



JABchem



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Past Papers Higher Chemistry

2012 Marking Scheme

Grade Awarded	Mark Required (/100)	% candidates achieving grade
A	75+	31.4%
B	62+	25.2%
C	49+	22.4%
D	42+	9.3%
No award	<42	11.7%

Section:	Multiple Choice	Extended Answer
Average Mark:	27.3 /40	36.4 /60

2012 Higher Chemistry Marking Scheme

MC Qu	Answer	% Pupils Correct	Reasoning								
1	D	76	Isotopes have <u>same</u> Atomic number but <u>different</u> Mass number Number of protons Number of neutrons								
2	C	55	<table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="width: 20%;">Reaction</th> <th>Conclusion</th> </tr> </thead> <tbody> <tr> <td>(i)</td> <td>W is the least reactive as it is the only metal which doesn't react with acid</td> </tr> <tr> <td>(ii)</td> <td>Z is most reactive as it is the only one not reduced when heated with carbon</td> </tr> <tr> <td>(iii)</td> <td>X is more reactive than Y as X displaces Y from a solution of Y nitrate</td> </tr> </tbody> </table>	Reaction	Conclusion	(i)	W is the least reactive as it is the only metal which doesn't react with acid	(ii)	Z is most reactive as it is the only one not reduced when heated with carbon	(iii)	X is more reactive than Y as X displaces Y from a solution of Y nitrate
Reaction	Conclusion										
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(ii)	Z is most reactive as it is the only one not reduced when heated with carbon										
(iii)	X is more reactive than Y as X displaces Y from a solution of Y nitrate										
3	D	33	<input checked="" type="checkbox"/> A Neon atoms are neutral and not positively charged <input checked="" type="checkbox"/> B Fluoride F ⁻ ions are negatively charged <input checked="" type="checkbox"/> C Sodium atoms are neutral and not positively charged <input checked="" type="checkbox"/> D Aluminium Al ³⁺ ions have electron arrangement 2,8 and are positively charged								
4	C	58	<input checked="" type="checkbox"/> A sodium carbonate is soluble ∴ no precipitate is formed <input checked="" type="checkbox"/> B potassium chloride is soluble ∴ no precipitate is formed <input checked="" type="checkbox"/> C magnesium carbonate is insoluble ∴ precipitate is formed <input checked="" type="checkbox"/> D potassium sulphate is soluble ∴ no precipitate is formed								
5	B	62	0.8mol of H ₂ remaining ∴ 0.2mol of H ₂ has reacted $ \begin{array}{ccc} \text{H}_2 + \text{I}_2 & \longrightarrow & 2\text{HI} \\ \begin{array}{c} 1\text{mol} \\ 0.2\text{mol} \end{array} & & \begin{array}{c} 2\text{mol} \\ 0.4\text{mol} \end{array} \end{array} $								
6	D	79	<input checked="" type="checkbox"/> A Activation Energy E _a is not altered by changes in temperature <input checked="" type="checkbox"/> B The Enthalpy Change ΔH is independent of the temperature it takes place at <input checked="" type="checkbox"/> C Activation Energy E _a is not altered by changes in temperature <input checked="" type="checkbox"/> D Increasing temperature means more collision with energy greater than E _a								
7	C	65	$ \begin{array}{ccccccc} \text{CaCO}_3 & + & 2\text{HNO}_3 & \longrightarrow & \text{Ca(NO}_3)_2 & + & \text{H}_2\text{O} & + & \text{CO}_2 \\ 1\text{mol} & & 2\text{mol} & & 1\text{mol} & & 1\text{mol} & & 1\text{mol} \\ 0.05\text{mol} & & 0.1\text{mol} & (0.1\text{mol HNO}_3 \text{ required but only } 0.08\text{mol HNO}_3 \text{ available}) & & & & & \\ 0.04\text{mol} & & 0.08\text{mol} & & 0.04\text{mol} & & 0.04\text{mol} & & 0.04\text{mol} \\ \text{(required)} & & \text{(available)} & & & & & & \end{array} $								
8	B	79	<input checked="" type="checkbox"/> A distance x represents the activation energy for the forward reaction <input checked="" type="checkbox"/> B distance y represents the enthalpy change R→P for the forward reaction <input checked="" type="checkbox"/> C distance x+y represents the activation energy for the reverse reaction <input checked="" type="checkbox"/> D distance x-y does not represent anything on this graph								
9	A	44	$ \begin{array}{ccccccc} 5\text{N}_2\text{O}_4 & + & 4\text{CH}_3\text{NHNH}_2 & \longrightarrow & 4\text{CO}_2 & + & 12\text{H}_2\text{O} & + & 9\text{N}_2 & \Delta\text{H} = -5116\text{kJ} \\ 5\text{mol} & & 4\text{mol} & & & & & & & \\ 2\text{mol} & & 4\text{mol} \times \frac{2}{5} & & & & & & & \Delta\text{H} = -5116\text{kJ} \times \frac{2}{5} \\ & & = 1.6\text{mol (2mol available)} & & & & & & & = -2046.4\text{kJ} \end{array} $								
10	C	62	<input checked="" type="checkbox"/> A Electronegativity difference = 3.0 - 2.5 = 0.5 <input checked="" type="checkbox"/> B Electronegativity difference = 3.5 - 3.0 = 0.5 <input checked="" type="checkbox"/> C Electronegativity difference = 3.0 - 3.0 = 0 <input checked="" type="checkbox"/> D Electronegativity difference = 3.0 - 2.2 = 0.8								
11	D	78	<input checked="" type="checkbox"/> A Solid has high mpt and non-conductor as solid ∴ substance likely to be ionic <input checked="" type="checkbox"/> B Solid conducts as solid ∴ substance contains metallic bonding <input checked="" type="checkbox"/> C Solid conducts as solid ∴ substance contains metallic bonding <input checked="" type="checkbox"/> D Substances with lower mpt and non-conductor as a solid is likely to be non-polar covalent and have only London dispersion forces between molecules/atoms								
12	B	68	<input checked="" type="checkbox"/> A Hydrogen has a single covalent bond in the diatomic H ₂ molecule <input checked="" type="checkbox"/> B Helium is monatomic Noble Gas element in Group 0 <input checked="" type="checkbox"/> C Nitrogen has a triple covalent bond in the diatomic N ₂ molecule <input checked="" type="checkbox"/> D Sulphur has covalent bonds within the S ₈ molecule								

13	B	88	<input checked="" type="checkbox"/> A All ionic substances must be compounds <input checked="" type="checkbox"/> B Monatomic elements cannot be found in compounds as they do not form bonds <input checked="" type="checkbox"/> C Covalent Networks can be elements (e.g. diamond) and compounds (e.g. SiO ₂) <input checked="" type="checkbox"/> D Covalent molecules can be elements (e.g. H ₂) or compounds (e.g. H ₂ O)												
14	C	54	<input checked="" type="checkbox"/> A Hexane C ₆ H ₁₄ is a non-polar hydrocarbon ∴ ionic substances do not dissolve in it <input checked="" type="checkbox"/> B Benzene C ₆ H ₆ is a non-polar hydrocarbon ∴ ionic substances do not dissolve in it <input checked="" type="checkbox"/> C Methanol is polar (due to -OH bond) and polar ionic substances can dissolve in it <input checked="" type="checkbox"/> D Tetrachloromethane CCl ₄ is non-polar due to its tetrahedral shape												
15	D	75	1mol of any gas has the same volume (at the same conditions of temp and pressure) $1\text{mol gas} = 6.02 \times 10^{23} \text{ molecules}$ $0.1\text{mol O}_2 + 0.2\text{mol CO}_2 = 0.3\text{mol gas} = 6.02 \times 10^{23} \text{ molecules} \times 0.3/10$ $= 1.806 \times 10^{23} \text{ molecules}$												
16	B	74	<input checked="" type="checkbox"/> A gfm H ₂ =2g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{2}{2} = 1 \text{ mol}$ gfm N ₂ =28g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{14}{28} = 0.5 \text{ mol}$ <input checked="" type="checkbox"/> B gfm CH ₄ =16g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{32}{16} = 2 \text{ mol}$ gfm CO ₂ =44g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{88}{44} = 2 \text{ mol}$ <input checked="" type="checkbox"/> C gfm CO=28g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{7}{28} = 0.25 \text{ mol}$ gfm O ₂ =32g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{16}{32} = 0.5 \text{ mol}$ <input checked="" type="checkbox"/> D gfm HCl=36.5g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{10}{36.5} = 0.27 \text{ mol}$ gfm SO ₂ =64.1g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{10}{64.1} = 0.16 \text{ mol}$												
17	A	61	$2\text{NO}_{(g)} + \text{O}_{2(g)} \longrightarrow 2\text{NO}_{2(g)}$ <table style="width: 100%; border-collapse: collapse;"> <tr> <td style="text-align: center;">2mol</td> <td style="text-align: center;">1mol</td> <td style="text-align: center;">2mol</td> </tr> <tr> <td style="text-align: center;">1mol</td> <td style="text-align: center;">$\frac{1}{2}$mol</td> <td style="text-align: center;">1mol</td> </tr> <tr> <td style="text-align: center;">1vol</td> <td style="text-align: center;">$\frac{1}{2}$vol</td> <td style="text-align: center;">1vol</td> </tr> <tr> <td style="text-align: center;">1litre</td> <td style="text-align: center;">$\frac{1}{2}$litre</td> <td style="text-align: center;">1litre</td> </tr> </table>	2mol	1mol	2mol	1mol	$\frac{1}{2}$ mol	1mol	1vol	$\frac{1}{2}$ vol	1vol	1litre	$\frac{1}{2}$ litre	1litre
2mol	1mol	2mol													
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18	A	90	<input checked="" type="checkbox"/> A Heptane is straight and lacks rings or branches needed for unleaded petrol <input checked="" type="checkbox"/> B Branch-chain hydrocarbons are added to unleaded petrol to prevent autoignition <input checked="" type="checkbox"/> C Cyclo-ring hydrocarbons are added to unleaded petrol to prevent auto-ignition <input checked="" type="checkbox"/> D Aromatic hydrocarbons are added to unleaded petrol to prevent auto-ignition												
19	C	94	<input checked="" type="checkbox"/> A Butane C ₄ H ₁₀ is not the main constituent of biogas <input checked="" type="checkbox"/> B Ethane C ₂ H ₆ is not the main constituent of biogas <input checked="" type="checkbox"/> C Methane CH ₄ is the main constituent of biogas <input checked="" type="checkbox"/> D Propane C ₃ H ₈ is not the main constituent of biogas												
20	A	60	<input checked="" type="checkbox"/> A Reforming: straight molecules are converted into branched/ring structures <input checked="" type="checkbox"/> B Hydrogenation: the addition of hydrogen across a C=C double bond <input checked="" type="checkbox"/> C Dehydration: the removal of water leaving a C=C double bond behind <input checked="" type="checkbox"/> D Cracking: Larger hydrocarbons are broken into smaller hydrocarbons												
21	A	56	<input checked="" type="checkbox"/> A C ₆ H ₁₄ molecule shown is 2,3-dimethylbutane and has a different structure. <input checked="" type="checkbox"/> B Same molecule: -CH ₃ methyl group sticking off C ₂ of 5 carbon chain. <input checked="" type="checkbox"/> C Same molecule: -CH ₃ methyl group sticking off C ₂ of 5 carbon chain. <input checked="" type="checkbox"/> D Same molecule: -CH ₃ methyl group sticking off C ₂ of 5 carbon chain.												
22	B	92	<table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="text-align: center;">Alcohol <small>(old name: alkanols)</small></th> <th style="text-align: center;">Aldehyde <small>(old name: alkanals)</small></th> <th style="text-align: center;">Carboxylic Acids <small>(old name: alkanonic acids)</small></th> </tr> </thead> <tbody> <tr> <td style="text-align: center; vertical-align: middle;">-OH</td> <td style="text-align: center; vertical-align: middle;"> $\begin{array}{c} \text{O} \\ \parallel \\ \text{-C} \\ \diagdown \\ \text{H} \end{array}$ </td> <td style="text-align: center; vertical-align: middle;"> $\begin{array}{c} \text{O} \\ \parallel \\ \text{-C} \\ \diagdown \\ \text{OH} \end{array}$ </td> </tr> </tbody> </table>	Alcohol <small>(old name: alkanols)</small>	Aldehyde <small>(old name: alkanals)</small>	Carboxylic Acids <small>(old name: alkanonic acids)</small>	-OH	$\begin{array}{c} \text{O} \\ \parallel \\ \text{-C} \\ \diagdown \\ \text{H} \end{array}$	$\begin{array}{c} \text{O} \\ \parallel \\ \text{-C} \\ \diagdown \\ \text{OH} \end{array}$						
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23	B	74	<input checked="" type="checkbox"/> A Ethyl ethanoate ∴ ethanol C ₂ H ₅ OH gfm=46g and ethanoic acid CH ₃ COOH gfm=60g <input checked="" type="checkbox"/> B Propyl ethanoate ∴ propanol C ₃ H ₇ OH gfm=60g and ethanoic acid CH ₃ COOH gfm=60g <input checked="" type="checkbox"/> C Methyl propanoate ∴ methanol CH ₃ OH gfm=32g and propanoic acid C ₂ H ₅ COOH gfm=74g <input checked="" type="checkbox"/> D Ethyl propanoate ∴ ethanol C ₂ H ₅ OH gfm=46g and propanoic acid C ₂ H ₅ COOH gfm=74g												
24	A	78	<input checked="" type="checkbox"/> A Benzene C ₆ H ₆ and ethyne C ₂ H ₂ have the same ratio of carbon to hydrogen <input checked="" type="checkbox"/> B Benzene is not oxidised by hot copper (II) oxide <input checked="" type="checkbox"/> C Benzene (bpt. 80°C) is less volatile than ethanal (bpt. 20°C) <input checked="" type="checkbox"/> D Benzene is resistant to addition reactions as it has no C=C double bonds												

25	C	85	$ \begin{array}{c} \text{H} \\ \\ \text{H}-\text{C}\equiv\text{C}-\text{C}-\text{H} \\ \\ \text{H} \end{array} \xrightarrow{+\text{Br}_2} \begin{array}{c} \text{Br Br H} \\ \\ \text{H}-\text{C}=\text{C}-\text{C}-\text{H} \\ \\ \text{H} \end{array} \xrightarrow{+\text{Br}_2} \begin{array}{c} \text{Br Br H} \\ \\ \text{H}-\text{C}-\text{C}-\text{C}-\text{H} \\ \\ \text{Br Br H} \end{array} $ <p style="text-align: center;"> propyne 1,2-dibromopropene 1,1,2,2-tetrabromopropane </p> <p> <input checked="" type="checkbox"/> A addition of 1st Br₂ molecule produces 1,2-dibromopropene (2 Br atoms attached) <input checked="" type="checkbox"/> B 1,1,2,2-tetrabromopropane produced not 1,1,1,2-dibromopropene <input checked="" type="checkbox"/> C 1,2-dibromopropene is the 1st addition product of propyne with bromine <input checked="" type="checkbox"/> D Br atoms are added to adjacent carbons as they add across triple/double bonds </p>												
26	D	87	<input checked="" type="checkbox"/> A pentan-2-ol is a secondary alcohol and oxidises to a ketone <input checked="" type="checkbox"/> B pentan-3-ol is a secondary alcohol and oxidises to a ketone <input checked="" type="checkbox"/> C 2-methylbutan-2-ol is a tertiary alcohol and cannot be oxidised <input checked="" type="checkbox"/> D 2,2-dimethylpropan-1-ol: primary alcohol and can be oxidised to carboxylic acids												
27	D	83	Synthesis Gas is made by the steam reforming of methane and is a mixture of carbon monoxide and hydrogen.												
28	C	57	<input checked="" type="checkbox"/> A Aluminium is a metal and has delocalised electrons which allow it to conduct <input checked="" type="checkbox"/> B poly(ethyne) is a polymer which has delocalised electrons to allow conduction <input checked="" type="checkbox"/> C poly(ethenol) is a soluble polymer which does not conduct electricity <input checked="" type="checkbox"/> D Graphite has delocalised electrons which allow it to conduct electricity												
29	B	74	<input checked="" type="checkbox"/> A mixture of peptides is missing a W-V fragment <input checked="" type="checkbox"/> B all possible fragments from peptide are in this answer <input checked="" type="checkbox"/> C mixture of peptides is missing a X-W fragment <input checked="" type="checkbox"/> D mixture of peptides is missing a Z-X fragment												
30	A	56	<table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="width: 20%;">Enthalpy</th> <th style="width: 30%;">Definition</th> <th style="width: 30%;">Equation</th> <th style="width: 20%;">ΔH</th> </tr> </thead> <tbody> <tr> <td>Formation</td> <td>The formation of one mole of a substance from its elements in their natural state</td> <td>$2\text{Al} + 1\frac{1}{2}\text{O}_2 \rightarrow \text{Al}_2\text{O}_3$</td> <td>-1670 kJ mol⁻¹</td> </tr> <tr> <td>Combustion</td> <td>Energy change for the complete combustion of one mole of a substance</td> <td>$\text{Al} + \frac{3}{4}\text{O}_2 \rightarrow \frac{1}{2}\text{Al}_2\text{O}_3$</td> <td>-835 kJ mol⁻¹</td> </tr> </tbody> </table>	Enthalpy	Definition	Equation	ΔH	Formation	The formation of one mole of a substance from its elements in their natural state	$2\text{Al} + 1\frac{1}{2}\text{O}_2 \rightarrow \text{Al}_2\text{O}_3$	-1670 kJ mol ⁻¹	Combustion	Energy change for the complete combustion of one mole of a substance	$\text{Al} + \frac{3}{4}\text{O}_2 \rightarrow \frac{1}{2}\text{Al}_2\text{O}_3$	-835 kJ mol ⁻¹
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31	B	71	<input checked="" type="checkbox"/> A Higher temperature will get you to equilibrium quicker <input checked="" type="checkbox"/> B Equilibrium explains why both reactants are still found in the mixture. <input checked="" type="checkbox"/> C methanol is left as products break down as fast as reactants are being used up <input checked="" type="checkbox"/> D there is an excess of ethanoic acid as there is only 0.1mol of methanol												
32	D	83	<input checked="" type="checkbox"/> A Catalysts do not affect the position of equilibrium <input checked="" type="checkbox"/> B Catalysts do not change the enthalpy change <input checked="" type="checkbox"/> C Catalysts do not affect the position of equilibrium <input checked="" type="checkbox"/> D Catalysts do not change the enthalpy change or the position of equilibrium												
33	A	54	<input checked="" type="checkbox"/> A OH ⁻ ions will remove H ⁺ , equilibrium moves to right ∴ more ClO ⁻ produced <input checked="" type="checkbox"/> B adding H ⁺ ions moves equilibrium to the left ∴ concentration of ClO ⁻ decreases <input checked="" type="checkbox"/> C adding Cl ⁻ ions moves equilibrium to the left ∴ concentration of ClO ⁻ decreases <input checked="" type="checkbox"/> D K ⁺ ion and SO ₄ ²⁻ ions have no effect on the position of equilibrium												
34	C	54	<input checked="" type="checkbox"/> A x100 dilution of acid ∴ pH of acid increases and stays at pH=7 <input checked="" type="checkbox"/> B x100 dilution of alkali ∴ pH of alkali decreases and stays at pH=7 <input checked="" type="checkbox"/> C both acid and alkali are diluted and become pH=7 <input checked="" type="checkbox"/> D the diluted acid and alkali both have a pH=7												
35	B	62	<input checked="" type="checkbox"/> A Hydrochloric acid is a strong acid and fully ionises. <input checked="" type="checkbox"/> B Ethanoic acid is a weak acid and only partially ionises ∴ has fewest ions <input checked="" type="checkbox"/> C Sodium chloride is fully soluble and produces many ions <input checked="" type="checkbox"/> D Sodium hydroxide is a strong alkali and fully ionises												
36	D	65	Sodium propanoate is a salt made from weak acid (propanoic acid) and a strong alkali ∴ pH of salt in solution is alkaline (pH>7)												

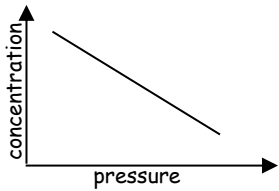
37	D	43	<p>Write down main species involved $\text{IO}_3^- \rightarrow \text{I}_2$</p> <p>Balance all atoms other than O and H $2\text{IO}_3^- \rightarrow \text{I}_2$</p> <p>Add H_2O to other side to balance O atoms $2\text{IO}_3^- \rightarrow \text{I}_2 + 6\text{H}_2\text{O}$</p> <p>Add H^+ to other side to balance H atoms $2\text{IO}_3^- + 12\text{H}^+ \rightarrow \text{I}_2 + 6\text{H}_2\text{O}$</p> <p>Add electrons to most positive side to balance charge $2\text{IO}_3^- + 12\text{H}^+ + 10\text{e}^- \rightarrow \text{I}_2 + 6\text{H}_2\text{O}$</p>																								
38	A	80	$\text{Ag}^+ + \text{e}^- \rightarrow \text{Ag} \quad \text{Au}^{3+} + 3\text{e}^- \rightarrow \text{Ag} \quad \text{Ni}^{2+} + 2\text{e}^- \rightarrow \text{Ni} \quad \text{Cu}^{2+} + 2\text{e}^- \rightarrow \text{Cu}$ <table style="width: 100%; border-collapse: collapse;"> <tbody> <tr> <td style="text-align: center;">1mol</td> <td style="text-align: center;">1mol</td> <td style="text-align: center;">3mol</td> <td style="text-align: center;">1mol</td> <td style="text-align: center;">2mol</td> <td style="text-align: center;">1mol</td> <td style="text-align: center;">2mol</td> <td style="text-align: center;">1mol</td> </tr> <tr> <td style="text-align: center;">96500C</td> <td style="text-align: center;">96500C</td> <td style="text-align: center;">289500C</td> <td style="text-align: center;">96500C</td> <td style="text-align: center;">193000C</td> <td style="text-align: center;">96500C</td> <td style="text-align: center;">193000C</td> <td style="text-align: center;">96500C</td> </tr> <tr> <td></td> <td style="text-align: center;">1mol</td> <td></td> <td style="text-align: center;">0.33mol</td> <td></td> <td style="text-align: center;">0.5mol</td> <td></td> <td style="text-align: center;">0.5mol</td> </tr> </tbody> </table>	1mol	1mol	3mol	1mol	2mol	1mol	2mol	1mol	96500C	96500C	289500C	96500C	193000C	96500C	193000C	96500C		1mol		0.33mol		0.5mol		0.5mol
1mol	1mol	3mol	1mol	2mol	1mol	2mol	1mol																				
96500C	96500C	289500C	96500C	193000C	96500C	193000C	96500C																				
	1mol		0.33mol		0.5mol		0.5mol																				
39	B	75	<input checked="" type="checkbox"/> A Fission is the splitting of a larger nucleus to form smaller nuclei <input checked="" type="checkbox"/> B Fusion is the joining up of smaller nuclei into a larger nucleus <input checked="" type="checkbox"/> C Proton Capture is when a proton is added to a nucleus <input checked="" type="checkbox"/> D Neutron Capture is when a neutron is added to a nucleus																								
40	C	53	<p>There are three different N_2 molecules formed from 2 isotopes of nitrogen:</p> $^{14}\text{N} \equiv ^{14}\text{N} \quad ^{14}\text{N} \equiv ^{15}\text{N} \quad ^{15}\text{N} \equiv ^{15}\text{N}$																								

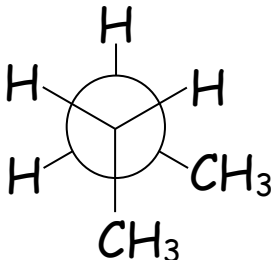
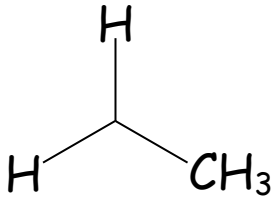
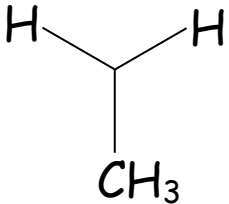
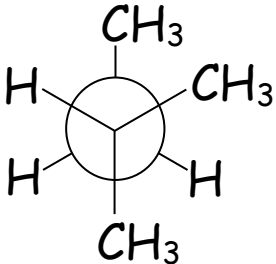
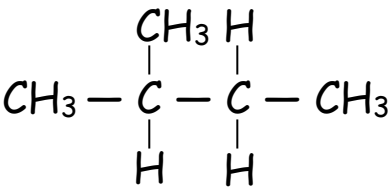
2012 Higher Chemistry Marking Scheme

Long Qu	Answer	Reasoning																											
1a	Boron or carbon	Covalent Networks are found in non-metal elements with high melting points: <table border="1" style="margin-left: auto; margin-right: auto; border-collapse: collapse; text-align: center;"> <thead> <tr> <th>Type</th> <th colspan="2">metal</th> <th colspan="6">non-metal</th> </tr> <tr> <th>Element</th> <th>Li</th> <th>Be</th> <th>B</th> <th>C</th> <th>N</th> <th>O</th> <th>F</th> <th>Ne</th> </tr> </thead> <tbody> <tr> <td>m.pt. (°C)</td> <td>181</td> <td>1287</td> <td>2075</td> <td>3825</td> <td>-210</td> <td>-219</td> <td>-220</td> <td>-249</td> </tr> </tbody> </table>	Type	metal		non-metal						Element	Li	Be	B	C	N	O	F	Ne	m.pt. (°C)	181	1287	2075	3825	-210	-219	-220	-249
Type	metal		non-metal																										
Element	Li	Be	B	C	N	O	F	Ne																					
m.pt. (°C)	181	1287	2075	3825	-210	-219	-220	-249																					
1b	Number of protons increases	Other acceptable answers: increased atomic number or greater/positive charge (pull) or greater pull on (outer) electrons																											
2a	To allow escape of gas from flask	Other acceptable answers: To prevent loss of any solution/spray/acid from flask or spurting or To stop any solids/liquids getting in/out																											
2b(i)	0.018	$\text{Rate} = \frac{\Delta \text{quantity}}{\Delta \text{time}} = \frac{0.18}{10} = 0.018 \text{ g s}^{-1}$																											
2b(ii)	0.393 (acceptable: 0.37 to 0.4)	mass of CO ₂ given off = 165.00g - 164.86g = 0.14g. gfm CO₂ = (1x12)+(2x16) = 44g $\text{no. of mol} = \frac{\text{mass}}{\text{gfm}} = \frac{0.14\text{g}}{44\text{g mol}^{-1}} = 0.00318\text{mol}$ $\text{CuCO}_3 + 2\text{HCl} \longrightarrow \text{CaCl}_2 + \text{H}_2\text{O} + \text{CO}_2$ <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">1mol 0.00318mol</div> <div style="text-align: center;">1mol 0.00318mol</div> </div> gfm CuCO ₃ = (1x63.5)+(1x12)+(3x16) = 123.5g $\text{mass} = \text{no. of mol} \times \text{gfm} = 0.00318\text{mol} \times 123.5\text{g mol}^{-1} = 0.393\text{g}$																											
3a	0.125	$\Delta H = cm\Delta T = 4.18 \times 0.5 \times 82 = -171.38\text{kJ}$ -1367kJ is released by burning 1 mol ethanol -171.38kJ released by $1 \text{ mol} \times \frac{-171.38}{-1367}$ $= 0.125\text{mol}$																											
3b	2 from:	<table border="1" style="border-collapse: collapse; text-align: center;"> <tr> <td style="padding: 2px;">heat lost to surroundings</td> <td style="padding: 2px;">incomplete combustion</td> <td style="padding: 2px;">loss (of ethanol) through evaporation</td> <td style="padding: 2px;">ethanol impure</td> </tr> </table>	heat lost to surroundings	incomplete combustion	loss (of ethanol) through evaporation	ethanol impure																							
heat lost to surroundings	incomplete combustion	loss (of ethanol) through evaporation	ethanol impure																										
4a	Equation showing:	${}_{38}^{89}\text{Sr} \rightarrow {}_{39}^{89}\text{Y} + {}_{-1}^0\text{e}$																											
4b(i)	no effect/no change	Half-life is not effected by physical state (solid/liquid/gas/solution), chemical state (atom/molecule/ion) or by changes of temperature																											
4b(ii)	5.56g	$\text{gfm } {}^{89}\text{SrCl}_2 = (1 \times 89) + (2 \times 35.5) = 89 + 71 = 160\text{g}$ 160g of ⁸⁹ SrCl ₂ contains 89g of ⁸⁹ Sr $10\text{g} \times \frac{89\text{g}}{160\text{g}} = 5.56\text{g}$																											
4c	$\frac{1}{4}$	<table border="1" style="border-collapse: collapse; text-align: center;"> <thead> <tr> <th>Time (days)</th> <th>Fraction remaining</th> </tr> </thead> <tbody> <tr> <td>0</td> <td>1</td> </tr> <tr> <td>14</td> <td>$\frac{1}{2}$</td> </tr> <tr> <td>28</td> <td>$\frac{1}{4}$</td> </tr> </tbody> </table>	Time (days)	Fraction remaining	0	1	14	$\frac{1}{2}$	28	$\frac{1}{4}$																			
Time (days)	Fraction remaining																												
0	1																												
14	$\frac{1}{2}$																												
28	$\frac{1}{4}$																												
5a	2.76×10^{21}	From graph: At voltage=20mV, volume = 110cm ³ 1 mol ethanol = 46g = 24 litres = 6.02×10^{23} molecules 0.110litres = 6.02×10^{23} molecules $\times \frac{0.110}{24}$ $= 2.76 \times 10^{21}$ molecules																											
5b	$\text{CH}_3\text{CH}_2\text{OH} + \text{O}_2 \longrightarrow \text{CH}_3\text{COOH} + \text{H}_2\text{O}$	$\begin{aligned} \textcircled{1} \quad & \text{O}_2 + 4\text{H}^+ + 4\text{e}^- \rightarrow 2\text{H}_2\text{O} \\ \textcircled{2} \quad & \text{CH}_3\text{CH}_2\text{OH} + \text{H}_2\text{O} \rightarrow \text{CH}_3\text{COOH} + 4\text{H}^+ + 4\text{e}^- \\ \textcircled{1+2} \quad & \text{O}_2 + \cancel{4\text{H}^+} + \cancel{4\text{e}^-} \rightarrow \cancel{2}^1\text{H}_2\text{O} \\ & \text{CH}_3\text{CH}_2\text{OH} + \cancel{\text{H}_2\text{O}} \rightarrow \text{CH}_3\text{COOH} + \cancel{4\text{H}^+} + \cancel{4\text{e}^-} \\ \text{overall} \quad & \text{CH}_3\text{CH}_2\text{OH} + \text{O}_2 \rightarrow \text{CH}_3\text{COOH} + \text{H}_2\text{O} \end{aligned}$																											

5c	Catalyst in different state to reactants	Catalyst		Description																
		Homogeneous	Catalyst in same state as reactants																	
		Heterogeneous	Catalyst in different state as reactants																	
6a	$\begin{array}{c} \text{H} & & \text{H} \\ & & \\ \text{H}-\text{C}-\text{S}-\text{S}-\text{C}-\text{H} \\ & & \\ \text{H} & & \text{H} \end{array}$	Methyl -CH ₃ groups must each be attached to a sulphur and this leaves a bond left for the sulphur to be attached to each other.																		
6b(i)	1.47x10 ⁻³	no. of mol Cl ₂ = volume x concentration = 0.0294 litres x 0.01 mol l ⁻¹ = 2.94x10 ⁻⁴ mol $\begin{array}{ccccccc} 4\text{Cl}_2 & + & \text{H}_2\text{S} & + & 4\text{H}_2\text{O} & \longrightarrow & \text{SO}_4^{2-} + 10\text{H}^+ + 8\text{Cl}^- \\ 4\text{mol} & & 1\text{mol} & & & & \\ 2.94 \times 10^{-4} \text{mol} & & 7.35 \times 10^{-5} \text{mol} & & & & \end{array}$ 50cm ³ water sample contains 7.35x10 ⁻⁵ mol of Cl ₂ 1000cm ³ water sample $7.35 \times 10^{-5} \text{ mol} \times \frac{1000}{50}$ $= 1.47 \times 10^{-3} \text{ mol l}^{-1}$																		
6b(ii)	2 marks for:	<u>First Mark:</u> <ul style="list-style-type: none"> Permanent dipole-permanent dipole attractions or polar-polar attractions/forces (½ mark) weak intermolecular bonds/forces (½ mark) <u>Second Mark:</u> If permanent dipole-permanent dipole attractions mentioned for 1 st Mark: <ul style="list-style-type: none"> Mention of difference in electronegativities or indication of polar bonds or indication of permanent dipole (1mark) If London dispersion forces mentioned for 1 st Mark: <ul style="list-style-type: none"> instantaneous dipoles or temporary dipoles or uneven distribution of electrons or electron wobbles (1mark) 																		
7a	w=9 z=6 y=2 z=2	<table border="1"> <tr> <td>Element</td> <td>C</td> <td>H</td> <td>N</td> <td>O</td> </tr> <tr> <td>Number</td> <td>9</td> <td>6</td> <td>2</td> <td>2</td> </tr> <tr> <td>Reasoning</td> <td>6 carbons in benzene ring 1 carbon in -CH₃ group 2 carbons in N=C=O groups</td> <td>3 hydrogens on benzene ring plus three in -CH₃ group</td> <td>2 nitrogens off benzene ring</td> <td>2 oxygens in molecule</td> </tr> </table>	Element	C	H	N	O	Number	9	6	2	2	Reasoning	6 carbons in benzene ring 1 carbon in -CH ₃ group 2 carbons in N=C=O groups	3 hydrogens on benzene ring plus three in -CH ₃ group	2 nitrogens off benzene ring	2 oxygens in molecule			
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7b	monomers have added across the (N=C) double bond	Other acceptable answers: no elimination of a small molecule (such as water) or only one product molecule formed or joined across the (N=C) double bond																		
7c	Dotted line between N and H between chains.	Hydrogen bonding takes place between:																		
		N-H bonds		O-H bonds	H-F bonds															
8a	Amide link or Peptide link	<table border="1"> <tr> <td> $\begin{array}{c} \text{O} & \text{H} \\ & \\ -\text{C}- & \text{N}- \end{array}$ Amide link </td> <td> $\begin{array}{c} \text{O} & \text{H} \\ & \\ -\text{C}- & \text{N}- \end{array}$ Peptide link </td> </tr> </table>				$\begin{array}{c} \text{O} & \text{H} \\ & \\ -\text{C}- & \text{N}- \end{array}$ Amide link	$\begin{array}{c} \text{O} & \text{H} \\ & \\ -\text{C}- & \text{N}- \end{array}$ Peptide link													
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8b	diagram showing:	$\begin{array}{c} \text{O} & & \text{NH}_2 & & \text{O} \\ & & & & \\ \text{HO}-\text{C} & - & \text{CH} & - & \text{CH}_2 & - & \text{C} \\ & & & & & & \text{OH} \end{array}$																		
8c	essential	Amino acids which cannot be made by the body are described as essential amino acids and must be consumed through our diet.																		
8d	one answer from:	Use a condenser	wet paper towel	cold finger test tube																
9a	Answer to include:	(½ mark) Use Bromine solution (½ mark) Oleic acid decolourises or stearic acid does not decolourise																		

9b	Octadec-9,12,15-trienoic acid	<h2 style="text-align: center;">Octadec-9,12,15-trienoic acid</h2> <div style="display: flex; justify-content: space-around; font-size: small;"> 18 Carbon Main Chain Position of C=C double bonds Three C=C double bonds Carboxyl -COOH group on C₁ </div>												
9c		Sodium hydroxide NaOH will neutralise carboxylic acids to form salts. The ionic portion of the molecule is ionic and water soluble. The hydrocarbon part of the molecule is non-polar and will be soluble in fat.												
10a	air	Raw Materials include: <table border="1" style="margin-left: auto; margin-right: auto; text-align: center;"> <tr> <td>air</td> <td>water</td> <td>methane (natural gas)</td> <td>sodium chloride (salt)</td> <td>calcium carbonate (chalk)</td> </tr> </table>	air	water	methane (natural gas)	sodium chloride (salt)	calcium carbonate (chalk)							
air	water	methane (natural gas)	sodium chloride (salt)	calcium carbonate (chalk)										
10b	methylmethanoate													
10c	70%	$1\text{mol HCOOH} = (2 \times 1) + (1 \times 12) + (2 \times 16) = 2 + 12 + 32 = 46\text{g}$ $1\text{mol HCONH}_2 = (3 \times 1) + (1 \times 12) + (1 \times 16) + (1 \times 14) = 3 + 12 + 16 + 14 = 45\text{g}$ $\text{HCOOH} + \text{NH}_3 \rightleftharpoons \text{HCONH}_2 + \text{H}_2\text{O}$ <table style="width: 100%; text-align: center;"> <tr> <td>1mol</td> <td></td> <td>1mol</td> </tr> <tr> <td>46g</td> <td></td> <td>45g</td> </tr> <tr> <td>1.38g</td> <td></td> <td>$45\text{g} \times \frac{1.38}{46}$</td> </tr> <tr> <td></td> <td></td> <td>= 1.35g (Theoretical)</td> </tr> </table> $\% \text{ Yield} = \frac{\text{Actual}}{\text{Theoretical}} \times 100 = \frac{0.935}{1.35} \times 100 = 70\%$	1mol		1mol	46g		45g	1.38g		$45\text{g} \times \frac{1.38}{46}$			= 1.35g (Theoretical)
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		= 1.35g (Theoretical)												
11a(i)	3-methylbutan-2-ol	<h2 style="text-align: center;">3-methylbutan-2-ol</h2> <div style="display: flex; justify-content: space-around; font-size: small;"> Methyl -CH₃ side groups on C₃ Four carbons on main chain Single bonds on main chain -OH hydroxyl group on C₂ </div>												
11a(ii)	Diagram of 2-methylpentan-1-ol													
11b(i)	$4\text{BF}_3 + 3\text{NaBH}_4$ <div style="text-align: center;">↓</div> $2\text{B}_2\text{H}_6 + 3\text{NaBF}_4$	$4\text{BF}_3 + 3\text{NaBH}_4 \longrightarrow 2\text{B}_2\text{H}_6 + 3\text{NaBF}_4$												

11b(ii)	-2168	$\begin{array}{ll} \textcircled{1} & 2\text{B} + 3\text{H}_2 \rightarrow \text{B}_2\text{H}_6 \quad \Delta\text{H} = +36 \text{ kJ mol}^{-1} \\ \textcircled{2} & \text{H}_2 + \frac{1}{2}\text{O}_2 \rightarrow \text{H}_2\text{O} \quad \Delta\text{H} = -236 \text{ kJ mol}^{-1} \\ \textcircled{3} & 2\text{B} + 1\frac{1}{2}\text{O}_2 \rightarrow \text{B}_2\text{O}_3 \quad \Delta\text{H} = -1274 \text{ kJ mol}^{-1} \\ \textcircled{1} \times -1 & \text{B}_2\text{H}_6 \rightarrow 2\text{B} + 3\text{H}_2 \quad \Delta\text{H} = -36 \text{ kJ mol}^{-1} \\ \textcircled{2} \times 3 & 3\text{H}_2 + 1\frac{1}{2}\text{O}_2 \rightarrow 3\text{H}_2\text{O} \quad \Delta\text{H} = -858 \text{ kJ mol}^{-1} \\ \textcircled{3} & 2\text{B} + 1\frac{1}{2}\text{O}_2 \rightarrow \text{B}_2\text{O}_3 \quad \Delta\text{H} = -1274 \text{ kJ mol}^{-1} \\ \text{Add} & \text{B}_2\text{H}_6 + 3\text{O}_2 \rightarrow \text{B}_2\text{O}_3 + 3\text{H}_2\text{O} \quad \Delta\text{H} = -2168 \text{ kJ mol}^{-1} \\ \textcircled{1}' + \textcircled{2}' + \textcircled{3} & \end{array}$																																																
11c	143 444 kJ	$\begin{array}{l} 1\text{mol B}_5\text{H}_9 = (5 \times 10.8) + (9 \times 1) = 54 + 9 = 63\text{g} \\ 1\text{mol } 63\text{g} = -9037\text{kJ} \\ 1000\text{g} = -9037 \text{ kJ} \times \frac{1000}{63} \\ = 143\,444 \text{ kJ} \end{array}$																																																
12a	One answer from:	<table border="1"> <tbody> <tr> <td>To allow the potato discs/catalase to reach the pH of the buffer.</td> <td>To allow buffer to soak/diffuse into the potato disc.</td> <td>To allow the enzyme/potato to reach the same pH as the surrounding solution.</td> <td>To allow the enzyme/potato to acclimatise.</td> </tr> </tbody> </table>	To allow the potato discs/catalase to reach the pH of the buffer.	To allow buffer to soak/diffuse into the potato disc.	To allow the enzyme/potato to reach the same pH as the surrounding solution.	To allow the enzyme/potato to acclimatise.																																												
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12b	hydrogen peroxide	<p>The enzyme (catalase) catalyses the breakdown of hydrogen peroxide:</p> $\text{H}_2\text{O}_2 \longrightarrow \text{H}_2\text{O} + \frac{1}{2}\text{O}_2$																																																
12c	One answer from:	<table border="1"> <tbody> <tr> <td>the enzyme is denatured</td> <td>enzymes work best at an optimum pH</td> <td>the enzyme changes its shape</td> </tr> <tr> <td>enzyme is destroyed</td> <td>too acidic for enzyme to function</td> <td>enzyme has stopped working</td> </tr> </tbody> </table>	the enzyme is denatured	enzymes work best at an optimum pH	the enzyme changes its shape	enzyme is destroyed	too acidic for enzyme to function	enzyme has stopped working																																										
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13a	1.89g	$\begin{array}{l} Q = It = 5 \times (32 \times 60) = 9600\text{C} \\ 2\text{F}^- \longrightarrow \text{F}_2 + 2\text{e}^- \\ \begin{array}{ccc} & 1\text{mol} & 2\text{mol} \\ & 38\text{g} & 193000\text{C} \\ \frac{9600}{193000} \times 38\text{g} & & 9600\text{C} \\ = 1.89\text{g} & & \end{array} \end{array}$																																																
13b(i)	exothermic or heat given out	<p>From Graph: Decrease in Temperature gives higher concentration of C_2F_4</p> <p>\therefore Decrease in temperature favours forward reaction which forms C_2F_4</p> <p>A decrease in temperature always favours the exothermic reaction</p> <p>\therefore Forward Reaction (Formation of C_2F_4) is exothermic</p>																																																
13b(ii)		<ul style="list-style-type: none"> Increasing pressure favours the pressure-reducing reaction Reverse reaction reduces pressure (2mol of gas \rightarrow 1mol of gas) <p>\therefore Increase in pressure favours reverse reaction</p> <p>\therefore Increase in pressure decreases concentration of C_2F_4</p>																																																
13c	depletion/break down of the ozone layer	Chlorofluorocarbons breaks down ozone, allowing harmful u.v. light through to cause skin cancer.																																																
14a(i)	1×10^{-5} or 10^{-5}	<table border="1"> <tbody> <tr> <td>pH</td> <td>0</td> <td>1</td> <td>2</td> <td>3</td> <td>4</td> <td>5</td> <td>6</td> <td>7</td> <td>8</td> <td>9</td> <td>10</td> <td>11</td> <td>12</td> <td>13</td> <td>14</td> </tr> <tr> <td>$[\text{H}^+]$ (mol l⁻¹)</td> <td>1</td> <td>10^{-1}</td> <td>10^{-2}</td> <td>10^{-3}</td> <td>10^{-4}</td> <td>10^{-5}</td> <td>10^{-6}</td> <td>10^{-7}</td> <td>10^{-8}</td> <td>10^{-9}</td> <td>10^{-10}</td> <td>10^{-11}</td> <td>10^{-12}</td> <td>10^{-13}</td> <td>10^{-14}</td> </tr> <tr> <td>$[\text{OH}^-]$ (mol l⁻¹)</td> <td>10^{-14}</td> <td>10^{-13}</td> <td>10^{-12}</td> <td>10^{-11}</td> <td>10^{-10}</td> <td>10^{-9}</td> <td>10^{-8}</td> <td>10^{-7}</td> <td>10^{-6}</td> <td>10^{-5}</td> <td>10^{-4}</td> <td>10^{-3}</td> <td>10^{-2}</td> <td>10^{-1}</td> <td>1</td> </tr> </tbody> </table>	pH	0	1	2	3	4	5	6	7	8	9	10	11	12	13	14	$[\text{H}^+]$ (mol l ⁻¹)	1	10^{-1}	10^{-2}	10^{-3}	10^{-4}	10^{-5}	10^{-6}	10^{-7}	10^{-8}	10^{-9}	10^{-10}	10^{-11}	10^{-12}	10^{-13}	10^{-14}	$[\text{OH}^-]$ (mol l ⁻¹)	10^{-14}	10^{-13}	10^{-12}	10^{-11}	10^{-10}	10^{-9}	10^{-8}	10^{-7}	10^{-6}	10^{-5}	10^{-4}	10^{-3}	10^{-2}	10^{-1}	1
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14a(ii)	Answer to include:	<table border="1"> <tbody> <tr> <td>1st Mark</td> <td>$\frac{1}{2}$ mark for $\text{NH}_3(\text{aq}) + \text{H}_2\text{O}(\text{l}) \rightleftharpoons \text{NH}_4^+(\text{aq}) + \text{OH}^-(\text{aq})$</td> </tr> <tr> <td>2nd Mark</td> <td>$\frac{1}{2}$ mark for OH^- ions are removed by pairing up with NH_4^+ ions</td> </tr> <tr> <td>1st Mark</td> <td>$\frac{1}{2}$ mark for $\text{H}_2\text{O}(\text{l}) \rightleftharpoons \text{H}^+(\text{aq}) + \text{OH}^-(\text{aq})$</td> </tr> <tr> <td>2nd Mark</td> <td>$\frac{1}{2}$ mark for water molecule dissociates increasing H^+ ion concentration</td> </tr> </tbody> </table>	1 st Mark	$\frac{1}{2}$ mark for $\text{NH}_3(\text{aq}) + \text{H}_2\text{O}(\text{l}) \rightleftharpoons \text{NH}_4^+(\text{aq}) + \text{OH}^-(\text{aq})$	2 nd Mark	$\frac{1}{2}$ mark for OH^- ions are removed by pairing up with NH_4^+ ions	1 st Mark	$\frac{1}{2}$ mark for $\text{H}_2\text{O}(\text{l}) \rightleftharpoons \text{H}^+(\text{aq}) + \text{OH}^-(\text{aq})$	2 nd Mark	$\frac{1}{2}$ mark for water molecule dissociates increasing H^+ ion concentration																																								
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14b	Answer to include:	<p>Description of a large volume/number of moles of gas is produced <u>or</u></p> <p>There is an increase in the number of moles of gas <u>or</u></p> <p>Oxygen gas is produced which can support combustion <u>or</u></p> <p>It is an oxidising agent</p>																																																

15a(i)	To keep the current/amps constant	The current must be kept constant throughout the experiment as a constant current value must be used in the equation $Q=It$	
15a(ii)	Current ($\frac{1}{2}$) Time ($\frac{1}{2}$)	The Current (measured in Amps) and the Time (measured in seconds) are used in the equation $Q=It$ to calculate the Charge (measured in Coulombs)	
15b(i)	SO_2 can be recycled/reused back into step 1	Recycling of chemicals in a manufacturing process can reduce costs of chemical used in the process.	
15b(ii)	$\text{H}_2\text{O} \rightarrow \text{H}_2 + \frac{1}{2}\text{O}_2$	<p> $\textcircled{1} \quad \text{I}_2 + \text{SO}_2 + 2\text{H}_2\text{O} \rightarrow 2\text{HI} + \text{H}_2\text{SO}_4$ $\textcircled{2} \quad \quad \quad \quad \quad 2\text{HI} \rightarrow \text{I}_2 + \text{H}_2$ $\textcircled{3} \quad \quad \quad \quad \quad \text{H}_2\text{SO}_4 \rightarrow \text{SO}_2 + \text{H}_2\text{O} + \frac{1}{2}\text{O}_2$ </p> <p>add $\text{I}_2 + \text{SO}_2 + 2\text{H}_2\text{O} \rightarrow 2\text{HI} + \text{H}_2\text{SO}_4$ $2\text{HI} \rightarrow \text{I}_2 + \text{H}_2$ $\text{H}_2\text{SO}_4 \rightarrow \text{SO}_2 + \text{H}_2\text{O} + \frac{1}{2}\text{O}_2$</p> <p>$\textcircled{1}+\textcircled{2}+\textcircled{3}$ final equation $\text{H}_2\text{O} \rightarrow \text{H}_2 + \frac{1}{2}\text{O}_2$</p>	
16a			
		Carbon closest to eye has $-\text{CH}_3$ methyl group at 4 o'clock position	Carbon furthest from eye has $-\text{CH}_3$ methyl group at 6 o'clock position
16b	2-methylbutane		
		Newman Projection of molecule	Structural formula of molecule