Correct final colour of indicator</li></ul>	<ul> <li>or add PCI<sub>5</sub> (or similar reagent)</li> <li>or steamy/ acidic fumes produced</li> </ul>	<i>or</i> aq FeCl₃ <i>or</i> red colour
I	Titration procedure – 7 marks		
T1	Makes up accurately a <b>known solutio</b> Detail needed: weighed mass of KOH		[1]
T2	Quantitative <b>dilution</b> of acid chosen us <i>The dilution must make the concentrat</i>		sk <b>[1]</b>
Т3	Correct equation for neutralisation rea	action used	[1]
T4	Uses equation to <b>justify mass of alka</b> <i>T4 cannot be awarded if the acid used</i>		[1]
Т5	Outline description of use of burette an	nd pipette in titration procedure	[1]
Т6	At least two <b>consistent titres</b> (or withi <b>and</b> suitable <b>indicator</b> chosen <b>and</b> co	,	[1]
Τ7	<b>Specimen calculation</b> to determine $\%$ 56 (or 56.1) for the $M_r$ of KOH must be		[1]

#### S Safety, Sources and QWC - 4 marks

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#### S1 Safety: Solid/pure potassium hydroxide is corrosive/ burns the skin: wear gloves or face mask or if spilt on skin wash with plenty of water [1] Hazard and precaution are both required No credit for references to hazard of the acids

#### S2 Two sources quoted in the text or at end of Plan.

S3 **QWC**: text is legible and spelling, punctuation and grammar are accurate [1] Accept no more than five different errors in legibility, spelling, punctuation or grammar.

#### S4 **QWC**: information is organised clearly and coherently

- Is a word count given and within the limits 450 1050 words?
- Is scientific language used correctly and are chemical formulae correct? •
- Are the four tests described in a logical sequence without excessive repetition?

#### Practical Test (B)

#### Part 1 [14 marks] Presentation of titration data

#### Six bullets correct = 2 marks: five bullets correct = 1 mark

- Table grid drawn (at least three lines) and all burette data is shown in the table, including • first/trial.
- Correctly labelled table (initial, final and difference aw) for burette data
- Three (or more) titres are shown
- All "accurate" burette data and subtractions are quoted to two decimal places, ending in .00 or .05
- No readings recorded above 50 cm<sup>3</sup>
- All subtractions are correct (these must be checked and indicated by a dot) Two subtraction errors count as two bullet errors, ie scores 0 for presentation

#### Self-consistency of titres

#### Four bullets correct = 2 marks: three bullets correct = 1 mark

- The titres for any two accurate experiments are within 0.20 cm<sup>3</sup>. For this bullet, assume that the first reading is a trial, whether labelled as such or not.
- The ticked titres (or the titres used to calculate the mean) are all within 0.10 cm<sup>3</sup> If three titres are ticked, award the self-consistency on the spread of **all** three.
- Two or three titres are ticked
- **Units**, cm<sup>3</sup> or ml, must be given somewhere (*once in the table is sufficient*).

#### Mean titre correctly calculated

- The mean should normally be calculated using the **closest** two accurate titres. However, a candidate may use the trial/first reading if appropriate, without penalty.
- The mean must be quoted to 1 or 2 d.p (but **not** to 3 d.p. except for 0.025 and 0.075). "Rounding up" must be done correctly if it is needed.

[1]

[2]

[2]

[1]

[1]

[7]

#### Accuracy

The supervisor's mean titre is rounded to nearest 0.05  $cm^3$ 

Compare the mean supervisor's titre with the candidate's mean titre (T).

Put " $\delta = \_$ " on the script to show the difference between these two titres.

Use the chart below to award the mark out of 7 for accuracy.

<i>T</i> is within <b>0.25</b> cm <sup>3</sup> of mean supervisor's value	[7 marks]
<i>T</i> is within <b>0.40</b> cm <sup>3</sup> of mean supervisor's value	[6]
<i>T</i> is within <b>0.60</b> cm <sup>3</sup> of mean supervisor's value	[5]
<i>T</i> is within <b>0.80</b> cm <sup>3</sup> of mean supervisor's value	[4]
<i>T</i> is within <b>1.00</b> cm <sup>3</sup> of mean supervisor's value	[3]
<i>T</i> is within <b>1.20</b> cm <sup>3</sup> of mean supervisor's value	[2]
<i>T</i> is within <b>1.50</b> cm <sup>3</sup> of mean supervisor's value	[1 mark]

#### Spread penalty

("Spread" is defined by the titres used by the candidate to calculate his/her mean) If these titres have a spread =  $0.40 \text{ cm}^3$  or more, deduct 1 accuracy mark. If these titres have a spread =  $0.60 \text{ cm}^3$  or more, deduct 2 marks. If these titres have a spread =  $1.00 \text{ cm}^3$  or more, deduct 3 marks (max spread penalty)

#### <u>Safety</u>

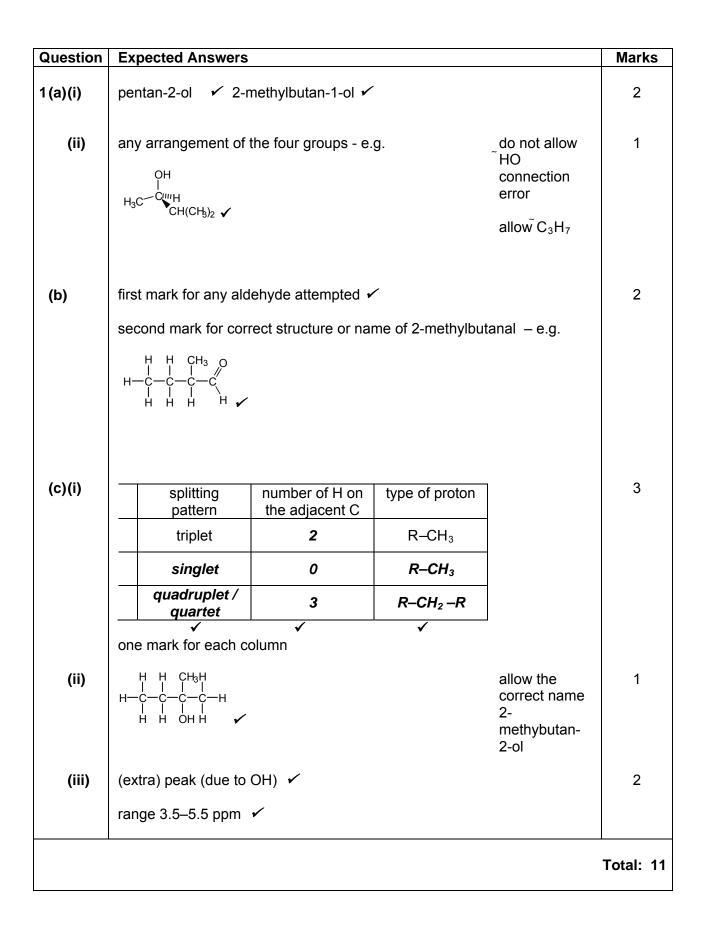
Dilut	ed <b>W</b> will have <u>no</u> hazard symbol	[1]	
or ad	ing the acid/ adding water reduces the [level of] hazard dding water makes the acid safer ( <i>owtte</i> ) <i>second mark is independent of the first</i>	[1]	
Part 2 [10 marks]			
	k <b>ecf</b> from one part of the question to the next, but <b>not</b> within a part. In a mark is awarded for the answer, it must be quoted to 3 significant figures		
(a)	Dilution factor was ${}^{250}/{}_{10}$ = 25, so concentration = ${}^{1.00}/{}_{25}$ = 0.0400 <i>Minimum "proof" needed is figures 1.0, 10 and 250 suitably used.</i>	[1]	
(b)	Answer <b>(a)</b> is multiplied by <sup>mean titre</sup> / <sub>1000</sub> This is a method mark; check that the candidate's mean titre was actually used	[1]	
	Correct answer was obtained from candidate's own mean titre	[1]	
(c)	$Ca(OH)_2(aq) + 2HX(aq) \rightarrow CaX_2(aq) + 2H_2O(I)$	[1]	
	State symbols are correct (mark is conditional on $H_2O$ shown as by-product)	[1]	
(d)	Answer <b>(b)</b> is multiplied by 0.5	[1]	

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(e)	Answer (d) is multiplied by 40 ( $or^{1000}/_{25}$ )	[1]
	[Ca(OH) <sub>2</sub> ] correctly calculated Candidate's answer should be 20 x <b>(b)</b> unless ecf has been applied	[1]
(f)	<b>2 marks</b> $M_r$ of calcium hydroxide = 74.1	[1]
	Mass = answer (e) x 74.1 Use of 74 as the $M_r$ can score 1 mark out of 2 in (f)	[1]
Part	3	[6 marks]
(a)	White/milky precipitate/solid formed	[1]
(b)	Precipitate disappears <i>or</i> solid dissolves <i>or</i> colourless solution formed <i>or</i> [mixture] goes clear	[1]
(c)	Identity: HX is HCI or hydrochloric acid	[1]
	<b>Reason:</b> Chlorides give a white precipitate with AgNO <sub>3</sub> /Ag <sup>+</sup> , soluble in a Reference to correct evidence from <b>both</b> tests is required for this mark	ammonia [1]
(d)	AgNO <sub>3</sub> (aq) + HX ( <i>or</i> HCI)(aq) → AgX ( <i>or</i> AgCI)(s) + HNO <sub>3</sub> (aq) <i>or</i> Ag <sup>+</sup> (aq) + Cl <sup>-</sup> (aq) → AgCl(s) For getting all formulae correct <i>Mark ecf if HX was identified as HBr or HI</i>	[1]
	For absence of any balancing figures and all state symbols correct	[1]
Part	4: (Skill E)	[14 marks]
(a)(i	) To ensure that <u>all</u> <b>HX</b> reacts	[1]
(a)(i	i)Number of moles of HX used = 0.01(0) mol	[1]
	Number of moles of Mg used = $^{0.15}/_{24.3}$ = 0.0062 mol	[1]
	Use of 2:1 mole reacting ratio to show clearly that Mg is in excess <i>There are alternative methods of doing this calculation.</i>	[1]
(a)(i	ii) There would be solid/magnesium left in the tube at the end of the expe	riment. [1]
(b)	No of moles of HX used = 0.01	[1]
	Volume of gas produced = $0.005 \times 24000 = 120 \text{ cm}^3$	[1]
	This volume exceeds the capacity of the cylinder, so too much acid was <i>Maximum 1 mark awarded for the calculation if 1:1 mole ratio is used There are alternative methods of doing this calculation.</i>	used [1]

(c)(i) 4 marks (Mark the best two strands given by the candidate)

C1	Gas would escape before the bung could be fitted	[1]
C2	Put one reagent into an ignition tube inside the flask (or similar precaution)	[1]
D1	Reaction might still be taking place [after one minute]	[1]
D2	Take final reading when fizzing/bubbling stops <i>or</i> wait until the volume of gas [in cylinder] stops increasing	[1]
E1	Use a burette instead of a measuring cylinder for gas collection Allow use of a gas syringe for E1, but this improvement cannot score E2	[1]
E2	Burette is more accurately calibrated/manufactured <b>or</b> has narrower stem <b>or</b> burette can be read to 0.05/0.1 cm <sup>3</sup> [whereas measuring cylinder to 1 cm <sup>3</sup> ]	[1]
F1	Remove oxide layer on Mg before use by scraping/ sandpapering	[1]
F2	The MgO would react with the acid without producing any gas	[1]
(c)(ii)	Gas collection experiment was carried out only once / was not repeated	[1]
	Titration was repeated to obtain consistent results/titres so it is [more] reliable The reliability must be clearly linked to the consistency (not to accuracy)	[1]

## 2814 Chains, Rings and Spectroscopy



Question	Expected Answers		Marks
2(a)(i)	orange solid / ppt (when added to) 2,4-DNPH / Brady's Reagent ✓		
(ii)	(when warmed) with ammoniacal silver nitrate (solution) / Tollens' Reagent ✓ no silver (mirror) / reaction ✓	or no green colour with acidified dichromate	1
		allow 'no reaction' only if a suitable reagent used	2
(iii)	H OH $H N O\delta^{-}$ $\delta^{+}$ $\delta^{+}$ $\delta^{+}$		3
	dipole on C=O ✓ curly arrow breaking bond ✓ curly arrow from lone pair on N in the direction of	the carbonyl C ✓	
(iv)	$M_{\rm r}$ caprolactam = 113 $\checkmark$		3
	mass of cyclohexanone needed = $98 / M_r = 0.867$	7 tonnes or ecf $\checkmark$	
scale for 99% yield = mass * 100/99 = <b>0.88</b> tonnes or ecf ✓ (or any number rounding to 0.876 tonnes)			
(b) a single molecule with six carbon atoms in a straight chain $\checkmark$		ght chain 🖌	2
	correct end groups on a straight chain ✓- e.g.		
	H <sub>2</sub> N COOH / H <sub>2</sub> N	COCI	
			Total: 11

Question	Expected Answers	Marks
6(a)(i)	same structure / displayed formula / order of bonds different 3-D / spatial arrangement </td <td>1</td>	1
(ii)	но сно	2
	chiral centre circled  C=C double bond circled	
(b)(i)	substitution: phenol / benzene / arene ✓(allow C=C, but not just 'double bond')	2
(ii)	$C_{17}H_{20}O_5 + 4[H] \longrightarrow C_{17}H_{24}O_5 \checkmark$ (or correct structure)	1
(iii)	correct products of hydrolysis (in acid or alkali conditions) $\checkmark \checkmark \checkmark$ $Na^+ = 0$ $Na^+ = 0$ $Na^+ = 0$ CHO $Na^+$ is not essential,	3
	ionisation of phenolic group <b>and</b> carboxylate group for alkali conditions ✓ but do not allow covalent Na-O	
(c)(i)	304 🗸	1
(ii)	1000–1300 (cm <sup>-1</sup> ) due to C–O ✓ 1680–1750 (cm <sup>-1</sup> ) due to C=O ✓ 3230–3550 (cm <sup>-1</sup> ) due to O-H ✓	3
(d)	only one (stereo)isomer is active / effective / needed AW 🖌	2
	natural oil contains one optical/stereoisomer isomer and lab synthesis contains a mixture ✓ AW	

Question	Expected Answers	Mark	S
4(a)(i)	LiAlH <sub>4</sub> (in ether) / Na in ethanol $\checkmark$	1	
(ii)	reduction/hydrogenation /redox 🖌 🛛 allow addition	1	
(b)(i)	$C_2H_5NH_3^+$ Cl + and - not essential, but	both or 1	
(ii)	$C_2H_5NH_3^+ CH_3COO \checkmark$ neither must be present	1	
(c)	5	be 4 blied	
	lone pair on N is (partially) delocalised / incorporated into refe	ore any rence to inductive	
	ethviamine	ct for nylamine)	
	comparison to explain relative basicity discussion of negative charge / electron density on the nitrogen ✓		
	discussion of the ease of donation/availability of the lone pair of electrons on the N ✓ any 4 out of 5 marks		
QWC	mark for well organised response with correct use of <b>one</b> of the t <b>inductive effect</b> , <b>delocalised, mesomeric</b>	erms: 1	
(d)	correct formula of 4-nitrobenzoic acid as starting material 🖌	6	
	correct formula of 4-aminobenzoic acid or ethyl-4-nitrobenzoate	✓	
	reduction of nitro group (reflux with) Sn + (conc) HCl ✓		
	$(1 \text{ Cond} \text{ With}) \text{ Cond} + (2 \text{ Cond}) \text{ Cond} + 6[\text{H}] \longrightarrow \text{H}_2\text{N} - (2 \text{ Cond} + 2 \text{ Cond}$	2H <sub>2</sub> O	
1	esterification	✓	
	(reflux / distil / heat with) ethanol + (conc.) $H_2SO_4$		
	$H_2N$ COOH + $C_2H_5OH$ $\rightarrow$ $H_2N$ COOC <sub>2</sub> $H_5$ +	- H <sub>2</sub> O ✓	
<i>.</i>	(steps can be in either order)		
(e)			
	$\checkmark$	2	
		Total:	

Question	Expected Answers	Marks
5(a)(i)	addition 🖌	1
(ii)	$\mathbf{n}_{H} \stackrel{H}{\underset{H}{\longrightarrow}} \mathbf{n}_{H} \stackrel{H}{\underset{H}{\longrightarrow}} \mathbf{n}_{H} \stackrel{H}{\underset{H}{\longrightarrow}} \mathbf{n}_{H} = \mathbf{n}_{H} \stackrel{H}{\underset{H}{\longrightarrow}} \mathbf{n}_{H} \stackrel{H}{\underset{H}{\underset{H}{\longrightarrow}} \mathbf{n}_{H} \stackrel{H}{\underset{H}{\underset{H}{\underset{H}{\longrightarrow}} \mathbf{n}_{H} \stackrel{H}{\underset{H}{\underset{H}{\underset{H}{\longrightarrow}} \mathbf{n}_{H} \stackrel{H}{\underset{H}{\underset{H}$	2
(iii)	allow ecf from (ii)	2
	benzene rings on every second carbon in random directions $\checkmark$	
	diagram also shows correct use of 3-D bonds $\checkmark$	
(iv)	isotactic has side chains on the same side AW $\checkmark$	1
(b)	$H_2N \longrightarrow NH_2 H_2N \longrightarrow SO_3 Na^{+} \checkmark N^{\parallel} N^{\parallel} SO_3 Na^{+} \checkmark$	5
	add the amine to HCI + NaNO <sub>2</sub> /HNO <sub>2</sub> $\checkmark$	
	temp < 10 °C ✓	
	add to alkaline phenol 🖌	
Total:		otal: 11

Question	Expected Answers	Marks
6(a)(i)		1
(ii)	$\begin{array}{rrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrr$	2
(b)	ester eg $C_6H_5COCI + C_2H_5OH \longrightarrow C_6H_5COOC_2H_5 + HCIalcohol / correct name/formula of a suitable example \checkmarkequation \checkmark$	4
	amide eg $C_6H_5COCI + C_2H_5NH_2 \longrightarrow C_6H_5CONHC_2H_5 + HCI$ $C_6H_5COCI + 2NH_3 \longrightarrow C_6H_5CONH_2 + NH_4CI$	
	amine / ammonia / correct name/formula of a suitable example 🖌 equation 🖌	
(c) (i)	FeCl <sub>3</sub> / AICl <sub>3</sub>	1
(ii)	electrophilic substitution	1
		Total: 9

Question	Expected Answers	Marks
7(a)	<b>cyclohexene</b> (electrophilic) addition $\checkmark$ C <sub>6</sub> H <sub>10</sub> + Br <sub>2</sub> $\longrightarrow$ C <sub>6</sub> H <sub>10</sub> Br <sub>2</sub> $\checkmark$	8
	<b>benzene</b> (electrophilic) substitution $\checkmark$ $C_6H_6 + Br_2 \longrightarrow C_6H_5Br + HBr \checkmark$ needs a / halogen carrier / catalyst/ Fe / FeBr <sub>3</sub> / AlBr <sub>3</sub> etc $\checkmark$	
	relative reactivites benzene is: (more) stable / less reactive / less susceptible to electrophiles ✓ AW ora	
	( )-electron electron density in benzene is low ora $\checkmark$	
	comparison of the relative ability to attract /polarise the electrophile or relative ability to donate an electron pair $\checkmark$	
	<b>QWC</b> mark for at least two sentences/bullet points in context with correct spelling, punctuation and grammar	1
(b)(i)	$^{30.5}/_{12.0} = ^{1.7}/_{1.0} = ^{67.8}/_{79.9} =$	2
	2.54 1.7 0.85 🗸	
	3 : 2 : 1	
	empirical formula = $C_3H_2Br$	
(ii)	( $M_r$ of C <sub>3</sub> H <sub>2</sub> Br = 117.9 = ${}^{235.8}/_{2}$ , so) molecular formula = C <sub>6</sub> H <sub>4</sub> Br <sub>2</sub> $\checkmark$	3
	Br Br Br Br Br Br	
	one correct structure 🖌 all three correct 🖌	
(c)	any valid structure for 1,2,3,4,5,6-hexabromocyclohexane – e.g.	2
	Br Br Br	
	$C_6H_6 + 3Br_2 \longrightarrow C_6H_6Br_6 \checkmark$	
		otal: 16
	Т	otal: 10

# 2815/01 Trends and Patterns

C	Question				Marks	Additional Guidance
1	(a)		1s <sup>2</sup> 2s <sup>2</sup> 2p <sup>6</sup> 3s <sup>2</sup> 3p <sup>6</sup> 3d <sup>9</sup> (1)	1		
	(b)	(i)	Correct formula of a copper(II) complex ion e.g. $CuCl_4^{2^-} / [Cu(NH_3)_4(H_2O)_2]^{2^+} / [Cu(H_2O)_6]^{2^+} (1)$	1		
		(ii)	Correct colour (1) e.g. $CuCl_4^2$ green/yellow, $[Cu(NH_3)_4(H_2O)_2]^{2^+}$ dark blue and $[Cu(H_2O)_6]^{2^+}$ blue	1	Allow ecf from a known copper compound	
		(iii)	Coordinate bond / dative bond (1) Lone pair donated by ligand / lone pair accepted by copper (1)	2		
	(c)	(i)	Blue precipitate / blue solid	1	Can get credit for ppt from state symbol of correct product in part (ii)	
		(ii)	$Cu^{2^+} + 2OH^- \rightarrow Cu(OH)_2 / [Cu(H_2O)_6]^{2^+} + 2OH^-$ → $Cu(OH)_2 + 6H_2O / [Cu(H_2O)_6]^{2^+} + 2OH^- \rightarrow$ $Cu(H_2O)_4(OH)_2 + 2H_2O (1)$	1	Allow correct multiples Ignore state symbol	
	(d)	(i)	Mole ratio C:Cu:K:N = 0.0320:0.00800:0.0240:0.0320 (1) K <sub>3</sub> CuC <sub>4</sub> N <sub>4</sub> (1)	2	Allow the four masses ÷ appropriate A <sub>r</sub> if mole ratio not calculated Allow any order of atoms Can award formula mark if given in part (ii) Allow ecf from wrong mole ratio	
		(ii)	[Cu(CN) <sub>4</sub> ] <sup>3-</sup> / CuC <sub>4</sub> N <sub>4</sub> <sup>3-</sup> (1)	1	Allow any order of atoms with or without brackets Allow ecf from wrong formula	
			Total	10		

Q	luesti	on	Expected Answers	Marks	Additional Guidance
2	(a)	(i)	Magnesium 2.8 or diagram with 8 or 0 electrons in outer shell and two chloride ions of 2.8.8 or diagram with 8 electrons in outer shell (1) Correct charge Mg <sup>2+</sup> and Cl <sup>-</sup> (1) Mark independently from dot and cross	2	Allow one mark for a correct magnesium ion with charge / two correct chloride ions with charge <b>Not</b> $[Cl_2]^{2-}$
		(ii)	Giant structure (1) Strong attraction between ions / lots of energy needed to overcome attraction between ions / lots of energy needed to break ionic bonds (1)	2	Allow strong ionic bonds
	(b)		Si + $2Cl_2 \rightarrow SiCl_4$ (1)	1	Allow any correct multiples Ignore state symbols
	(c)		Correct outer shell of phosphorus (five bond pairs) (1) Rest of structure correct (1) <b>independent</b> of outer shell of phosphorus mark so that PCl <sub>4</sub> , PCl <sub>3</sub> or PCl <sub>6</sub> could get this mark if all chlorine atoms are correct	2	Rest of structure of mark
	(d)	(i)	PCI <sub>5</sub> + 4H <sub>2</sub> O → H <sub>3</sub> PO <sub>4</sub> + 5HCI / PCI <sub>5</sub> + H <sub>2</sub> O → POCI <sub>3</sub> + 2HCI (1)	1	Allow any correct multiples Allow H <sup>+</sup> and Cl <sup>-</sup> instead of HCl Ignore state symbols
		(ii)	Hydrolysis (1)	1	
	(e)		Covalent (1)	1	Allow simple molecule / dative bond / has considerable covalent character
	(f)	(i)	(orange, brown, rusty or yellow) to (blood) red (1)	1	Allow (deep) red Not green / blue to red Not precipitate
		(ii)	Ligand substitution / ligand exchange (1)	1	Allow ligand displacement / ligand replacement
			Total	12	

Question	Expected Answers	Marks	Additional Guidance
3 (a)	$2Cr^{3+} + 3H_2O_2 + 10OH^- \rightarrow 2CrO_4^{2-} + 8H_2O$ Correct reactants and products (allow e <sup>-</sup> and OH <sup>-</sup> on <b>both</b> left and right) and correct molar ratio of $Cr^{3+}$ and $H_2O_2$ (1); Balanced (1)	2	For the second mark the OH <sup>-</sup> and e <sup>-</sup> must be cancelled down
(b)	Moles $MnO_4^- = 0.000463 (1)$ Moles $Fe^{2^+} = 5 \times moles MnO_4^- / 0.002315 (1)$	4	Allow ecf within the question
	<i>M</i> <sub>r</sub> = 392 / 391.8 (1)		ecf is 0.907 ÷ moles of Fe <sup>2+</sup> <b>Allow</b> three marks for 392 / 391.8 with no working
	$x = 6$ (1) <b>dependent</b> on $M_r$ given		ecf is ( <i>M</i> <sub>r</sub> – 283.8) ÷ 18 <b>Allow</b> one mark for 6 with no working
	Total	6	

(	Question		Expected Answers	Marks	Additional Guidance
4	(a)	(i)	Any three from Magnesium ion smaller than barium ion / magnesium ion has a higher charge density / ora (1);	3	Must use correct particle but only penalise once
			Magnesium ion is more polarising / ora (1): Magnesium ion distorts the carbonate ion more than barium ion / ora (1); So carbon–oxygen bond (in carbonate) is weaker / so covalent bond in carbonate is weaker (1)		Assume carbonate refers to an ion Allow CO <sub>3</sub> <sup>-</sup> Only need one comparative statement in the last three marking points
		(ii)	Contains Na <sup>+</sup> which has a low charge density / contains Na <sup>+</sup> which is not very polarising (1)	1	Ignore Na <sup>+</sup> is a large ion Allow Na <sup>+</sup> is a larger ion than Mg <sup>2+</sup> Allow Na <sup>+</sup> has a smaller charge than Mg <sup>2+</sup>

Question	Expected Answers	Marks	Additional Guidance
4 (b)	Definitions $2Na^{+} + O^{2^{-}} \rightarrow Na_2O(1)$ $2Na + \frac{1}{2}O_2 \rightarrow Na_2O(1)$	3	If given state symbols must be <b>correct</b>
	Lattice enthalpy is the enthalpy change when one mole (of ionic solid) is made from its constituent gaseous ions but formation from its constituent elements (1)		Allow energy released Not energy absorbed Allow states from equations
	Born-Haber cycle Correct state symbols for the formulae given (1);	6	Formula must have correct state symbol at least once in the cycle
	Correct formula (1);		Formulae given must be correct but there can be a formula missing
	Labelling of enthalpy changes Three correct labels (1) but five correct labels (2) but all labels correct (3)		Allow ecf from the cycle drawn Allow conventional symbols e.g. $\Delta H_{\rm f}$
	Expression or statement in words or symbols to show how lattice enthalpy is calculated (1)		Allow ecf from cycle drawn

Question	Expected Answers	Marks	Additional Guidance
4	Lattice enthalpy magnesium oxide, sodium chloride and potassium bromide (1) Any two from Comparison of charge density or ionic radius of cation (1) e.g. ionic radius decreases from K <sup>+</sup> , Na <sup>+</sup> to Mg <sup>2+</sup> / charge density increases from K <sup>+</sup> , Na <sup>+</sup> to Mg <sup>2+</sup> (1) Comparison of charge density or ionic radius of anion (1) e.g. e.g. ionic radius decreases from Br <sup>-</sup> , Cl <sup>-</sup> to O <sup>2-</sup> / charge density increases from Br <sup>-</sup> , Cl <sup>-</sup> to O <sup>2-</sup> (1) Comparison of charge on ions (1) Na <sup>+</sup> but Mg <sup>2+</sup> / O <sup>2-</sup> but Cl <sup>-</sup>	3	<b>Allow</b> Mg <sup>2+</sup> but Na <sup>+</sup> (1) and Na <sup>+</sup> is smaller than K <sup>+</sup> (1) <b>Allow</b> O <sup>2-</sup> but Cl <sup>-</sup> (1) and Cl <sup>-</sup> is smaller than Br <sup>-</sup> (1)
	Quality of Written Communication (1) At least two complete sentences with correct spelling, punctuation and grammar that address the question set	1	
	Total	17	

# 2815/02 Biochemistry

Question	Expected Answers	Mark
1 (a)(i)	Diagram must include –CO–NH- (or displayed version)√ Amide/peptide link ✓	2
(ii)	Phenylalanine and alanine. $\checkmark$ Any logical reasoning eg a dipeptide has general formula H <sub>2</sub> NCHRCONHCHRCOOH which leaves C <sub>7</sub> H <sub>8</sub> – can only be those two. $\checkmark$	2
(b)	<ul> <li>Any three examples from the following four using the amino acids provided in each case @ 2 marks each.</li> <li>Ionic(accept electrostatic) ✓ between COO<sup>-</sup> and NH<sub>3</sub><sup>+</sup> ✓</li> <li>van der Waals/Instantaneous dipole–induced dipole ✓ between any two sidechains from alanine and phenylalanine ✓ Not 'hydrophobic bonding'.</li> <li>Covalent or disulphide or C-S-S-C ✓ between two cysteines. ✓</li> <li>Accept hydrogen bonding ✓ between two suitable groups</li> </ul>	6
	<ul> <li>such as COOH and NH<sub>2</sub> ✓ as long as these groups are not ionised.</li> <li>Where candidate uses both ionised, for ionic, and non-ionised, for hydrogen bonding allow both bonding type marks but only one example mark – max 3 marks from 4.</li> </ul>	1
	Diagrams, if given, must be correct.	
	QWC. Award the mark for correct reference to three types of interaction, and one appropriate amino acids (by name or diagram)	

Question	Expected Answers	Mark
2 (a)(i)	Only one optical isomer of tyrosine is the right shape to fit in the active site AW. $\checkmark$	1
(ii)	COOH becomes COO and NH <sub>2</sub> becomes NH <sub>3</sub> <sup>+</sup> and no change to phenol $\checkmark$	1
(iii)	$\frac{\rm NH_{3}^{\pm} group \ becomes \ uncharged \ (\rm NH_{2}), \ losing \ \underline{ionic} \ attraction. \checkmark}{\rm No \ mark \ for \ protonating \ COO-}$	1
(b)(i)	Binds/ <b>fits</b> to active site <b>reversibly/</b> instead of tyrosine✓ AW	1
(ii)	initial rate initial rate tyrosine concentration One mark for each curve. ✓ Labels wrong way round on correct graphs, max 1 mark.	2

Question	Expected Answers	Mark
3 (a)	CTATGC ✓	1
(b)	Three each $\checkmark$ of AT and CG links. (3 x 2) + (3 x 3) = 15 $\checkmark$ Correct answer(15) earns both marks	2
(c)	CUAUGC ✓	1
(c) (d)	<ul> <li>CUAUGC ✓</li> <li>Six points from:</li> <li>Base triplet on m-RNA attract t-RNA molecule with the complementary triplet ✓ (Not codon etc without reference somewhere to base triplets)</li> <li>Hydrogen bond forms between t-RNA and m-RNA ✓</li> <li>Each t-RNA carries its appropriate amino acid AW✓</li> <li>Two t-RNA molecules are attached to m-RNA at a time ✓</li> <li>Bond between t-RNA and its amino acid breaks ✓</li> <li>The amino acid is then joined, to the growing peptide/polypeptide chain by a peptide/amide/C-N link ✓</li> <li>by a condensation reaction / mention of peptidyl transferase</li> <li>Hydrogen bonds between the t-RNA and m-RNA break. ✓</li> <li>The pair UA has two hydrogen bonds and CG three ✓</li> <li>Start codon is AUG/ for methionine ✓</li> <li>Marks may be found on a suitable diagram.</li> </ul>	6

Question	Expected Answers	Mark
4 (a)(i)	( $\alpha$ ) means that the O on the right hand carbon (C1) points	1
	down/is below plane of ring✓. (pyranose) means a ring of five carbon atoms and one oxygen.✓AW	1
(ii)	HOCH₂CHOHCHOHCHOHCHOHCHO or fully displayed. Ignore stereochemistry.✓	1
(b) (i)	$\begin{array}{c ccccc} CH_2OH & CH_2OH \\ H & C & O & H & H & C & O & H \\ \hline & & & & & & & \\ H & & & & & & \\ C & & & & & & & \\ H & & & & & & \\ H & & & &$	
	<ul> <li>✓ for glycosidic link C-O-C (which may be skeletal)</li> <li>✓ for the rest including stereochemistry</li> </ul>	2
(ii)		
(c)(i)	maltose✓	1
	<ul> <li>Because the necessary OH groups are not available o hydrogen bond to water ✓ being</li> <li>used in glycosidic links ✓</li> <li>used in hydrogen bonding between chains. ✓</li> <li>Any two points</li> </ul>	2
(ii)	<ul> <li>Any three points from:</li> <li>A linear polymer (not helical)√</li> <li>The polymer molecules lie alongside each other and hydrogen bond across√</li> <li>Diagram or explanation of the hydrogen bonds involved O-H···O √</li> <li>The molecules/ fibrils form bundles of fibres which have tensile strength √</li> </ul>	3

#### 2815/02

Questio	n	Expected Answers	Mark
5	(a)	Three 🗸	1
	(b)(i)	$ \begin{array}{c}                                     $	2
		$H_{2}C \longrightarrow O \longrightarrow C_{17}H_{29}  \checkmark \text{ for ester group } \checkmark \text{ for rest}$	
	(ii)	alkene/C=C ✓ and ester✓ (not ecf from (i))	2
	(c)	Hydrocarbon chains in triglyceride molecules $\checkmark$ Can form van der Waals $\checkmark$ forces with non polar solvents.	2
		Allow the marks for arguments based on the van der Waals forces being similar in separate triglyceride and solvent, and in a mixture of the two. This presents no energy barrier to mixing.	

# 2815/04 Methods of Analysis and Detection

Questio	n	Expected Answers	Marks
1(a)	(i)	$C_3H_7^{79}Br^+$ /CH <sub>3</sub> CH <sup>79</sup> BrCH <sub>3</sub> ✓ C <sub>3</sub> H <sub>7</sub> <sup>81</sup> Br <sup>+</sup> /CH <sub>3</sub> CH <sup>81</sup> BrCH <sub>3</sub> ✓	2
	(ii)	1:1 ✓	1
(b)		29 /43/ 93 / 95 🗸	1
		$H_{c} H_{b} H_{a} / H_{c} H_{b} H_{a} / H_{c} H_{b} H_{a} / H_{c} H_{b} H_{a} / H_{c} H_{c} H_{b} H_{a} / H_{c} H_{c} H_{b} H_{a} / H_{c} H_{c} H_{b} H_{a} / H_{c} H_{c} H_{b} H_{a} / H_{c} H_{c} H_{c} H_{b} H_{a} / H_{c} H_{c} H_{c} H_{c} H_{b} H_{a} / H_{c} $	2
	(ii)	Adjacent C has 2 Hs/protons/next to a CH₂ ✓	1
	(iii)	number of peaks = 2 $\checkmark$ relative peak areas = 6 : 1 $\checkmark$	2
		Total	9

Question	Expected Answers	Marks
2 (a)	<ul> <li>(electrons excited to higher energy levels) and fall back to lower</li> <li>levels ✓</li> <li>&amp; emit energy/light ✓</li> </ul>	2
(b)	excited electrons fall back to different energy levels / in a series, each electron falls back to the same energy level $\checkmark$	1
(c)		2
	n =2B	
	n = 1	
	1 mark for each transition. If arrows point upwards, credit one mark only	
(d)	uses correct formulae $f = c/\lambda$ and $E = hf$ or $E = hc/\lambda$ $\checkmark$	4
	$f = 3.29 \times 10^{15}$ (calculator value 3.289473684 x 10 <sup>15</sup> ) $\checkmark$	
	$E = h \times f = 2.18 \times 10^{-18} J / 2.18 \times 10^{-21} kJ \text{ (calculator value 2.180921053 \times 10^{-18})}$	
	1313 / 1310 / 1312.9 🗸	
	(allow ecf throughout the calculation)	
	Total	9

Ques	stion	Expected Answers	Marks
3(a)	(i)	$R_{\rm f} = {\rm distance \ moved \ by \ solute/spot} \qquad \checkmark$ distance moved by solvent	1
	(ii)	they have similar $R_{\rm f}$ values/similar solubilities (AW) $\checkmark$	1
	(iii)	run another chromatogram with a different solvent/use two way chromatography ✓	2
		because it is unlikely that two substances have similar $R_{\rm f}$ values in two different solvents $\checkmark$	
	(iv)	distribution/relative solubility of the solute/spot/component / differences in solubility ✓	2
		between the mobile and the stationary phase $\checkmark$	
(b)		uses (an applied) voltage $\checkmark$ movement depends on mass & charge $\checkmark$ pH controlled by buffer $\checkmark$ pH must be controlled because charge changes with pH $\checkmark$ example of how charge changes by showing cation (-NH <sub>3</sub> <sup>+</sup> ) at low pH or anion (-COO <sup>-</sup> ) at high pH $\checkmark$ pH changes may affect direction of movement of amino acids $\checkmark$ <b>6</b> marking points - maximum = <b>5</b>	5
014/0			1
QWC		At least two sentences that shows legible text with accurate spelling, punctuation and grammar so that the meaning is clear.	1
		Total	12

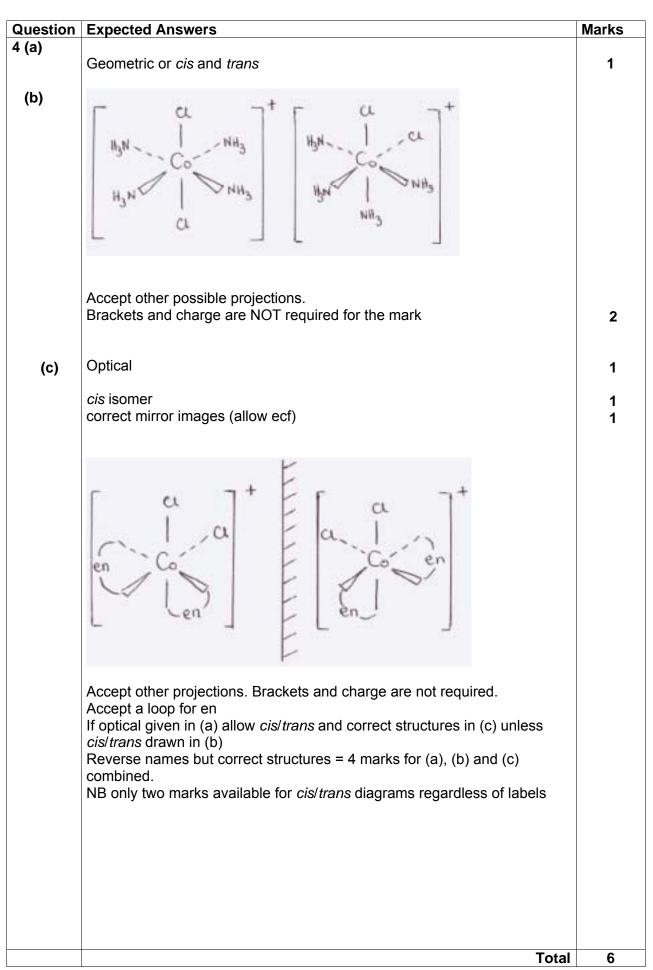
Question	Expected Answers	Marks
4(a)	$CH_3CH_2NH_2$ yes due to lone pair on N $\checkmark$	4
	$C_6H_5CH_3$ yes due to delocalised benzene ring $\checkmark$	
	$CH_3CH_2CH_3$ no because contains no lone pairs or multiple bonding/is a saturated molecule $\checkmark$	
	$CH_3CH_2CN$ yes due to <u>CEN/ unsaturated/<math>\pi</math>-bonding</u> $\checkmark$	
(b) (i)	more conjugation / larger chromophore $\checkmark$ decreases energy gap $\checkmark$ shifts absorption (to longer $\lambda$ ) into visible region $\checkmark$	3
(ii)	likely to absorb in UV region ✓	1
	Total	8

Question	Expected Answers	Marks
5(a) (i)	$M$ = 164 (from mass spec) $\checkmark$	2
	empirical mass of $C_5H_6O$ = 82 , therefore molecular formula = $C_{10}H_{12}O_2$ $\checkmark$	
(ii)	carbonyl identified from IR (at about 1700cm⁻¹) ✓	5
	ring identified from nmr: $C_6H_5$ at $\delta$ = 7.3 ppm or from peak <i>m</i> / <i>e</i> = 77 in mass spectrum $\checkmark$	
	methyl groups identified from nmr: 2 x CH <sub>3</sub> s at $\delta$ = 1.2 ppm $\checkmark$	
	Single H identified from nmr by peak at $\delta$ = 4.0 ppm/ doublet at $\delta$ = 1.2 ppm indicates next to a CH $\checkmark$	
	$ \begin{array}{c} & & \\ & & $	
	Total	7

# **2815/06 Transition Elements**

Question	Expected Answers	Marks
1 (a) (i)	+6 / 6+ / 6	1
(ii)	$H_2$ $H_2$ $H_1$ $H_1$ $H_1$ $H_2$ $H_2$ $H_1$ $H_2$ $H_2$ $H_1$ $H_2$ $H_2$ $H_1$ $H_2$ $H_2$ $H_2$ $H_1$ $H_2$ $H_2$ $H_2$ $H_1$ $H_2$ $H_2$ $H_2$ $H_1$ $H_2$	
	$H_2$ and $H^+$ (state symbols not required) Voltmeter and salt bridge [if no liquid levels lose this mark] Both Platinum electrodes labelled $Cr_2O_7^{2-} / Cr^{3+} (/ H^+ not needed for mark)$ Temp 298K or 25°C + Pressure 1 Atm / 100 kPa/ 101 KPa / 1 bar / 10 <sup>5</sup> Pa + concentration 1 mol dm <sup>-3</sup> (can take from diagram) or equimolar mixture of $Cr_2O_7^{2-} / Cr^{3+}$	1 1 1 1 1
(b)	$E^{\Theta}$ would be lower than +1.33 V Equilibrium would move from right to left / backwards	1 1
	Total	8
2 (a)	$[Co(H_2O)_6]^{3^+}$ $E^{\Theta}$ for $[Co(H_2O)_6]^{3^+} + e^- \Rightarrow [Co(H_2O)_6]^{2^+}$ is more positive This reaction is more likely to proceed from left to right / $[Co(H_2O)_6]^{3^+}$ is more likely to accept electrons / be reduced	1 1 1
(b)	Pink	1
(c)	$2[Co(NH_3)_6]^{3^+}$ + Fe $\rightarrow 2[Co(NH_3)_6]^{2^+}$ + Fe <sup>2+</sup> idea that $E^{\Theta}$ is positive / idea of cobalt complex equilibrium tending to go from left to right whilst Fe <sup>2+</sup> / Fe equilibrium goes from right to left / calculation of cell emf ((+)0.55V).	1
	Total	6

Question	Expected Answers	Marks
3 (a) (i)	1s <sup>2</sup> 2s <sup>2</sup> 2p <sup>6</sup> 3s <sup>2</sup> 3p <sup>6</sup> 3d <sup>10</sup>	1
(ii)	Colour requires electrons in d-orbitals and a space for promotion / colour results from transitions of electrons between d-orbitals / colour results	
	from a partially filled d-subshell Cu <sup>+</sup> has a full set of 3d orbitals / is 3d <sup>10</sup> / has a full d-subshell	1
(b)		-
	Energy levels of d-orbitals are split	1
	Different ligands interact differently with d-orbitals / $\Delta E$ changes for different ligands	1
	Size of $\Delta {\pmb E}$ determines frequency / wavelength / colour of visible light absorbed	1
(c)	Disproportionation or explained eg copper(I) goes to copper(II) and copper / copper(I) is oxidised and reduced.	1
	$2CuCl \rightarrow CuCl_2 + Cu / 2Cu^+ \rightarrow Cu^{2+} + Cu$	1
	$CuCl_2$ / $Cu^{2+}$ / $[Cu(H_2O)_6]^{2+}$ and $CuCl_4^{2-}$ is blue-green and Cu is reddishpink (both needed for 1 mark)	1
(d)	Moles $S_2 O_3^{2-} = 0.00198$ mol	1
	1 mole $S_2O_3^{2-}$ = 1 mole $Cu^{2+}$ / 25 cm <sup>3</sup> $Cu^{2+}$ contains 0.00198 moles	1
	250 cm <sup>3</sup> of Cu <sup>2+</sup> contains 0.0198 moles	1
	0.0198 moles Cu has a mass of 0.0198 x 63.5 g = 1.26 g (1.2573)	1
	% Cu = 1.26/1.65 x 100 = 76.2% (allow 76.0% - 76.4%)	1
	Allow ecf after each stage of the calculation	
	Total	14



Question	Expected Answers	Marks
5	Co-ordination number is the number of dative covalent / co-ordinate bond formed (with central transition metal / ion)	ls 1
	Hyn Pt a $\begin{bmatrix} a \\ c \\$	
	Suitable charge / brackets needed for these examples	
	Square planar tetrahedral octahedral	
	90 <sup>°</sup> 109.5 <sup>°</sup> 90 <sup>°</sup>	
	Two marks for each type with suitable example and correct name of shap and bond angle. Clear 3-D diagram with correct bond angle for a correct complex will receive 2 marks	
	$[Co(H_2O)_6]^{2+} + 4Cl^- \Rightarrow [CoCl_4]^{2-} + 6H_2O / other suitable correct equations$	6 1
	Shape changes from octahedral to tetrahedral	1
	Co-ordination number changes from 6 to 4 } } 1 mark for both Charge changes from +2 to -2 }	1
	(Mark for co-ordination number and charge can be taken from equation	
	Quality of Written Communication:	
	1 mark awarded for the correct use in context of at least 2 of the following terms;	
	Square planar, tetrahedral, octahedral, dative, covalent, co-ordinate,	1
	Tot	al 11

# 2816/01 Unifying Concepts in Chemistry/ Experimental Skills 2 Written Paper

Question	Expected Answers	Marks
1(a)	$\mathcal{K}_{c} = \frac{[H_{2}][I_{2}]}{[HI]^{2}} \checkmark$	1
1(b)(i)	HI: 0.28 $\checkmark$ H <sub>2</sub> : 0.11 $\checkmark$	2
1(b)(ii)	Use of $K_{c} = \frac{0.11 \times 0.11}{0.28^{2}}$ to generate a calculated value $\checkmark$	3
	= 0.15 $\checkmark$ (2 significant figures) (calc. value: 0.154336735) no units $\checkmark$	
	There must be some response here, not left blank. If [HI] = 0.39 mol $dm^{-3}$ (common mistake),	
	$K_c = 0.07955292571$ (calc value) = 0.080 to 2 sig figs Do <b>NOT</b> accept 0.08 mol dm <sup>-3</sup> (1 significant figure)	
1(c)	$K_c$ doesn't change $\checkmark$ Composition stays the same <b>OR</b> equilibrium does not move $\checkmark$	2
1(d)	$K_c$ increases $\checkmark$ (forward) reaction is endothermic <b>OR</b> reverse reaction is exothermic $\checkmark$	2
1(e)	I : CI = $\frac{78.15}{127}$ : $\frac{21.85}{35.5}$ OR 0.615 : 0.615 $\checkmark$ A: ICI OR any multiple, eg I <sub>2</sub> CI <sub>2</sub> , etc $\checkmark$	5
	ICI with no working scored 2 marks.	
	HI + $CI_2 \longrightarrow ICI + HCI \checkmark$ <b>ACCEPT</b> 2HI + $CI_2 \longrightarrow 2ICI + H_2$ Accept multiples from identification of <b>A</b> .	
	Accept equation based on an incorrect formula for A but ONLY if a compound of I and CI	
	<b>B</b> : $I_2CI_6 \checkmark$ $2HI + 4CI_2 \longrightarrow I_2CI_6 + 2HCI \checkmark$ <b>ACCEPT</b> $2HI + 3CI_2 \longrightarrow I_2CI_6 + H_2 \checkmark$ Accept equation based on an incorrect formula for B but ONLY if a compound of I and CI	
	Total:	15

Question	Expected Answers	Marks
2(a)	3 🗸	1
2(b)	$k = \frac{6.90 \times 10^{-7}}{(2.80 \times 10^{54})^2 \times 1.44 \times 10^{53}} \checkmark$ = 6.11 x 10 <sup>3</sup> $\checkmark$ (calculator 6.111819728 × 10 <sup>3</sup> ) units: dm <sup>6</sup> mol <sup>-2</sup> s <sup>-1</sup> $\checkmark$ ACCEPT 6.1 × 10 <sup>3</sup> up to calculator value If expression is upside down, calculated value = 1.636173913 1.6 up to calculator value would score 1 mark for the numerical value ECF units dm <sup>-6</sup> mol <sup>2</sup> s <sup>1</sup>	3
	If square is missed, calculated value = $1.711309524$ 1.7 up to calculator value would score 1 mark for the numerical value <b>ECF</b> units dm <sup>3</sup> mol <sup>-1</sup> s <sup>-1</sup>	
2(c)(i)	Curve downwards with slope gradually levelling off $\checkmark$	1
2(c)(ii)	Measure its gradient <b>OR</b> slope $\checkmark$ (Tangent) at $t = 0$ <b>OR</b> at start $\checkmark$ <i>Either mark could be from triangle shown on graph with y/x</i>	2
2(c)(iii)	Half-life is constant ✓	1
2(d)(i)	Curve upwards with slope gradually getting steeper $\checkmark$	1
2(d)(ii)	rate × 9 <b>OR</b> $3^2 \checkmark$ order = 2 (with respect to NO) $\checkmark$ Each marking point is independent	2
2(d)(iii)	rate × $2^2$ × 3 = × 12 ✓	1
	Total:	12

Question	Expected Answers	Marks
3(a)	pK <sub>a</sub> = 2.82 ✓	1
	calculated value = 2.823908741	
	ACCEPT 2.8 up to calculator value	
3(b)(i)	$K_{a} = \frac{[H^{+}][HSO_{3}^{\$}]}{[H_{2}SO_{3}]} \checkmark$	1
	$K_a = \frac{1}{[H_aSO_a]}$	
3(b)(ii)	, [H <sup>+</sup> ] <sup>2</sup>	3
	$1.50 \times 10^{33} \approx \frac{[\text{H}^+]^2}{0.0265} \checkmark (=' \text{ sign is acceptable})$	
	0.0200	
	$[H^+] = \sqrt{1.50 \times 10^{\frac{53}{3}} \times 0.0265} = 6.30 \times 10^{-3} \text{ mol dm}^{-3} \checkmark$	
	$pH = -log[H^+] = -log 6.30 \times 10^{-3} = 2.20 \checkmark$	
	(Stand alone mark; ie pH –log(0.0265) = 1.58 can be awarded 1 mark)	
	If all figures kept in calculator, value = 2.200331434	
	ACCEPT 2.2 up to calculator value	
	If no square root, pH =. 4.40	
3(b)(iii)	a small amount of second dissociation	1
-()	<b>OR</b> it is a diprotic acid $\checkmark$	-
	<b>ACCEPT</b> equilibrium concentration $H_2SO_3$ is less than the initial	
	concentration.	
3(c)(i)	ionic product (of water) ✓	1
3(c)(ii)	$K_{\rm w} = [{\rm H}^+] [{\rm OH}^-] \checkmark$	1
o( 1)	×	0
3(d)	$[H^+] = \frac{1.0 \times 10^{514}}{0.0265}$ <b>OR</b> 3.77 × 10 <sup>-13</sup> <b>OR</b> pOH –log(0.0265) = 1.58 \checkmark	2
	$pH = -log(3.77 \times 10^{-13})$ <b>OR</b> 14 - 1.58 = 12.42 $\checkmark$	
	calculated value = 12.42324587	
2(-)	ACCEPT 12.4 up to calculator value	4
3(e)	$\mathbf{C}$ : KHSO <sub>3</sub> $\checkmark$	4
	$KOH + H_2SO_3 \longrightarrow KHSO_3 + H_2O\checkmark$	
	<b>D</b> : $K_2SO_3 \checkmark$	
	$2KOH + H_2SO_3 \longrightarrow K_2SO_3 + 2H_2O /$	
	$KOH + KHSO_3 \longrightarrow K_2SO_3 + H_2O\checkmark$	
	If <b>C</b> and <b>D</b> are the wrong way around award 3 max by <b>ECF</b>	
	If H <sub>2</sub> SO <sub>4</sub> used throughout, award 3 max by ECF	4.4
	Total:	14

Question		Marks
4(a)	$C_2H_4O_3 \checkmark$	1
4(b)	Stage 1:CICH2COOH + 2NaOH $\longrightarrow$ HOCH2COONa + NaCI + H2Oscores two marks $\checkmark \checkmark$	3
	CICH <sub>2</sub> COOH + NaOH $\longrightarrow$ HOCH <sub>2</sub> COONa + HCI scores one mark $\checkmark$	
	CICH <sub>2</sub> COOH + NaOH $\longrightarrow$ CICH <sub>2</sub> COONa + H <sub>2</sub> O scores one mark $\checkmark$	
	CICH <sub>2</sub> COOH + NaOH $\longrightarrow$ HOCH <sub>2</sub> COOH + NaCl scores one mark $\checkmark$	
	Stage 2: HOCH <sub>2</sub> COONa + H <sup>+</sup> $\longrightarrow$ HOCH <sub>2</sub> COOH + Na <sup>+</sup> $\checkmark$ ACCEPT ECF from CICH <sub>2</sub> COONa forming CICH <sub>2</sub> COOH	
4(c)	buffer minimises <b>OR</b> resists pH changes ✓	
	$HOCH_2COOH \Rightarrow HOCH_2COO^- + H^+ \checkmark$	2
	For explanation below, accept HA and A <sup>-</sup> OR other weak acid added alkali reacts with H <sup>+</sup> / H <sup>+</sup> + OH <sup>-</sup> $\rightarrow$ H <sub>2</sub> O $\checkmark$ $\rightarrow$ HOCH <sub>2</sub> COO <sup>-</sup> / Equil $\rightarrow$ right (to counteract change) $\checkmark$	
	HOCH <sub>2</sub> COO <sup>−</sup> reacts with added acid or H <sup>+</sup> $\checkmark$ → HOCH <sub>2</sub> COOH / Equil → left (to counteract change) $\checkmark$	4
	$[H^+] = 10^{-pH} = 10^{-4.4} = 3.98 \times 10^{-5} \checkmark$	
	$\frac{[\text{HOCH}_2\text{COOH}]}{[\text{HOCH}_2\text{COO}^{\$}]} = \frac{[\text{H}^*]}{K_a}$	
	<b>OR</b> $\frac{[\text{HOCH}_2\text{COO}^{\$}]}{[\text{HOCH}_2\text{COOH}]} = \frac{K_a}{[\text{H}^+]} \checkmark$	
	$\frac{[\text{HOCH}_2\text{COOH}]}{[\text{HOCH}_2\text{COO}^{\$}]} = \frac{3.98 \times 10^{\$}}{1.48 \times 10^{\$}} \text{ OR } 0.27$ OR	
	$\frac{[\text{HOCH}_2\text{COO}^{\$}]}{[\text{HOCH}_2\text{COOH}]} = \frac{1.48 \times 10^{\$4}}{3.98 \times 10^{\$5}} \text{ OR } 3.7 \checkmark$	3
	<b>QWC:</b> Buffer explanation includes discussion of equilibrium shift $\checkmark$	1

Question	Expected Answers	Marks
4(d)	moles $CO_2 = \frac{5.119}{44} = 0.116 \checkmark$	
	moles H <sub>2</sub> O = $\frac{1.575}{18}$ = 0.0875 <b>OR</b> moles H = 2 x 0.0875 = 0.175 $\checkmark$	
	moles $A = \frac{4.362}{150} = 0.0291 \checkmark$	
	Molar ratio <b>A</b> : C : H = 1 : $\frac{0.116}{0.0291}$ : $\frac{0.175}{0.0291}$ = 1 : 4 : 6 $\checkmark$	
	$O = 150 - (4 \times 12 + 6 \times 1) = 96$	
	moles $O = \frac{96}{16} = 6$	
	molecular formula = $C_4 H_6 O_6 \checkmark$	
		5
	ACCEPT suitable alternatives methods	
	e.g.	
	moles C = $\frac{5.119}{44}$ = 0.116 $\checkmark$	
	moles H = $2 \times \frac{1.575}{18} = 0.175 \checkmark$	
	(mass C = 1.396 g; mass H = 0.175 g; mass O = 2.791 g)	
	moles O = $\frac{2.791}{16}$ = 0.174 $\checkmark$	
	empirical formula = C : H : O = $C_2H_3O_3 \checkmark$	
	molecular formula = $C_4H_6O_6$ (related to 150) $\checkmark$	
	Total:	19

# 2816/03 Unifying Concepts in Chemistry/ Experimental Skills 2 Practical Examination

Skill P: 16 marks (out of up to 19 available)

The "expected" solution to the problem involves the following two stage procedure.  $KMnO_4$  titration (for  $H_2C_2O_4$ ) followed by adding Mg and gas collection, or precipitation (for both acids)

### T Titration - 6 marks

- T1 Equation for the redox reaction [1]  $2MnO_4^- + 16H^+ + 5C_2O_4^{2-} \rightarrow 2Mn^{2+} + 10CO_2 + 8H_2O_2$ or  $2KMnO_4 + 3H_2SO_4 + 5H_2C_2O_4 \rightarrow 2MnSO_4 + 10CO_2 + 8H_2O + K_2SO_4$ Use KMnO<sub>4</sub> of concentration 0.0250 - 0.0500 mol dm<sup>-3</sup> in the burette T2 and a verbal statement that KMnO<sub>4</sub> is an oxidising agent (*owtte*) [1] If 10 cm<sup>3</sup> pipette is used for acid, then concentration is 0.0100 - 0.0200 mol dm<sup>-3</sup> Concentration does not need to be justified, but must give a titre 15 - 40 cm<sup>3</sup> T3 Pipette a known volume of acid mixture into a [conical] flask and acidify with extra sulphuric acid (or reference to sulphuric acid in the mixture) and heat the mixture. [1] Τ4 No indicator is required (implied) and end colour change in titration is from colourless/decolorised to [pale] pink [1] Allow pink or pale purple as the final colour T5 Titrate until two consistent/concordant accurate titres are obtained [1] Accept "titres within 0.1 cm<sup>3</sup>" (unit needed). T6 Calculate concentration of  $H_2C_2O_4$  in the solution from specimen titration data. [1] Gas collection – 9 marks G G1 Pipette known volume of acid mixture and add Mg [1] No mark for selection of any other metal or of a metal carbonate Thereafter mark ecf for other metals or a metal carbonate or NaHCO<sub>3</sub> G2 Precaution: Use Mg in excess and in the form of powder/turnings or sandpaper/remove oxide layer before use **Reason:** to ensure that both acids in the mixture react completely or speed reaction up or reference to MgO reacting to produce no gas [1] Two precautions and one reason are required G3 Equations for reaction of **both** acids [1] Mg + H<sub>2</sub>SO<sub>4</sub>  $\rightarrow$  MgSO<sub>4</sub> + H<sub>2</sub> and Mg + H<sub>2</sub>C<sub>2</sub>O<sub>4</sub>  $\rightarrow$  MgC<sub>2</sub>O<sub>4</sub> + H<sub>2</sub> No G3 if metal carbonate produces an insoluble sulphate (eq  $CaCO_3$ ) G4 Calculation of suitable maximum volume of acid mixture to use [1]
- G5 Calculation of suitable minimum mass of Mg (or other reagent) to use [1]

Calculation must relate explicitly to the capacity of the syringe/collecting vessel.

G6	Collect gas in a gas syringe/ inverted burette/ inverted measuring cylinder <b>and</b> use of inner tube (or similar arrangement) to keep reagents apart at start Marks can be scored from a neat labelled diagram, if drawn If a metal carbonate was used, collection must be in a gas syringe	[1]
G7	Record final volume of gas (when fizzing ceases/when syringe stops moving) Visual observation to indicate completion of reaction is required	[1]
G8	Repeat procedure and take mean of gas volumes/until volumes are consistent	[1]
G9	Specimen calculation of the concentration of $H_2SO_4$ in mixture Calculation must use volume of gas collected and the answer from strand T	[1]
<u>P</u>	Precipitation (P) – 7 marks (maximum)	
P1	Add <u>excess</u> aqueous barium nitrate/chloride to known volume (measured by pipette/ burette) of mixed acid.	[1]
P2	Equation for precipitation reaction(s), with state symbols <b>and</b> justification of volume ( <b>or</b> mass) of aq barium compound used	[1]
P3	Filter using reduced pressure filtration with pre-weighed filter paper	[1]
P4	<ul> <li>Two precautions from</li> <li>Wash all solid on to the filter funnel with distilled water</li> <li>Use of fine grade (or multiple sheets of) filter paper</li> <li>Wash solid on filter paper [to remove impurities]</li> </ul>	
	<ul> <li>Repeat whole experiment and take mean results</li> </ul>	[1]
P5	Dry in desiccator <b>or</b> in an oven until constant mass is achieved	[1]
P6	Specimen calculation of concentration of sulphuric acid $M_r$ of sulphate (for Ba =233; Ca = 136) must be shown in working	[1]
P7	A clear discussion of the solubility of barium ethanedioate or calculation of mass of barium ethanedioate precipitated	[1]
<u>s</u>	Safety, Sources and QWC - 4 marks	
S1	<b>Safety</b> : Sulphuric acid is irritant/corrosive ( <i>depending on context of use in Plan</i> ): use safety spectacles or lab coat <b>or</b> wash with plenty of water if spilt	[1]
S2	<ul> <li>Two sources quoted in the text or at end of Plan.</li> <li>Book references must have chapter or page numbers</li> <li>Internet reference must go beyond the first slash of web address</li> <li>Accept one reference to "Hazcards" without any qualification</li> </ul>	[1]
S3	<b>QWC</b> : text is legible and spelling, punctuation and grammar are accurate Allow no more than five different errors in legibility, spelling, punctuation or grammar.	[1]

# 2816/03

### S4 **QWC**: information is organised clearly and coherently

- Is a word count given and within the limits 450 1050 words?
- Are scientific language used and chemical formulae quoted correctly?

**Mark Scheme** 

Are both experiments described in a logical sequence?

### Practical Test (B)

### Page 3 : Part 1

### Presentation of titration data

- Table grid <u>drawn</u> (at least three lines) **and** all burette data shown in the table, including first/trial.
- <u>One</u> correctly labelled table three burette volumes clearly indicated and recorded *"Reversed" data (eg final = 0.00 and initial = 27.30) forfeits this bullet*
- Three (or more) sets of titration readings are shown
- All "accurate" burette data are quoted to two decimal places, ending in .00 or .05 The 2 d.p. rule does not apply to the trial (if labelled as trial) but it does apply to "0.00 cm<sup>3</sup>"
- No readings recorded above 50 cm<sup>3</sup>
- All subtractions are shown and are correct

#### Six bullets correct = 2 marks Five bullets correct = 1 mark

#### Self-consistency of titres

- The two selected titres are within 0.10 cm<sup>3</sup> at the **first** end-point
- The two selected **total** titres are within 0.10 cm<sup>3</sup> at the **second** end-point If 3 titres are used to calculate mean, assess on the spread of **all** three accurate titres. For incorrect subtraction, award these marks using the candidate's own values. The candidate may choose the trial titre if its value makes it appropriate to do so
- Units, cm<sup>3</sup> or ml, must be given somewhere (once in the table is sufficient).

#### Three bullets correct = 2 marks: two bullets correct = 1 mark

#### Mean titres (A, B and C) calculated

- A may be calculated using either the closest two titres or the two accurate titres. A candidate may use three titres if all are within 0.10 cm<sup>3</sup>.
   e.g. if titres are 26.00, 26.10, 26.00 - means of 26.00, 26.03, or 26.05 are all correct Give the candidate the benefit of any legitimate doubt.
- **B** may be calculated using either the closest two titres **or** the two accurate titres. A candidate may use different experiments to calculate **A** and **B**
- Correct subtraction to give C
- Units are shown at least once in this section
- The means must be quoted correctly to 1 or 2 d.p (or to 3 d.p. only for .025 or .075)

#### Five bullets correct = 2 marks: four bullets correct = 1 mark

#### Accuracy

Write down the supervisor's mean titres (A and B) next to the candidate's table. Check the candidate's mean titres (A and B).

- Use the candidate's own means to assess accuracy, where suitable.
- Round the candidate's mean to the nearest 0.05 cm<sup>3</sup>
- Compare the mean supervisor's titres with the candidate's titres.
- Use the conversion chart below to award the marks for accuracy.

#### [2]

[10]

[2]

# [17 marks]

[2]

The accuracy for each mean titre is marked out of 5 (A is phenolphthalein end-point and B is methyl orange end-point)

A/B is within 0.25 cm <sup>3</sup> of mean supervisor's value	[5]
A/B is within 0.40 cm <sup>3</sup> of mean supervisor's value	[4]
A/B is within 0.60 cm <sup>3</sup> of mean supervisor's value	[3]
A/B is within 0.90 cm <sup>3</sup> of mean supervisor's value	[2]
A/B is within 1.20 cm <sup>3</sup> of mean supervisor's value	[1]

# Spread penalty

("Sp If the	read" is defined by the titres used by the candidate to calculate the mean) e titres have a spread of 0.40 cm <sup>3</sup> or more, deduct 1 mark. e titres have a spread of 0.90 cm <sup>3</sup> or more, deduct 2 marks (max on each titration	n).
	ety nolphthalein contains ethanol since it has flammable hazard symbol henolphthalein, since methyl orange contains water which is not flammable.	[1]
Page	<u>es 4 - 6: Part 2</u>	[13 marks]
	k <b>ecf</b> wherever possible from one part of an answer to the next. wers should be quoted to <b>three</b> significant figures (if the answer is a marking poi	nt)
(a)	Concentration = $2.50 \times 10/250$ = 0.100 mol dm <sup>-3</sup>	[1]
(b)	Answer (a) is multiplied by <sup>difference between mean titres</sup> / <sub>1000</sub> This is a method mark but the correct mean <b>C</b> must be used	[1]
(c)	Answer quoted must be the same as that given in <b>(b)</b>	[1]
(d)	<b>1 mark</b> Second chemical equation shows that 1 $\underline{mol}$ Na <sub>2</sub> CO <sub>3</sub> gives 1 $\underline{mol}$ NaHCO <sub>3</sub>	[1]
(e)(i)	) Answer is 40 times that given in (c)	[1]
(e)(i	<b>i) 2 marks</b> <i>M</i> <sub>r</sub> of sodium carbonate = 106	[1]
	Correct answer obtained by candidate <i>No</i> ecf allowed	[1]
(f)(i)	Volume correctly calculated, $(A - C)$ , to one <b>or</b> two d.p.	[1]
(f)(ii	) No of moles of HCI used = <sup>answer (i) x answer (a)</sup> / <sub>1000</sub>	[1]
(f)(ii	i)Concentration of NaOH = no of moles of HCI x 40	[1]
	Concentration of NaOH used, correctly worked out	[1]
(f)(iv	<i>r)M</i> <sub>r</sub> of NaOH is 40	[1]
	Mass of NaOH (= 40 x iii), correctly calculated	[1]

		,
Page	<u>es 7 + 8: Part 3</u> (Skill E)	[14 marks]
(a)(i)	) 3 marks	
	Correct volumes used in both calculations – 10 $\text{cm}^3$ and 250 $\text{cm}^3$	[1]
	% error in pipette = 0.30 % ( <i>or</i> 99.70% accuracy)	[1]
	% error in vol flask = 0.080 % ( <i>or</i> 99.92% accuracy), so flask is more accura	ate [1]
(ii)	2 marks	
	Burette reads to 0.05/ 0.1 cm <sup>3</sup> so is less accurate than pipette [reads to 0.03	B] <b>[1]</b>
	Burette must be read twice, so error is increased/ accuracy reduced further	[1]
(b)	2 marks	
	Idea that reliable end points are consistent <i>(or vice versa)</i> This mark not awarded if accuracy ideas are discussed (apparatus or indica	<b>[1]</b> tor)
	Candidate uses his/her own titre values appropriately to decide which end-p reliable <b>or</b> less reliable ( <b>or</b> justifies that both were equally reliable)	oint was more [1]
(c)	<b>5 marks (max)</b> – but only 4 on question paper <i>Mark the best two strands</i>	
	• Methyl orange <b>or</b> phenolphthalein end-point is indistinct /difficult to jud	ge [1]
	Colour change at the end of each stage is gradual <i>Reference to specific end colour of chosen indicator is needed</i>	[1]
	Double indicator titration involves more burette readings	[1]
	There is cumulative error or increased % error in the titration	[1]
	Sensible attempt made to calculate a combined % error in expt	[1]
	• Volume C is small/ <u>only</u> c13 cm <sup>3</sup> [compared to a normal mean titre]	[1]
	The $\underline{\%}$ error in measuring this volume is higher than in a titre of about	25 cm <sup>3</sup> [1]
	Indicators interfere with each other	[1]
(d)	3 marks (max) Credit any three points from the following	
	Sodium carbonate is a base/ alkali <b>or</b> both chemicals contain OH <sup>-</sup> ions	[1]
	The pH of the mixed solution is higher than that of NaOH alone	[1]
	However, the effect on pH is small since $Na_2CO_3$ is a weak alkali or $CO_3^{2-} + H_2O \Leftrightarrow HCO_3^{-} + OH^{-}$	[1]
	Correct concentration of NaOH corresponding to any stated pH value <i>or</i> equation(s) linking the pH to the hydroxide ion concentration <i>or</i> logarithmic scale increases % error in measurement of concentration	[1]

# **Grade Thresholds**

# Advanced GCE Chemistry (3882/7882) January 2009 Examination Series

### Unit Threshold Marks

Unit		Maximum Mark		b	С	d	е	u
2811	Raw	60	51	45	40	35	30	0
	UMS	90	72	63	54	45	36	0
2812	Raw	60	49	44	39	34	29	0
	UMS	90	72	63	54	45	36	0
2813A	Raw	120	98	88	78	69	60	0
	UMS	120	96	84	72	60	48	0
2813B	Raw	120	98	88	78	69	60	0
	UMS	120	96	84	72	60	48	0
2813C	Raw	120	94	84	74	64	55	0
	UMS	120	96	84	72	60	48	0
2814	Raw	90	73	65	57	49	42	0
	UMS	90	72	63	54	45	36	0
2815A	Raw	90	74	68	62	56	50	0
	UMS	90	72	63	54	45	36	0
2815C	Raw	90	75	68	61	54	47	0
	UMS	90	72	63	54	45	36	0
2815E	Raw	90	77	70	63	56	49	0
	UMS	90	72	63	54	45	36	0
2816A	Raw	120	100	90	80	70	61	0
	UMS	120	96	84	72	60	48	0
2816B	Raw	120	100	90	80	70	61	0
	UMS	120	96	84	72	60	48	0
2816C	Raw	120	92	82	72	62	53	0
	UMS	120	96	84	72	60	48	0

#### **Specification Aggregation Results**

Overall threshold marks in UMS (ie after conversion of raw marks to uniform marks)

	Maximum Mark	Α	В	C	D	E	U
3882	300	240	210	180	150	120	0
7882	600	480	420	360	300	240	0

The cumulative percentage of candidates awarded each grade was as follows:

	A	В	С	D	E	U	Total Number of Candidates
3882	10.0	35.9	60.7	84.1	98.4	100	1206
7882	14.9	49.3	78.0	92.9	98.1	100	288

#### 1694 candidates aggregated this series

For a description of how UMS marks are calculated see: <u>http://www.ocr.org.uk/learners/ums\_results.html</u>

Statistics are correct at the time of publication.

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