



JABchem



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Past Papers Advanced Higher Chemistry

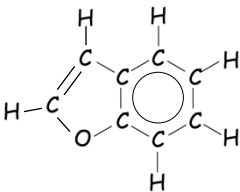
2018 Marking Scheme

Grade Awarded	Mark Required		% candidates achieving grade
	(/130)	%	
A	88+	67.7%	31.5%
B	73+	56.2%	27.6%
C	59+	45.4%	23.2%
D	52+	40%	8.1%
No award	<52	<40%	9.6%

Section:	Multiple Choice	Extended Answer	Project
Average Mark:	18.8 /30	39.3 /70	19.1 /30

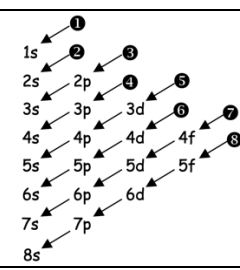
2018 Adv Higher Chemistry Marking Scheme


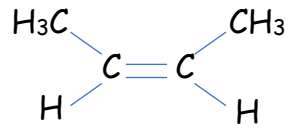

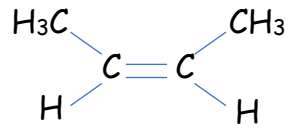

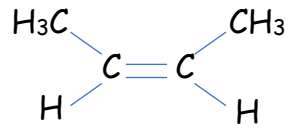

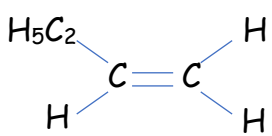
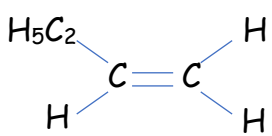
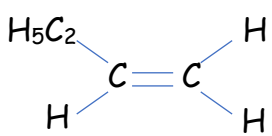
MC Qu	Answer	% Pupils Correct	Reasoning
1	A		<input checked="" type="checkbox"/> A Beta particles which are electrons emitted from the nucleus <input checked="" type="checkbox"/> B Gamma rays are a form of electromagnetic radiation with wavelength & frequency <input checked="" type="checkbox"/> C Infrared is a form of electromagnetic radiation with wavelength & frequency <input checked="" type="checkbox"/> D ultraviolet is a form of electromagnetic radiation with wavelength & frequency
2	D		<input checked="" type="checkbox"/> A s-block is found in groups 1+2 of the periodic table <input checked="" type="checkbox"/> B p-block is found in groups 2 through to group 0 of the periodic table <input checked="" type="checkbox"/> C d-block is found between groups 2+3 of the periodic table (transition metals) <input checked="" type="checkbox"/> D f-block is the two groups at the bottom of periodic table (Actinides and Lanthanides)
3	A		The d-orbital shown (d_{xy}) in the question will hold a maximum of two electrons.
4	D		<p style="text-align: center;">trigonal planar tetrahedral</p>
5	B		Ni atom: $1s^2 2s^2 2p^6 3s^2 3p^6 3d^8 4s^2$ ∴ Ni^{2+} ion: $1s^2 2s^2 2p^6 3s^2 3p^6 3d^8$ <input checked="" type="checkbox"/> A Ni^{2+} ions have 8 electrons in the 3d shell as 4s electrons are removed before 3d <input checked="" type="checkbox"/> B The six electrons occupy the three lower 3d orbitals after 3d splits into 2 levels <input checked="" type="checkbox"/> C The ligands in the complex split the 3d orbitals into two levels. <input checked="" type="checkbox"/> D Ni^{2+} ions have 8 electrons in the 3d shell as 4s electrons are removed before 3d
6	C		<input checked="" type="checkbox"/> A Oxidation number of Mn in MnO_4^- = +7 <input checked="" type="checkbox"/> B Oxidation number of Mn in MnO_4^{2-} = +6 <input checked="" type="checkbox"/> C Oxidation number of Mn in MnO_4^{3-} = +5 <input checked="" type="checkbox"/> D Oxidation number of Mn in MnO_2 = +4
7	A		<input checked="" type="checkbox"/> A increase in temperature decreases K and decreases concentration of SO_2 <input checked="" type="checkbox"/> B increase in temperature favours reverse endothermic reaction ∴ less products <input checked="" type="checkbox"/> C increase in temperature favours reverse endothermic reaction ∴ K decreases <input checked="" type="checkbox"/> D increase in temperature favours reverse endothermic reaction ∴ K decreases
8	D		$ \begin{array}{ccccccc} H_2CO_3 & + & CN^- & \rightleftharpoons & HCN & + & HCO_3^- \\ \text{Acid} & & \text{Base} & & \text{Conjugate Acid} & & \text{Conjugate Base} \\ \text{Donates } H^+ & & \text{Accepts } H^+ & & \text{Formed when Base accepts } H^+ & & \text{Formed when Acid loses } H^+ \end{array} $
9	B		$pH = 8.5 \therefore -\log_{10}[H^+] = 8.5 \therefore \log_{10}[H^+] = -8.5 \therefore [H^+] = 10^{-8.5} = 3.16 \times 10^{-9} \text{ mol l}^{-1}$ $[H^+][OH^-] = 10^{-14} \therefore [OH^-] = \frac{10^{-14}}{[H^+]} = \frac{1 \times 10^{-14}}{3.16 \times 10^{-9}} = 3.16 \times 10^{-6} \text{ mol l}^{-1}$
10	C		<input checked="" type="checkbox"/> A Catalysts do not change the position of equilibrium <input checked="" type="checkbox"/> B H_3O^+ ions are a product so adding sulphuric acid will shift equilibrium to left <input checked="" type="checkbox"/> C NaOH neutralises H_3O^+ ions so removing a product and shifts equilibrium to right <input checked="" type="checkbox"/> D $C_3H_7COO^-$ ions are a product so adding sulphuric acid will shift equilibrium to left
11	D		<input checked="" type="checkbox"/> A sodium sulphate solution is neutral pH=7 (strong acid v strong alkali) <input checked="" type="checkbox"/> B lithium chloride solution is neutral pH=7 (strong acid v strong alkali) <input checked="" type="checkbox"/> C ammonium nitrate solution is acidic pH<7 (strong acid v weak alkali) <input checked="" type="checkbox"/> D potassium propanoate solution is alkaline pH>7 (weak acid v strong alkali) H^+ ions join up with propanoate ions to form molecules of propanoic acid. Water molecules then split into ions to replace H^+ ions but concentration of OH^- builds up as H^+ is removed.

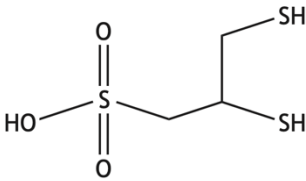
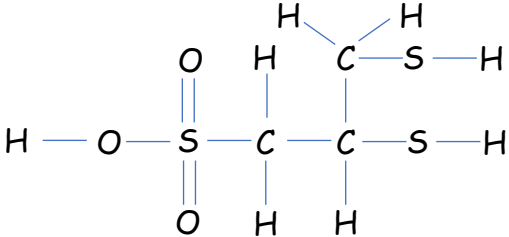
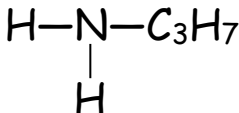
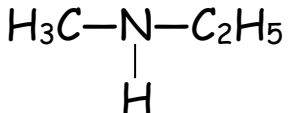
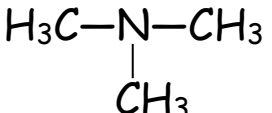
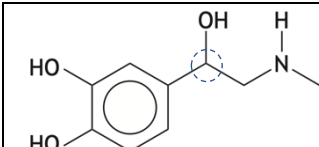
12	B	<input checked="" type="checkbox"/> A sodium chloride is made from a strong alkali so no buffer forms <input checked="" type="checkbox"/> B buffers form when salt of a weak alkali (ammonium chloride) dissolves in a weak acid (ammonia) <input checked="" type="checkbox"/> C sodium hydroxide is a strong acid and no buffer forms <input checked="" type="checkbox"/> D sodium hydroxide is a strong acid and no buffer forms															
13	D	$\Delta G = \Delta H - T\Delta S \therefore \Delta G - \Delta H = -T\Delta S$ If $\Delta G - \Delta H$ is approximately zero then $-T\Delta S$ must also be approximately zero. This means ΔS must be approximately zero then there must be little change to disorder during the reaction. <input checked="" type="checkbox"/> A CO_2 gas released which increases disorder and increases ΔS <input checked="" type="checkbox"/> B Two gases formed from a solid which increases disorder and increases ΔS <input checked="" type="checkbox"/> C H_2 gas released which increases disorder and increases ΔS <input checked="" type="checkbox"/> D Solid and an ion turning into a solid and an ion keeps disorder level (ΔS) similar															
14	B	<input checked="" type="checkbox"/> A The overall order is the sum of the individual orders \therefore overall order = $1+2 = 3$ <input checked="" type="checkbox"/> B This reaction must have second step as equal no. of moles of P + Q are used up in the equation but the rate determining step has two particles of Q reacting with one particle of P. A second particle of P must react in the other step. <input checked="" type="checkbox"/> C Rates of reaction always decrease as concentration of reactants decreases <input checked="" type="checkbox"/> D As P is first order then doubling [P] will double the rate of reaction															
15	B	<table border="1"> <thead> <tr> <th>Bond</th> <th>C-H</th> <th>C-C</th> <th>C=C</th> </tr> </thead> <tbody> <tr> <td>Type of hybridisation</td> <td>sp^3 hybridisation</td> <td>sp^3 hybridisation</td> <td>sp hybridisation</td> </tr> </tbody> </table>	Bond	C-H	C-C	C=C	Type of hybridisation	sp^3 hybridisation	sp^3 hybridisation	sp hybridisation							
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16	C	 <p style="text-align: right;">formula $\text{C}_8\text{H}_6\text{O}$ $\text{gfm} = (8 \times 12) + (6 \times 1) + (1 \times 16) = 96 + 6 + 16 = 118\text{g}$</p>															
17	A	<input checked="" type="checkbox"/> A W and X in same position in both diagrams but Z and Y in opposite positions <input checked="" type="checkbox"/> B XYZ on bottom are in same anti-clockwise as comparison diagram. <input checked="" type="checkbox"/> C XYZ on bottom are in same anti-clockwise as comparison diagram. <input checked="" type="checkbox"/> D XYZ on bottom are in same anti-clockwise as comparison diagram.															
18	A	<input checked="" type="checkbox"/> A sodium + butan-1-ol react to form sodium butoxide $\text{CH}_3\text{CH}_2\text{CH}_2\text{CHO}^- \text{Na}^+ + \text{H}_2$ <input checked="" type="checkbox"/> B sodium + butanoic acid would form sodium butanoate + hydrogen <input checked="" type="checkbox"/> C sodium hydroxide does not react with alcohols <input checked="" type="checkbox"/> D sodium hydroxide + butanoic acid would form sodium butanoate + hydrogen															
19	C	<input checked="" type="checkbox"/> A Hydration: Adding water across a double or triple bond <input checked="" type="checkbox"/> B Oxidation: Increasing the oxygen : hydrogen ratio in a compound <input checked="" type="checkbox"/> C Hydrolysis: spitting into two molecules with water added at break <input checked="" type="checkbox"/> D Hydrogenation: Adding hydrogen across a double bond or triple bond															
20	B	<table border="1"> <thead> <tr> <th>Element</th> <th>Cu</th> <th>O</th> </tr> </thead> <tbody> <tr> <td>Mass</td> <td>16g</td> <td>2g</td> </tr> <tr> <td>No. of moles (divide % by gfm)</td> <td>$\frac{16}{63.5}$ = 0.252</td> <td>$\frac{2}{16}$ = 0.125</td> </tr> <tr> <td>Mole ratio (divide through by smallest value)</td> <td>$\frac{0.252}{0.125}$ = 2.02</td> <td>$\frac{0.125}{0.125}$ = 1.00</td> </tr> <tr> <td>Round to Whole Number</td> <td>2</td> <td>1</td> </tr> </tbody> </table>	Element	Cu	O	Mass	16g	2g	No. of moles (divide % by gfm)	$\frac{16}{63.5}$ = 0.252	$\frac{2}{16}$ = 0.125	Mole ratio (divide through by smallest value)	$\frac{0.252}{0.125}$ = 2.02	$\frac{0.125}{0.125}$ = 1.00	Round to Whole Number	2	1
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21	C	<input checked="" type="checkbox"/> A gfm $\text{CH}_3\text{OCOCH}_3 = (3 \times 12) + (6 \times 1) + (2 \times 16) = 36 + 6 + 32 = 74\text{g}$ <input checked="" type="checkbox"/> B gfm $\text{CH}_3\text{CH}_2\text{COOH} = (3 \times 12) + (6 \times 1) + (2 \times 16) = 36 + 6 + 32 = 74\text{g}$ <input checked="" type="checkbox"/> C gfm $\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{NH}_2 = (4 \times 12) + (11 \times 1) + (1 \times 14) = 48 + 11 + 14 = 73\text{g}$ <input checked="" type="checkbox"/> D gfm $\text{CH}_3\text{CH}(\text{OH})\text{CH}_2\text{CH}_3 = (4 \times 12) + (10 \times 1) + (1 \times 16) = 48 + 10 + 16 = 74\text{g}$															
22	C	<input checked="" type="checkbox"/> A Absorption peak at $3100 - 3000\text{cm}^{-1}$ due to C-H stretch in benzene ring <input checked="" type="checkbox"/> B Absorption peak at $2962 - 2853\text{cm}^{-1}$ due to C-H stretch in an alkane <input checked="" type="checkbox"/> C No absorption peak at $1730 - 1717\text{cm}^{-1}$ as there is no C=O aromatic ester in eugenol <input checked="" type="checkbox"/> D Absorption peak at $1150 - 1070\text{cm}^{-1}$ due to C-O stretch in alkyl ether															

23	A	<input checked="" type="checkbox"/> A Agonist: Binds to the receptor and causes an internal response in the cell <input checked="" type="checkbox"/> B Antagonist: Bind to the receptor but does not cause internal response in the cell <input checked="" type="checkbox"/> C Inhibitor: Block the substrate from entering the receptor <input checked="" type="checkbox"/> D Receptor: Protein in membrane of cells that allows molecules to bind with it.
24	B	no. of mol = volume x concentration = 0.05litres x 2 mol l ⁻¹ = 0.1mol concentration = $\frac{\text{no. of mol}}{\text{volume}} = \frac{0.1 \text{ mol}}{0.25\text{litres}} = 0.4 \text{ mol l}^{-1}$
25	B	<input checked="" type="checkbox"/> A CH ₃ CH ₂ CH ₂ CH ₂ CHO is an aldehyde and is miscible with water <input checked="" type="checkbox"/> B CH ₃ CH ₂ OCH ₂ CH ₂ CH ₃ is an ether and is immiscible with water <input checked="" type="checkbox"/> C CH ₃ CH ₂ CH ₂ CH ₂ COOH is a carboxylic acid and is miscible with water <input checked="" type="checkbox"/> D CH ₃ CH(OH)CH ₂ CH ₂ CH ₃ is an alcohol and is miscible with water
26	D	<input checked="" type="checkbox"/> A This is a step in a recrystallisation technique <input checked="" type="checkbox"/> B This is a step in a recrystallisation technique <input checked="" type="checkbox"/> C This is a step in a recrystallisation technique <input checked="" type="checkbox"/> D This is a step in gravimetric analysis to show all dissolved ion has precipitated.
27	B	<input checked="" type="checkbox"/> A Purification should give a melting point over a narrower range <input checked="" type="checkbox"/> B Recrystallisation raises the melting point and narrows the temperature range <input checked="" type="checkbox"/> C Impurities lower the melting point so purification raises the melting point <input checked="" type="checkbox"/> D Impurities lower the melting point so purification raises the melting point
28	C	<input checked="" type="checkbox"/> A Desiccators remove moisture from the atmosphere not oxygen <input checked="" type="checkbox"/> B Heating the sample removes water from the sample, not the desiccator. <input checked="" type="checkbox"/> C While sample is cooling in desiccator, moisture cannot be reabsorbed by the sample <input checked="" type="checkbox"/> D Desiccators do not prevent decomposition of sample, probably in the heating stage
29	D	<input checked="" type="checkbox"/> A The distance moved by the solvent does not alter the value of the R _f value <input checked="" type="checkbox"/> B The sample will move the same distance regardless of concentration <input checked="" type="checkbox"/> C The length of TLC plate is not a factor in the R _f value <input checked="" type="checkbox"/> D The solvent used decides how far the sample moves and the R _f value.
30	C	<input checked="" type="checkbox"/> A Distillation separates the chemicals while refluxing returns the chemical to flask <input checked="" type="checkbox"/> B Distillation separates the chemicals while refluxing returns the chemical to flask <input checked="" type="checkbox"/> C Coldest water (from tap) should be nearest hot vapours entering condenser <input checked="" type="checkbox"/> D If the water travels against the flow in condenser arm then the hottest water is nearest the flask and the vapour will travel further up the condenser

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Long Qu	Answer	Reasoning																																																																																																						
1a(i)	sodium	From page 15 of data booklet: FLAME COLOURS <table border="1"> <tr> <td>Element</td> <td>Barium</td> <td>Calcium</td> <td>Copper</td> <td>Lithium</td> <td>Potassium</td> <td>Sodium</td> <td>Strontium</td> </tr> <tr> <td>Wavelength /nm</td> <td>554</td> <td>620</td> <td>522</td> <td>671</td> <td>405</td> <td>589</td> <td>650</td> </tr> <tr> <td>Colour</td> <td>green</td> <td>orange-red</td> <td>blue-green</td> <td>crimson</td> <td>lilac</td> <td>orange-yellow</td> <td>red</td> </tr> </table>	Element	Barium	Calcium	Copper	Lithium	Potassium	Sodium	Strontium	Wavelength /nm	554	620	522	671	405	589	650	Colour	green	orange-red	blue-green	crimson	lilac	orange-yellow	red																																																																														
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1a(ii)	425nm	$E = \frac{L \times h \times c}{\lambda} \therefore \lambda = \frac{L \times h \times c}{E} = \frac{6.02 \times 10^{23} \text{ mol}^{-1} \times 6.63 \times 10^{-34} \text{ J s} \times 3 \times 10^8 \text{ m s}^{-1}}{282 \times 1000 \text{ J mol}^{-1}}$ $= 4.25 \times 10^{-7} \text{ m}$ $= 425 \text{ nm}$																																																																																																						
1a(iii)	87	Relative Intensity 375 \longleftrightarrow 435mg kg ⁻¹ Relative Intensity 75 \longleftrightarrow 435mg kg ⁻¹ $\times \frac{75}{375}$ $= 87 \text{ mg kg}^{-1}$																																																																																																						
1b(i)	Orbitals fill up in order of increasing energy	<u>Aufbau Principle:</u> Electrons fill up in order of increasing energy: 1s 2s 2p 3s 3p 4s 3d 4p 5s 4d 5p 6s 4f 5d 6p 7s 5f 6d 7p 8s 																																																																																																						
1b(ii)	<table border="1"> <tr> <th>n</th> <th>l</th> <th>m_l</th> <th>m_s</th> </tr> <tr> <td>3</td> <td>1</td> <td>-1 or 0 or +1</td> <td>+$\frac{1}{2}$ or -$\frac{1}{2}$</td> </tr> </table>	n	l	m _l	m _s	3	1	-1 or 0 or +1	+ $\frac{1}{2}$ or - $\frac{1}{2}$	<table border="1"> <tr> <th>n</th> <th>Principal Quantum Number</th> <th colspan="4">n = Electron Shell Number</th> </tr> <tr> <th>l</th> <th>Angular Momentum Quantum Number</th> <th colspan="4">l = Type/Shape of subshell where l goes from 0 up to n-1</th> </tr> <tr> <td></td> <td></td> <td>l = 0</td> <td>l = 1</td> <td>l = 2</td> <td>l = 3</td> </tr> <tr> <td></td> <td></td> <td>s subshell</td> <td>p subshell</td> <td>d subshell</td> <td>f subshell</td> </tr> <tr> <th>m_l</th> <th>Magnetic Quantum Number</th> <th colspan="8">m_l is the orientation of the subshell where m_l goes from -l to +l</th> </tr> <tr> <td></td> <td></td> <th>Value of l</th> <th>Subshell type</th> <th colspan="5">Value of m_l</th> </tr> <tr> <td></td> <td></td> <td>0</td> <td>s</td> <td></td> <td></td> <td>0</td> <td></td> <td></td> <td></td> </tr> <tr> <td></td> <td></td> <td>1</td> <td>p</td> <td></td> <td>-1</td> <td>0</td> <td>1</td> <td></td> <td></td> </tr> <tr> <td></td> <td></td> <td>2</td> <td>d</td> <td></td> <td>-2</td> <td>-1</td> <td>0</td> <td>1</td> <td>2</td> </tr> <tr> <td></td> <td></td> <td>3</td> <td>f</td> <td>-3</td> <td>-2</td> <td>-1</td> <td>0</td> <td>1</td> <td>2</td> <td>3</td> </tr> <tr> <th>m_s</th> <th>Spin Quantum Number</th> <td colspan="8">m_s is the spin direction of an electron and has either value of +$\frac{1}{2}$ or -$\frac{1}{2}$</td> </tr> </table>	n	Principal Quantum Number	n = Electron Shell Number				l	Angular Momentum Quantum Number	l = Type/Shape of subshell where l goes from 0 up to n-1						l = 0	l = 1	l = 2	l = 3			s subshell	p subshell	d subshell	f subshell	m _l	Magnetic Quantum Number	m _l is the orientation of the subshell where m _l goes from -l to +l										Value of l	Subshell type	Value of m _l							0	s			0						1	p		-1	0	1					2	d		-2	-1	0	1	2			3	f	-3	-2	-1	0	1	2	3	m _s	Spin Quantum Number	m _s is the spin direction of an electron and has either value of + $\frac{1}{2}$ or - $\frac{1}{2}$							
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2a(i)	$K_a = \frac{[\text{HOOCCH}_2\text{CH}(\text{OH})\text{COO}^-][\text{H}_3\text{O}^+]}{[\text{HOOCCH}_2\text{CH}(\text{OH})\text{COOH}]}$	H ₂ O is both a reactant and the solvent $\therefore [\text{H}_2\text{O}] = 1$ $K_a = \frac{[\text{HOOCCH}_2\text{CH}(\text{OH})\text{COO}^-]^1 \times [\text{H}_3\text{O}^+]^1}{[\text{HOOCCH}_2\text{CH}(\text{OH})\text{COOH}]^1 \times [\text{H}_2\text{O}]^1} = \frac{[\text{HOOCCH}_2\text{CH}(\text{OH})\text{COO}^-][\text{H}_3\text{O}^+]}{[\text{HOOCCH}_2\text{CH}(\text{OH})\text{COOH}]}$																																																																																																						
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2b(i)	2.90	$\begin{aligned} \text{pH} &= \frac{1}{2}\text{p}K_a - \frac{1}{2}\log_{10} c \\ &= -\frac{1}{2}\log_{10} K_a - \frac{1}{2}\log_{10} c \\ &= -\frac{1}{2}\log_{10}(3.2 \times 10^{-4}) - \frac{1}{2} \times \log_{10}(0.0051) \\ &= (-\frac{1}{2} \times -3.49) - (\frac{1}{2} \times -2.29) \\ &= 1.75 - (-1.15) \\ &= 2.90 \end{aligned}$																																																																																																						
2b(ii)	Hydrogen bonding between chains	Each unit within a pectin chain has two hydroxyl -OH groups which would allow hydrogen bonding between the pectin chains and thicken the jam.																																																																																																						
3a	All four of the following required:	<table border="1"> <tr> <td>Dissolve sodium carbonate in deionised water</td> <td>Transfer the solution and the rinsings</td> <td>Use of a 250cm³ standard/volumetric flask</td> <td>Make up to the line flask with deionised water</td> </tr> </table>	Dissolve sodium carbonate in deionised water	Transfer the solution and the rinsings	Use of a 250cm ³ standard/volumetric flask	Make up to the line flask with deionised water																																																																																																		
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3b(i)	0.0348mol	$\text{Average titre} = \frac{19.5 + 19.4}{2} = 19.45 \text{ cm}^3$ $\text{no. of mol} = \text{volume} \times \text{concentration} = 0.01945 \text{ litres} \times 0.358 \text{ mol l}^{-1} = 6.97 \times 10^{-3} \text{ mol}$ $\text{Na}_2\text{CO}_3 + 2\text{HCl} \longrightarrow 2\text{NaCl} + \text{H}_2\text{O} + \text{CO}_2$ $\begin{array}{ccc} 1\text{mol} & & 2\text{mol} \\ 3.48 \times 10^{-3} \text{ mol} & & 6.97 \times 10^{-3} \text{ mol} \end{array}$ $25\text{cm}^3 \text{ sodium carbonate solution} \longleftrightarrow 3.48 \times 10^{-3} \text{ mol}$ $250\text{cm}^3 \text{ sodium carbonate solution} \longleftrightarrow 3.48 \times 10^{-2} \text{ mol}$																					
3b(ii)	7	$\text{gfm Na}_2\text{CO}_3 = (2 \times 23) + (1 \times 12) + (3 \times 16) = 46 + 12 + 48 = 106 \text{ g}$ $\text{mass} = \text{no. of mol} \times \text{gfm} = 0.0348 \times 106 = 3.69 \text{ g}$ <table border="0" style="width: 100%;"> <tr> <td style="width: 33%;">mass of water =</td> <td style="width: 33%;">mass of hydrated sodium carbonate =</td> <td style="width: 33%;">mass of unhydrated sodium carbonate =</td> </tr> <tr> <td></td> <td>8.10g</td> <td>3.69g</td> </tr> <tr> <td>mass of water =</td> <td>4.41g</td> <td></td> </tr> </table> <table border="1" style="width: 100%; text-align: center;"> <thead> <tr> <th>Chemical</th> <th>Na₂CO₃</th> <th>H₂O</th> </tr> </thead> <tbody> <tr> <td>Mass</td> <td>3.69g</td> <td>4.41g</td> </tr> <tr> <td>no. of mol (divide by gfm)</td> <td>$\frac{3.69}{106} = 0.0348$</td> <td>$\frac{4.41}{18} = 0.245$</td> </tr> <tr> <td>Divide by smallest value</td> <td>$\frac{0.0348}{0.0348} = 1$</td> <td>$\frac{0.235}{0.0348} = 7.04$</td> </tr> </tbody> </table> <p style="text-align: center;">Formula of hydrated sodium carbonate: Na₂CO₃·7H₂O</p>	mass of water =	mass of hydrated sodium carbonate =	mass of unhydrated sodium carbonate =		8.10g	3.69g	mass of water =	4.41g		Chemical	Na ₂ CO ₃	H ₂ O	Mass	3.69g	4.41g	no. of mol (divide by gfm)	$\frac{3.69}{106} = 0.0348$	$\frac{4.41}{18} = 0.245$	Divide by smallest value	$\frac{0.0348}{0.0348} = 1$	$\frac{0.235}{0.0348} = 7.04$
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4d(i) PART A	-126 kJ mol ⁻¹	$\begin{aligned} \Delta H^\circ &= \Sigma \Delta H_f^\circ(\text{products}) - \Sigma \Delta H_f^\circ(\text{reactants}) \\ &= (1 \times -6.99) - ((1 \times 119) + (1 \times 0)) \\ &= -6.99 - (119 + 0) \\ &= -6.99 - 119 \\ &= -125.99 \text{ kJ mol}^{-1} \end{aligned}$																					
4d(i) PART B	-23.2 J K ⁻¹ mol ⁻¹	$\begin{aligned} \Delta G^\circ &= \Sigma \Delta G^\circ(\text{products}) - \Sigma \Delta G^\circ(\text{reactants}) \\ &= (1 \times -65.9) - ((1 \times 185) + (1 \times 0)) \\ &= -65.9 - (185 + 0) \\ &= -65.9 - 185 \\ &= -119.1 \text{ kJ mol}^{-1} \end{aligned}$ $\Delta G^\circ = \Delta H^\circ - T \Delta S^\circ \quad \therefore \Delta S^\circ = \frac{\Delta G^\circ - \Delta H^\circ}{-T} = \frac{-119.1 - 126}{-298} = -0.0232 \text{ kJ K}^{-1} \text{ mol}^{-1}$ $= -23.2 \text{ J K}^{-1} \text{ mol}^{-1}$																					

4d(ii)	5431	The reaction becomes thermodynamically feasible when $\Delta G^\circ = 0$ $\Delta G^\circ = \Delta H^\circ - T\Delta S^\circ = 0 \quad \therefore T\Delta S^\circ = \Delta H^\circ \quad \therefore T = \frac{\Delta H^\circ}{\Delta S^\circ} = \frac{-126 \times 1000 \text{ J mol}^{-1}}{-23.2 \text{ J K}^{-1} \text{ mol}^{-1}} = 5431\text{K}$				
5a	Conjugated system	A conjugated system is a section of a compound with alternating C=C double bonds and C-C single bonds.				
5b	Answer to include:	1 st Mark	Electrons move from HOMO to LUMO (Highest Occupied Molecular Orbital) (Lowest Unoccupied Molecular Orbital)			
		2 nd Mark	Absorption of light (from the visible part of the spectrum) means that light of the complementary colour is seen			
5c	Answer to include:	1 st Mark	One from:	less conjugation	a shorter sequence of alternating double and single bonds	a smaller chromophore
		2 nd Mark	Larger gap between HOMO and LUMO and greater energy (absorbed) as shorter wavelength has greater energy			
6a	One answer from:	Ligands donate pairs of electrons to metal atom (lone pairs or non-bonding pair)			Ligands form dative covalent bonds with metal ion	
6b(i)	$C_3H_8O_3S_3$	 Skeletal Formula		 Full Structural Formula		
6b(ii) PART A	Bidentate	An electron pair on each Sulphur in the thiol -SH groups donate one of their lone pairs to the Mercury ion and forms a dative covalent bond.				
6b(ii) PART B	4	Each DMPMS ligand molecule donates lone pairs of from two of its sulphur atoms to give four dative covalent bonds on the Mercury ion in the centre.				
6c(i)	Gravimetric	Gravimetric analysis involves measuring the mass accurately to calculate the number of moles of substances.				
6c(ii)	96.1%	$\text{no. of mol} = \frac{\text{mass}}{\text{gfm}} = \frac{4.82\text{g}}{388.7 \text{ g mol}^{-1}} = 0.0167\text{mol}$ 0.0167mol complex contains 0.0167mol Ni ions $\text{mass of Ni} = \text{no. of mol} \times \text{gfm} = 0.0167\text{mol} \times 58.7\text{g mol}^{-1} = 0.98\text{g}$ $\% \text{ mass} = \frac{\text{mass of Ni}}{\text{mass of Alloy}} \times 100 = \frac{0.98}{1.02} \times 100 = 96.1\%$				
7a(i)	Electrophilic Substitution	Adding onto a benzene ring is electrophilic substitution. The H on the benzene ring is joins to the Cl and the remainder of that molecule joins onto the benzene ring.				
7a(ii)	Secondary	Primary Amine	Secondary Amine	Tertiary Amine		
		 1 Carbon attached to the Nitrogen	 2 Carbons attached to the Nitrogen	 3 Carbons attached to the Nitrogen		
7a(iii)	lithium aluminium hydride	Step 3 converts a ketone into a secondary alcohol. This reaction is reduction and a reducing agent like lithium aluminium hydride $LiAlH_4$ will carry out this reaction.				
7a(iv)	Step 3				Groups attached to chiral carbon (circled) <ul style="list-style-type: none"> -OH group -H group -CH₂NHCH₃ group -C₆H₃(OH)₂ group 	

7b	0.15mg	500ppm = 500mg per litre ∴ 1 litre of adrenaline solution contains 500mg adrenaline 1000cm ³ of adrenaline solution contains 500mg adrenaline 0.3cm ³ of adrenaline solution contains 500mg adrenaline × 0.3/1000 = 0.15mg				
7c	Diagram showing:	Labelled start positions of the extract and pure samples on a horizontal line. This line must be above the level of the solvent.				
8a	0.97	In Step 1, 13% ethanol is produced 13cm ³ ethanol and 87cm ³ water 1cm ³ ethanol = 0.79g ∴ 13cm ³ ethanol = 10.27g 1cm ³ water = 1.00g ∴ 87cm ³ ethanol = 87.00g $d = \frac{m_1 + m_2}{V} = \frac{10.27 + 87.00}{100} = 0.9727 \text{ g cm}^3$				
8b	One answer from:	Boiling points are similar	(some) water evaporates at ethanol's boiling point	Any mention of attraction or forces between water and ethanol		
8c	One answer from:	Water molecules are smaller than ethanol molecules	ethanol molecules are too large to pass through	ethanol molecules are larger than water molecules	water molecules pass through but ethanol molecules cannot	water molecules are trapped in the sieve
8d	Open Question to include:	3 mark answer Demonstrates a <u>good</u> understanding of the chemistry involved. A good comprehension of the chemistry has provided in a logically correct, including a statement of the principles involved and the application of these to respond to the problem.		2 mark answer Demonstrates a <u>reasonable</u> understanding of the chemistry involved, making some statement(s) which are relevant to the situation, showing that the problem is understood.		1 mark answer Demonstrates a <u>limited</u> understanding of the chemistry involved. The candidate has made some statement(s) which are relevant to the situation, showing that at least a little of the chemistry within the problem is understood.
9a	(Base-induced) Elimination	Elimination reactions involve the removal of a small water leaving behind a C=C double bond. Elimination reactions are the opposite reactions to addition reaction.				
9b	Mechanism showing:	<div style="display: flex; justify-content: space-around;"> <div style="text-align: center;"> 1st Mark Curly arrow from double bond to H⁺ </div> <div style="text-align: center;"> 2nd Mark Correct carbocation drawn </div> <div style="text-align: center;"> 3rd Mark Curly arrow showing water attacking carbocation </div> <div style="text-align: center;"> 3rd Mark Curly arrow showing hydrogen being removed from water </div> </div>				
9c			Multiplet has 6 peaks Conclusion: 5 Hydrogen atoms on adjacent carbons Chemical Shift at 3.7ppm Conclusion: Group has oxygen in Alcohol/ether			
9d(i)	Rate = k [C ₄ H ₉ Br] [OH ⁻]	Order of 1-bromobutane is [C ₄ H ₉ Br] ¹		Rate = k [C ₄ H ₉ Br] ¹ [OH ⁻] ¹		
		Order of hydroxide ion is [OH ⁻] ¹		∴ Rate = k [C ₄ H ₉ Br] [OH ⁻]		

9d(ii)	$1.32 \times 10^{-4} \text{ l mol}^{-1} \text{ s}^{-1}$	$\text{rate} = k[\text{C}_4\text{H}_9\text{Br}][\text{OH}^-]$ $k = \frac{\text{rate}}{[\text{C}_4\text{H}_9\text{Br}] \times [\text{OH}^-]}$ $= \frac{3.3 \times 10^{-6} \text{ mol l}^{-1} \text{ s}^{-1}}{0.25 \text{ mol l}^{-1} \times 0.10 \text{ mol l}^{-1}}$ $= 1.32 \times 10^{-4} \text{ l mol}^{-1} \text{ s}^{-1}$			
10a	Delocalised electrons	Benzene rings have six delocalised electrons which provide the stability in a benzene molecule.			
10b	$1s^2 2s^2 2p^6 3s^2 3p^6 3d^7$	Co atom: $1s^2 2s^2 2p^6 3s^2 3p^6 3d^7 4s^2$ ∴ Co^{2+} ion: $1s^2 2s^2 2p^6 3s^2 3p^6 3d^7$			
10c(i) PART A	All four sections required:	solutions of known concentration are prepared	suitable filter/wavelength/colour used	Mention of a blank or solvent only measurement	Absorbance/transmittance of each solution is measured/plotted
10c(i) PART B	Both sections required:	The absorbance/transmittance of the unknown is measured		mention of using the graph to turn unknown's absorbance/transmittance back into concentration	
10c(ii)	Propan-1-ol	Propanal is an aldehyde and will reduce to form the primary alcohol propan-1-ol. Other reduction reactions: carboxylic acid \longrightarrow aldehyde \longrightarrow primary alcohol ketone \longrightarrow secondary alcohol			
10d	1	Every bond in the decamethylcobaltocene is identical i.e. aromatic carbon with a methyl $-\text{CH}_3$ group attached			