



# JABchem



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# Past Papers Higher Chemistry

# 2015 Marking Scheme

Grade Awarded	Mark Required		% candidates achieving grade
	(/120)	%	
A	82	68.3%	22.0%
B	68	56.7%	24.9%
C	54	45.0%	26.7%
D	47	39.2%	10.5%
No award	<47	<39.2%	15.9%

Section:	Multiple Choice	Extended Answer	Assignment
Average Mark:	13.7 /20	40.5 /80	12.1 /20

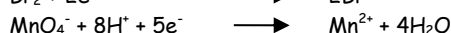
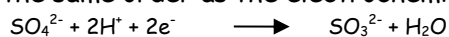
# 2015 CfE Higher Chemistry Marking Scheme

MC Qu	Answer	% Pupils Correct	Reasoning
1	D	84	<input checked="" type="checkbox"/> A Neon is a noble gas with a full outer shell and does not need to form ions. <input checked="" type="checkbox"/> B Neon is a noble gas and is monatomic not diatomic. <input checked="" type="checkbox"/> C Neon is a noble gas and does not need to form bonds to get a full outer shell. <input checked="" type="checkbox"/> D Nitrogen, oxygen, fluorine and neon are all gases at room temperature.
2	C	89 oldH=85	<input checked="" type="checkbox"/> A First ionisation energy forms positive ions not negative ions <input checked="" type="checkbox"/> B First ionisation energy removes one electron from a gaseous neutral atom <input checked="" type="checkbox"/> C One mole of electrons is removed from one mole of atoms in the gaseous state <input checked="" type="checkbox"/> D First ionisation removes one electron from gaseous single atoms
3	D	74	<input checked="" type="checkbox"/> A Electronegativity of C = 2.5 <input checked="" type="checkbox"/> B Electronegativity of N = 3.0 ∴ most attraction for bonding electrons <input checked="" type="checkbox"/> C Electronegativity of P = 2.2 <input checked="" type="checkbox"/> D Electronegativity of Si = 1.9 ∴ least attraction for bonding electrons
4	A	94	<input checked="" type="checkbox"/> A Covalent bonding is found inside molecules not between molecules <input checked="" type="checkbox"/> B Hydrogen bonding is an intermolecular force between molecules <input checked="" type="checkbox"/> C London dispersion forces are intermolecular forces between molecules <input checked="" type="checkbox"/> D permanent dipole - permanent dipole attractions are forces between molecules
5	C	80	<input checked="" type="checkbox"/> A Br <sub>2</sub> is non-polar as both atoms in molecule have same electronegativity <input checked="" type="checkbox"/> B CO <sub>2</sub> is non-polar due to the spacial arrangement of atoms in the molecule <input checked="" type="checkbox"/> C NH <sub>3</sub> is polar due to the large electronegativity difference of the N-H bond <input checked="" type="checkbox"/> D CH <sub>4</sub> is non-polar due to the similar electronegativities within the C-H bond
6	B	63	<input checked="" type="checkbox"/> A oils are reduced when hydrogen adds across the C=C double bond in an oil <input checked="" type="checkbox"/> B the oxidation of the C=C double bonds in an oil results in rancidity <input checked="" type="checkbox"/> C hydrolysis of an oil would produce three fatty acids and glycerol <input checked="" type="checkbox"/> D oils are already unsaturated due to the presence of C=C double bonds
7	A	29	<input checked="" type="checkbox"/> A propanol does not react with NaOH but sodium propanoate would form in the neutralisation reaction between alkali (NaOH) and acid (ethanoic acid) <input checked="" type="checkbox"/> B all ethanoic acid would react with NaOH to form a salt plus water <input checked="" type="checkbox"/> C no propylethanoate would be formed as H <sup>+</sup> ions are required to form esters <input checked="" type="checkbox"/> D ester theoretically formed would be propylethanoate not ethylpropanoate
8	B	50	Oil with lowest melting point will have the highest number of C=C double bonds ∴ highest number of C=C double bonds would react with the most iodine ∴ highest iodine number
9	B	79	<input checked="" type="checkbox"/> A Glycerol formed has three -OH hydroxyl bonds but no -COOH carboxyl groups <input checked="" type="checkbox"/> B Glycerol formed has three -OH groups on a different carbon each <input checked="" type="checkbox"/> C different fatty acids produced by hydrolysis of oils, not just C <sub>17</sub> H <sub>35</sub> COOH <input checked="" type="checkbox"/> D different fatty acids produced by hydrolysis of oils, not just C <sub>17</sub> H <sub>33</sub> COOH
10	A	58	<input checked="" type="checkbox"/> A oxidation: increase in oxygen : hydrogen ratio in molecule <input checked="" type="checkbox"/> B reduction: decrease in oxygen : hydrogen ratio in molecule <input checked="" type="checkbox"/> C hydrolysis: molecules splitting into smaller molecules with H <sub>2</sub> O added at break <input checked="" type="checkbox"/> D condensation: small molecules joining to form larger molecule with H <sub>2</sub> O removed
11	D	68	<input checked="" type="checkbox"/> A secondary alcohols oxidise to ketones, tertiary alcohols do not oxidise <input checked="" type="checkbox"/> B -CHO group is found in aldehydes not ketones <input checked="" type="checkbox"/> C carboxyl -COOH groups are found in carboxylic acids not ketones <input checked="" type="checkbox"/> D ketones will not oxidise by oxidising agents e.g. Fehling's solution
12	D	48	<input checked="" type="checkbox"/> A Carvone contains C=C double bonds ∴ Carvone decolourises bromine solution <input checked="" type="checkbox"/> B Carvone contains C=C double bonds ∴ Carvone decolourises bromine solution <input checked="" type="checkbox"/> C Carvone has ketone group ∴ does not oxidise with acidified dichromate solution <input checked="" type="checkbox"/> D Carvone decolourises bromine solution and is oxidised by acidified dichromate

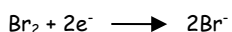
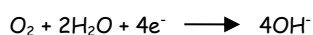
13	B	83	<input checked="" type="checkbox"/> A isoprene is a hydrocarbon ∴ no oxygen atoms <input checked="" type="checkbox"/> B isoprene is 2-methylbuta-1,3-diene C <sub>5</sub> H <sub>8</sub> <input checked="" type="checkbox"/> C isoprene is a 5 carbon unit of C <sub>5</sub> H <sub>8</sub> <input checked="" type="checkbox"/> D isoprene is a hydrocarbon ∴ no oxygen atoms
14	C	74	The -OH bond in erythromycin can be reacted with a carboxylic acid to form an ester by a condensation reaction.
15	B	69	2,2-dimethylpentan-1-ol has the molecular formula C <sub>7</sub> H <sub>15</sub> OH <input checked="" type="checkbox"/> A CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH(CH <sub>3</sub> )CH <sub>2</sub> OH has the molecular formula of C <sub>6</sub> H <sub>13</sub> OH <input checked="" type="checkbox"/> B (CH <sub>3</sub> ) <sub>3</sub> CCH(CH <sub>3</sub> )CH <sub>2</sub> OH has the molecular formula of C <sub>7</sub> H <sub>15</sub> OH ∴ isomer <input checked="" type="checkbox"/> C CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> OH has the molecular formula of C <sub>8</sub> H <sub>17</sub> OH <input checked="" type="checkbox"/> D (CH <sub>3</sub> ) <sub>2</sub> CHC(CH <sub>3</sub> ) <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> OH has the molecular formula of C <sub>8</sub> H <sub>17</sub> OH
16	A	58 oldH=64	$\Delta H_1 = \Delta H_2 + \Delta H_3 + \Delta H_4$ $\Delta H_4 = \Delta H_1 - \Delta H_2 - \Delta H_3$ $\Delta H_4 = -210 - (-50) - (-86)$ $\Delta H_4 = -74\text{kJ mol}^{-1}$ But $\Delta H$ for Z to Y = +74kJ mol <sup>-1</sup>
17	D	67 oldH=70	<input checked="" type="checkbox"/> A $2\text{I}_{(g)} + 2e^- \longrightarrow 2\text{I}^-_{(g)}$ should have $\Delta H = 2 \times -349\text{kJ} = -698\text{kJ}$ <input checked="" type="checkbox"/> B $2\text{I}_{(g)} + 2e^- \longrightarrow 2\text{I}^-_{(g)}$ should have $\Delta H = 2 \times -349\text{kJ} = -698\text{kJ}$ <input checked="" type="checkbox"/> C $\text{I}_{2(g)} \longrightarrow 2\text{I}_{(g)}$ should have $\Delta H = +243\text{kJ}$ <input checked="" type="checkbox"/> D all steps have the correct enthalpy changes: Enthalpy of sublimation $\text{I}_{2(s)} \longrightarrow \text{I}_{2(g)} \quad \Delta H = +60\text{kJ}$ Bond dissociation of I <sub>2</sub> $\text{I}_{2(g)} \longrightarrow 2\text{I}_{(g)} \quad \Delta H = +243\text{kJ}$ 2x electron affinity of iodine $2\text{I}_{(g)} + 2e^- \longrightarrow 2\text{I}^-_{(g)} \quad \Delta H = -698\text{kJ}$
18	A	89 oldH=87	<input checked="" type="checkbox"/> A at equilibrium, rate of forward reaction = rate of reverse reaction <input checked="" type="checkbox"/> B At equilibrium, the rates of the forward and reverse reactions are rarely equal <input checked="" type="checkbox"/> C Forward and reverse reactions continue at equilibrium and don't stop <input checked="" type="checkbox"/> D Adding a catalyst has no effect on the position of equilibrium
19	A	67 oldH=74	<input checked="" type="checkbox"/> A X-Y is energy difference between the reactants (R) and the activated complex <input checked="" type="checkbox"/> B Y-X would give a negative value but activation energy (E <sub>a</sub> ) must be positive <input checked="" type="checkbox"/> C Y-Z is the enthalpy change ( $\Delta H$ ) of the reverse reaction <input checked="" type="checkbox"/> D Z-Y is the enthalpy change ( $\Delta H$ ) of the forward reaction

20 C 38

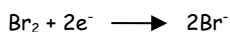
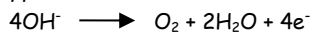
Necessary equations in the same order as the electrochemical series in data booklet



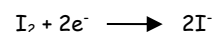
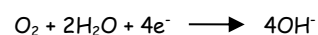
A  $\text{OH}^-$  would react with both  $\text{Br}_2$  or  $\text{I}_2$  as  $\text{OH}^-$  is above both on ECS



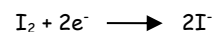
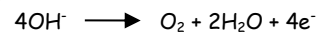
*upper reaction reverses*



*add together and cancel electrons*



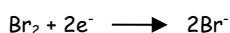
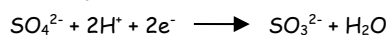
*upper reaction reverses*



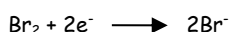
*add together and cancel electrons*



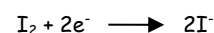
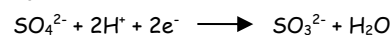
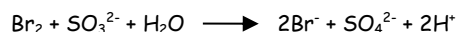
B  $\text{SO}_3^{2-}$  would react with both  $\text{Br}_2$  or  $\text{I}_2$  as  $\text{SO}_3^{2-}$  is above both on ECS



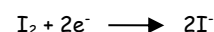
*upper reaction reverses*



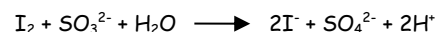
*add together and cancel electrons*



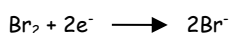
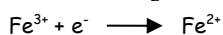
*upper reaction reverses*



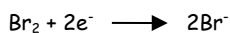
*add together and cancel electrons*



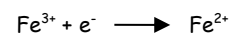
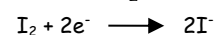
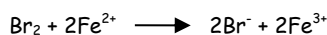
C  $\text{Fe}^{2+}$  is above  $\text{Br}_2$  so would react with  $\text{Br}_2$  but  $\text{Fe}^{2+}$  is below  $\text{I}_2$  so no reaction



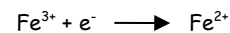
*upper reaction reverses*



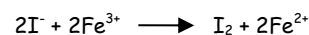
*Multiple to equalise electrons, add together & cancel electrons*



*upper reaction reverses*

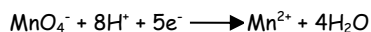
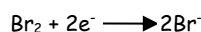


*Multiple to equalise electrons, add together & cancel electrons*

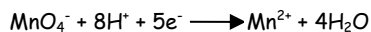


(This reaction is not the reaction of  $\text{I}_2$  with  $\text{Fe}^{2+}$ )

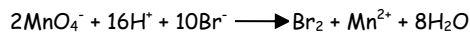
D  $\text{Mn}^{2+}$  would not react with either  $\text{Br}_2$  or  $\text{I}_2$  as  $\text{Mn}^{2+}$  is below them on ECS



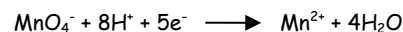
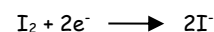
*upper reaction reverses*



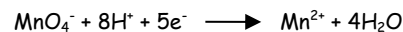
*add together and cancel electrons*



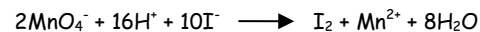
(This reaction is not the reaction of  $\text{Br}_2$  with  $\text{Mn}^{2+}$ )



*upper reaction reverses*

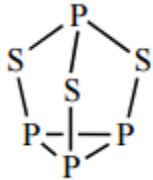
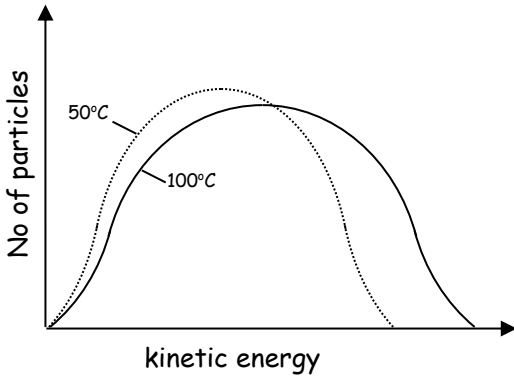
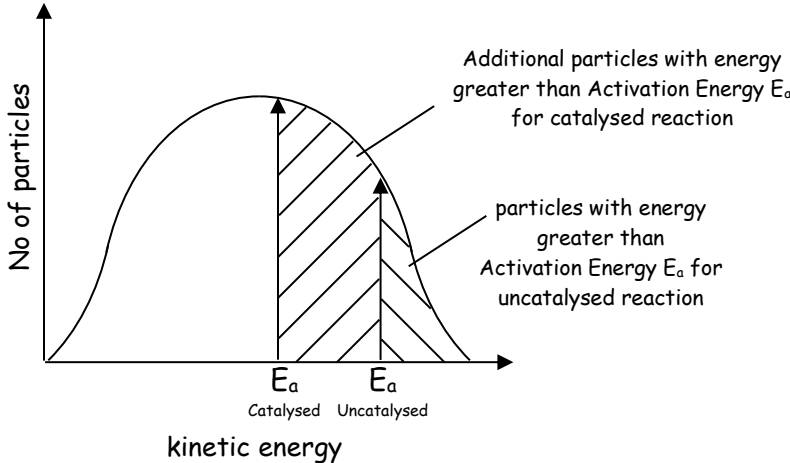


*add together and cancel electrons*

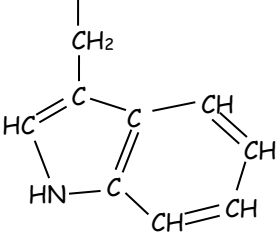
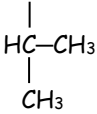
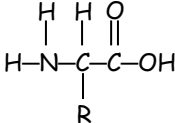
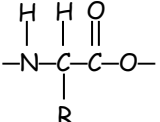
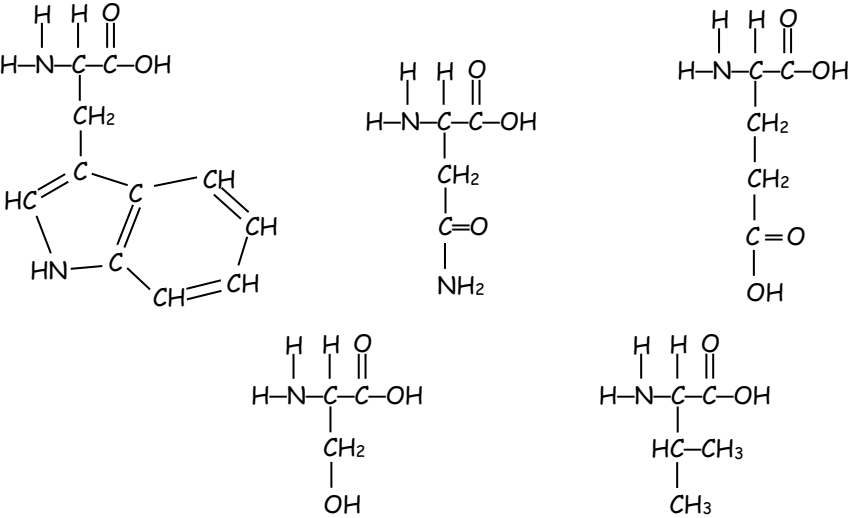


(This reaction is not the reaction of  $\text{I}_2$  with  $\text{Mn}^{2+}$ )

# 2015 CfE Higher Chemistry Marking Scheme

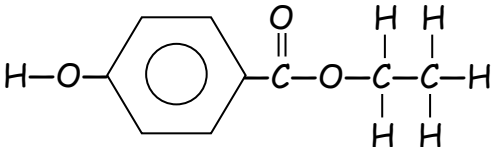
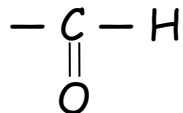
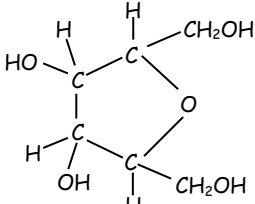
Long Qu	Answer	Reasoning						
1a	<table border="1"> <tr> <td>London dispersion forces</td> </tr> <tr> <td>covalent bonds</td> </tr> </table>	London dispersion forces	covalent bonds	<p>The covalent bonds inside the S<sub>8</sub> rings do not break as the sulphur melts. When sulphur S<sub>8</sub> melts, London dispersion forces between the S<sub>8</sub> rings must be overcome but as these are weak then sulphur has a lower melting point.</p> <p>Silicon dioxide is a covalent network. Covalent bonds must be broken before a covalent network will melt into a liquid resulting in a high melting point.</p>				
London dispersion forces								
covalent bonds								
1b(i)		<p>Any structure for P<sub>4</sub>S<sub>3</sub> that obeys the following valency rules:</p> <ul style="list-style-type: none"> <li>• 3 bonds per P atoms</li> <li>• 2 bonds per S atom</li> </ul>						
1b(ii)	Increased nuclear attraction/charge or more protons in sulphur nucleus	Sulphur and phosphorus are in the same period of the periodic table and the same shell is being filled with electrons. Sulphur has 16 protons which pull in the outer shell of the sulphur atom further than the 15 protons in a phosphorus nucleus would do. Electrons closer to the nucleus are harder to remove.						
1b(iii)	Answer to include:	<table border="1"> <tr> <td>1<sup>st</sup> Mark:</td> <td>Forces are stronger between sulphur than between phosphorus</td> </tr> <tr> <td>2<sup>nd</sup> Mark:</td> <td>London dispersion forces are the forces between the molecules.</td> </tr> <tr> <td>3<sup>rd</sup> Mark:</td> <td>Forces are stronger between Sulphur molecules of 8 atoms than between phosphorus molecules of 4 atoms.</td> </tr> </table>	1 <sup>st</sup> Mark:	Forces are stronger between sulphur than between phosphorus	2 <sup>nd</sup> Mark:	London dispersion forces are the forces between the molecules.	3 <sup>rd</sup> Mark:	Forces are stronger between Sulphur molecules of 8 atoms than between phosphorus molecules of 4 atoms.
1 <sup>st</sup> Mark:	Forces are stronger between sulphur than between phosphorus							
2 <sup>nd</sup> Mark:	London dispersion forces are the forces between the molecules.							
3 <sup>rd</sup> Mark:	Forces are stronger between Sulphur molecules of 8 atoms than between phosphorus molecules of 4 atoms.							
2a	45-46	<p>From graph: rate = 0.0022 s<sup>-1</sup></p> $\text{Rate} = \frac{1}{\text{time}} \therefore \text{time} = \frac{1}{\text{rate}} = \frac{1}{0.022} = 45\text{s}$						
2b(i)	Curve and peak drawn to left of original curve							
2b(ii)	Line drawn to the left of the E <sub>a</sub> line.	 <p>Additional particles with energy greater than Activation Energy E<sub>a</sub> for catalysed reaction</p> <p>particles with energy greater than Activation Energy E<sub>a</sub> for uncatalysed reaction</p>						
3a(i)	Diagram showing:	<p>1mark: workable apparatus for passing the steam through the strawberry gum (steam must pass through the strawberry gum leaves not just pass over)</p> <p>1mark: workable apparatus for condensing the steam and essential oil</p>						
3a(ii)	(fractional) distillation or chromatography	<p>Distillation separates chemicals with different boiling points.</p> <p>Chromatography separates chemicals due to difference in polarity or size.</p>						

3b(i)	Answer to include:	$\text{gfm cinnamic acid} = 148\text{g}$ $\text{no. of mol} = \frac{\text{mass}}{\text{gfm}} = \frac{6.5\text{g}}{148\text{g}} = 0.0439\text{mol}$ $\text{gfm methanol} = 32\text{g}$ $\text{no. of mol} = \frac{\text{mass}}{\text{gfm}} = \frac{2.0\text{g}}{32\text{g}} = 0.0625\text{mol}$ $\begin{array}{ccccccc} \text{cinnamic acid} & + & \text{methanol} & \longrightarrow & \text{methyl cinnamate} & + & \text{water} \\ 1\text{mol} & & 1\text{mol} & & 1\text{mol} & & 1\text{mol} \\ 0.0439\text{mol} & & 0.0439\text{mol} & & & & \\ & & (+0.0186\text{mol leftover}) & & & & \end{array}$ <p>Cinnamic acid is the limiting reactant as all 0.0439mol of cinnamic acid is used up in the reaction. 0.0186mol methanol is left over in the reaction.</p>								
3b(ii) Part A	52%	$\begin{array}{ccc} \text{Cinnamic acid} & + & \text{methanol} \longrightarrow \text{methyl cinnamate} & + & \text{water} \\ 1\text{mol} & & & & 1\text{mol} \\ 148\text{g} & & & & 162\text{g} \\ 6.5\text{g} & & & & 162\text{g} \times \frac{6.5}{148} \\ & & & & = 7.1\text{g (theoretical)} \end{array}$ $\% \text{ yield} = \frac{\text{actual}}{\text{theoretical}} \times 100 = \frac{3.7}{7.1} \times 100 = 52\%$								
3b(ii) Part B	£24.59	$\begin{array}{ccc} \text{Cinnamic acid} & + & \text{methanol} \longrightarrow \text{methyl cinnamate} & + & \text{water} \\ 6.5\text{g} & & & & 3.7\text{g (at 52\% yield)} \\ 6.5\text{g} \times \frac{100}{3.7} & & & & 100\text{g} \\ = 175.7\text{g} & & & & \\ 250\text{g cinnamic acid} & \text{costs} & \text{£35.00} & & \\ 175.7\text{g cinnamic acid} & \text{costs} & \text{£35.00} \times \frac{175.7}{250} & = & \text{£24.59} \end{array}$								
4a	Citronellol or geraniol or anisyl alcohol	Only peaks B (citronellol), C (geraniol) and E (anisyl alcohol) appear on all three chromatograms								
4b	Counterfeit perfumes have lower concentrations of compounds	The area under each peak is proportional to the quantity of that chemical in the sample. The smaller the peak, the lower the concentration of that chemical.								
4c(i)	Inert/does not react with molecules	Gas chromatography needs a carrier gas to flow through the separating column. The sample compounds pass through the column at different rates dependent on their attraction to the mobile phase (the helium gas) or the stationary phase (the contents of the column)								
4c(ii)	Size of molecules or temperature of column	The smaller molecules will pass through the column more quickly than larger molecules. Increasing the temperature in the column would increase the kinetic energy of the particles and they would pass through the column more quickly.								
4d(i)	Terpenes	Terpenes are formed when multiple units of isoprene join together.								
4d(ii) Part A	3,7-dimethylocta-1,6-dien-3-ol	$\text{3,7-dimethylocta-1,6-dien-3-ol}$ $\begin{array}{ccc} \text{-CH}_3 \text{ groups on} & \text{8 carbons with} & \text{-OH group} \\ \text{carbon C}_3 \text{ and C}_7 & \text{2xC=C on C}_1 \text{ and C}_6 & \text{on C}_3 \end{array}$								
4d(ii) Part B	3 carbons attached to the Carbon with the -OH group	<table border="1"> <thead> <tr> <th>Alcohol</th> <th>Description</th> </tr> </thead> <tbody> <tr> <td>Primary</td> <td>1 carbon directly attached to the carbon with the -OH group</td> </tr> <tr> <td>Secondary</td> <td>2 carbons directly attached to the carbon with the -OH group</td> </tr> <tr> <td>Tertiary</td> <td>3 carbons directly attached to the carbon with the -OH group</td> </tr> </tbody> </table>	Alcohol	Description	Primary	1 carbon directly attached to the carbon with the -OH group	Secondary	2 carbons directly attached to the carbon with the -OH group	Tertiary	3 carbons directly attached to the carbon with the -OH group
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4e	1.7g	<p>1kg body mass allowed 0.10mg coumarin</p> <p>75kg body mass allowed 0.10mg coumarin <math>\times \frac{75}{1}</math></p> <p style="text-align: center;">= 7.5mg coumarin = 0.0075g coumarin</p> <p>4.4g coumarin contained in 1000g cinnamon powder</p> <p>0.0075g coumarin contained in 1000g cinnamon powder <math>\times \frac{0.0075}{4.4}</math></p> <p style="text-align: center;">= 1.70g cinnamon powder</p>								
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6a	Heat breaks hydrogen bonds	Globular proteins have a very specific shape and hydrogen bonding is important in holding these proteins chains together. Hydrogen bonds break on heating, the shape of the protein unravels and won't return to the original shape when the temperature decreases again.		
6b(i)	One structure from:			
		These side groups are non-polar and predominantly hydrocarbon. Hydrocarbons are hydrophobic in nature. The other side groups shown are hydrophilic as they have polar groups e.g. -NH <sub>2</sub> , -OH and -COOH		
6b(ii)	50.5 ± 1°C	Highest Fluorescence = 2600 units } Halfway fluorescence value = 1737.5 Lowest Fluorescence = 875 units } ∴ temperature at 1737.5 units = 50.5°C		
6c(i)	Hydrolysis	When protein is broken down into amino acids in the body by digestive enzymes, the process breaks the protein into amino acids adding a water molecule across the break point each time. This is a hydrolysis reaction.		
6c(ii) Part A	5	Amino acids are monomers with an -NH <sub>2</sub> group and a -COOH group. All amino acids have this structure with only the side group R changing from amino acid to amino acid.		
6c(ii) Part B	One amino acid structure from:		When amino acids join together by a condensation reaction and a water molecule is removed between amino acids and the repeating unit is. There are five of these units in the diagram in the question	
6c(ii) Part B	One amino acid structure from:			
7a	88.5 litres	$1\text{mol CH}_3\text{OH} = (1 \times 12) + (4 \times 1) + (1 \times 16) = 12 + 4 + 16 = 32\text{g}$ $\text{no. of mol} = \frac{\text{mass}}{\text{gfm}} = \frac{118\text{g}}{32\text{g mol}^{-1}} = 3.69\text{mol}$ $2\text{CH}_3\text{OH} + 3\text{O}_2 \longrightarrow 2\text{CO}_2 + 4\text{H}_2\text{O}$ $\begin{array}{ccc} 2\text{mol} & & 2\text{mol} \\ 3.69\text{mol} & & 3.69\text{mol} \end{array}$ $\text{Volume} = \text{no. of mol} \times \text{Molar Volume} = 3.69\text{mol} \times 24\text{ litres mol}^{-1} = 88.5\text{litres}$		
7b(i) Part A	One answer from:	Thermometer touching bottom of beaker	Thermometer is directly above flame	Temperature rise recorded would be greater than expected
7b(i) Part B	One answer from:	Distance between flame and beaker	Height of wick in burner	Same draught proofing
				Beaker made of same material e.g. copper

7b(i) Part C	-288 kJ mol <sup>-1</sup>	$E_h = c \times m \times \Delta T$ $E_h = 4.18 \times 0.1 \times 23$ $E_h = 9.614 \text{ kJ}$ <p>1mol CH<sub>3</sub>OH = (1x12)+(4x1)+(1x16) = 12+4+16 = 32g</p> $\begin{array}{ccc} 1.07\text{g} & \leftrightarrow & 9.614\text{kJ} \\ 32\text{g} & \leftrightarrow & 9.614\text{kJ} \times 32/1.07 \\ & & = -287.52 \text{ kJ mol}^{-1} \end{array}$																				
7b(ii)	0.799	density = $\frac{\text{mass}}{\text{volume}} = \frac{19.98\text{g}}{25.0\text{cm}^3} = 0.799\text{g cm}^{-3}$																				
7c(i)	Heat produced by exothermic reactions will need to be removed	The heat produced by an exothermic reaction must be removed or the system will continually rise in temperature. Endothermic reactions must have heat energy continually supplied to system for the chemical reaction to take place. The design of the reaction chamber must take account of the supply or removal of heat energy.																				
7c(ii)	+191 kJ mol <sup>-1</sup>	<table border="1"> <thead> <tr> <th colspan="2">Bond Breaking Steps</th> <th colspan="2">Bond Forming Steps</th> </tr> </thead> <tbody> <tr> <td>3x C-H bonds</td> <td>3x +412kJ = 1236kJ</td> <td>3x H-H bonds</td> <td>3x -436kJ = 1308kJ</td> </tr> <tr> <td>1x C-O bond</td> <td>1x +360kJ = 360kJ</td> <td>2x C=O bonds</td> <td>2x -743kJ = 1486kJ</td> </tr> <tr> <td>3x O-H bonds</td> <td>3x +463kJ = 1389kJ</td> <td></td> <td></td> </tr> <tr> <td>Total</td> <td>= +2985kJ</td> <td>Total</td> <td>= 2794 kJ</td> </tr> </tbody> </table> <p>Enthalpy change = +2985 - 2794 = +191kJ mol<sup>-1</sup></p> $\Delta H = \Sigma \text{Bond enthalpies for bonds broken} - \Sigma \text{Bond enthalpies for bonds formed}$ $\Delta H = 2985 - 2794$ $\Delta H = +191 \text{ kJ mol}^{-1}$	Bond Breaking Steps		Bond Forming Steps		3x C-H bonds	3x +412kJ = 1236kJ	3x H-H bonds	3x -436kJ = 1308kJ	1x C-O bond	1x +360kJ = 360kJ	2x C=O bonds	2x -743kJ = 1486kJ	3x O-H bonds	3x +463kJ = 1389kJ			Total	= +2985kJ	Total	= 2794 kJ
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8a	Flow chart complete with:	<p>1<sup>st</sup> Mark: <span style="border: 1px solid black; padding: 2px;">ammonia</span> <span style="border: 1px solid black; padding: 2px;">calcium carbonate</span></p> <p style="margin-left: 150px;"><span style="border: 1px solid black; padding: 2px;">carbon dioxide</span> <span style="border: 1px solid black; padding: 2px;">calcium oxide</span></p> <p>2nd Mark: <span style="border: 1px solid black; padding: 2px;">sodium hydrogencarbonate</span> <span style="border: 1px solid black; padding: 2px;">ammonium chloride</span></p> <p style="margin-left: 150px;"><span style="border: 1px solid black; padding: 2px;">water</span></p> <p style="margin-left: 50px;"><span style="border: 1px solid black; padding: 2px;">sodium carbonate</span></p>																				
8b	Adding Na <sup>+</sup> shifts equilibrium to right	Brine contains Na <sup>+</sup> ions. Na <sup>+</sup> ions are a reactant in the reaction. When a reactant is added to a reaction at equilibrium, the equilibrium shifts to the right to make additional products and remove the additional Na <sup>+</sup> ions.																				
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10a(i)	24 hours allows time for all zinc to react No stopper allows gas to escape from flask	The reactants need time to react so leaving them over night for 24 hours allows more than sufficient time for every particle of reactants to react and become products. If a stopper were to prevent the hydrogen gas from escaping then the pressure would build up and the flask could explode.																				
10a(ii)	One answer from:	$\underbrace{\text{zinc ions} \quad \text{impurities} \quad \text{metal ions} \quad \text{salts}}_{\text{may be present in tap water}}$																				
10b(i)	Pipette	Measuring cylinders and beakers do not measure volumes accurately enough to prepare standard solutions. A pipette is more accurate at measuring exact volumes of liquid.																				
10b(ii)	10	$\text{mass} = \text{volume} \times \text{concentration (in g l}^{-1}\text{)} = 0.010 \text{ litres} \times 1 \text{ g l}^{-1} = 0.01\text{g} = 10\text{mg}$ $\text{concentration} = \frac{\text{mass}}{\text{volume}} = \frac{10\text{mg}}{1\text{litre}} = 10\text{mg l}^{-1}$																				
10c	4.6-4.8	A line of best fit should be drawn on the graph which ignores the obvious rogue result. When the absorbance at 0.3 is extrapolated from the drawn line, a concentration of 4.6-4.8 is obtained (not 4.0 which would be obtained using the rogue point)																				
11a	Carboxyl or carboxylic acid	<table border="1"> <tbody> <tr> <td style="text-align: center;">-O-H</td> <td style="text-align: center;"> <math>\begin{array}{c} \text{O} \\    \\ -\text{C}-\text{OH} \end{array}</math> </td> <td style="text-align: center;"> <math>\begin{array}{c} \text{O} \\    \\ -\text{C}-\text{H} \end{array}</math> </td> <td style="text-align: center;"> <math>\begin{array}{c} \text{O} \\    \\ \text{C}-\text{C}-\text{C} \end{array}</math> </td> </tr> <tr> <td style="text-align: center;">hydroxyl group</td> <td style="text-align: center;">carboxyl group</td> <td style="text-align: center;">aldehyde group</td> <td style="text-align: center;">ketone group</td> </tr> </tbody> </table>	-O-H	$\begin{array}{c} \text{O} \\    \\ -\text{C}-\text{OH} \end{array}$	$\begin{array}{c} \text{O} \\    \\ -\text{C}-\text{H} \end{array}$	$\begin{array}{c} \text{O} \\    \\ \text{C}-\text{C}-\text{C} \end{array}$	hydroxyl group	carboxyl group	aldehyde group	ketone group												
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11b	Esterification or condensation	Carboxylic acids react with alcohols to make esters by a condensation reaction with water removed as they join. Condensation reactions which form esters are also known as esterification reactions.												
11c	Diagram showing:													
11d	As molecular size increases, adsorption decreases	Problem Solving: drawing a conclusion from a table of results												
12a(i)	Answer to include:	<p>1<sup>st</sup> mark: rinse the burette with (thiosulphate) solution.            2<sup>nd</sup> and 3<sup>rd</sup> marks for two of the following points:</p> <table border="1" data-bbox="555 515 1492 593"> <tbody> <tr> <td>Fill burette above the scale with thiosulphate solution</td> <td>Filter funnel to be removed</td> <td>Tap opened/some solution drained to ensure no air bubbles</td> <td>(thiosulphate) solution run into scale reading</td> <td>Reading should be made from bottom of the meniscus</td> </tr> </tbody> </table>	Fill burette above the scale with thiosulphate solution	Filter funnel to be removed	Tap opened/some solution drained to ensure no air bubbles	(thiosulphate) solution run into scale reading	Reading should be made from bottom of the meniscus							
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12a(ii)	$2I^- \rightarrow I_2 + 2e^-$	<p>Redox equation: <math>NaOCl + 2I^- + 2H^+ \rightarrow I_2 + NaCl + H_2O</math></p> <p>Oxidation step: <math>2I^- \rightarrow I_2 + 2e^-</math></p> <p>Reduction step: <math>NaOCl + 2H^+ + 2e^- \rightarrow NaCl + H_2O</math></p>												
12a(iii)	$6.20 \times 10^{-5} \text{ mol l}^{-1}$	<p><math>S_2O_3^{2-}</math> no. of mol = concentration <math>\times</math> volume = <math>0.00100 \text{ litres} \times 0.0124 \text{ mol l}^{-1} = 1.24 \times 10^{-5} \text{ mol}</math></p> $I_2 + 2Na_2S_2O_3 \rightarrow 2NaI + Na_2S_4O_6$ <p style="text-align: center;"> <math>\begin{matrix} 1\text{mol} &amp; 2\text{mol} \\ 6.20 \times 10^{-6} \text{mol} &amp; 1.24 \times 10^{-5} \text{mol} \end{matrix}</math> </p> $NaOCl + 2I^- + 2H^+ \rightarrow I_2 + NaCl + H_2O$ <p style="text-align: center;"> <math>\begin{matrix} 1\text{mol} &amp; 1\text{mol} \\ 6.20 \times 10^{-6} \text{mol} &amp; 6.20 \times 10^{-6} \text{mol} \end{matrix}</math> </p> <p>100cm<sup>3</sup> of swimming pool water contains <math>6.20 \times 10^{-6} \text{ mol NaOCl}</math>            1000cm<sup>3</sup> of swimming pool water contains <math>6.20 \times 10^{-6} \text{ mol} \times \frac{1000}{100} = 6.20 \times 10^{-5} \text{ mol l}^{-1}</math></p>												
12b	44.4 litres	<p>45 000 litres of swimming pool water requires 400cm<sup>3</sup> hypochlorite to raise by 1ppm            45 000 litres of swimming pool water requires 800cm<sup>3</sup> hypochlorite to raise by 2ppm            2500000 litres of pool water requires <math>800 \text{ cm}^3 \times \frac{2500000}{45000} = 44444 \text{ cm}^3 = 44.4 \text{ litres}</math></p>												
12c(i)	Answer to include:	<p>1<sup>st</sup> mark: ammonia is polar and trichloramine is non-polar            2<sup>nd</sup> mark: electronegativity difference is bigger in N-H bond than N-Cl bond  <math>\therefore</math> NH<sub>3</sub> is polar and NCl is non-polar</p>												
12c(ii)	Substance with unpaired electron	Free radicals are very reactive particles which have an unpaired electron. The free radical will react with a large variety of substances to achieve the pairing of the unpaired electron.												
12a(iii)	propagation	<table border="1" data-bbox="651 1496 1396 1720"> <thead> <tr> <th>Step</th> <th>Reactants (before Arrow)</th> <th>Products (after Arrow)</th> </tr> </thead> <tbody> <tr> <td>Initiation</td> <td>No free radicals on Left Hand Side</td> <td>Free radicals on Right Hand Side</td> </tr> <tr> <td>Propagation</td> <td colspan="2">Free Radicals found on both sides of arrow</td> </tr> <tr> <td>Termination</td> <td>Free radicals on Left Hand Side</td> <td>No free radicals on Right Hand Side</td> </tr> </tbody> </table>	Step	Reactants (before Arrow)	Products (after Arrow)	Initiation	No free radicals on Left Hand Side	Free radicals on Right Hand Side	Propagation	Free Radicals found on both sides of arrow		Termination	Free radicals on Left Hand Side	No free radicals on Right Hand Side
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13a		Aldehyde groups contain a carbonyl (C=O) group with a hydrogen atom attached. The carbon in the aldehyde group is always carbon number one in any numbering system assigned to an aldehyde when naming the compound.												
13b		Problem Solving Question												