

# GCE

# **Chemistry B**

Unit H433/01: Fundamentals of chemistry

Advanced GCE

## Mark Scheme for June 2018

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This mark scheme is published as an aid to teachers and students, to indicate the requirements of the examination. It shows the basis on which marks were awarded by examiners. It does not indicate the details of the discussions which took place at an examiners' meeting before marking commenced.

All examiners are instructed that alternative correct answers and unexpected approaches in candidates' scripts must be given marks that fairly reflect the relevant knowledge and skills demonstrated.

Mark schemes should be read in conjunction with the published question papers and the report on the examination.

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Annotations available in RM Assessor

Annotation	Meaning
<ul> <li>Image: A set of the set of the</li></ul>	Correct response
×	Incorrect response
	Omission mark
000	Benefit of doubt given
CON	Contradiction
12	Rounding error
BT	Error in number of significant figures
	Error carried forward
<b>—</b>	Level 1
<u></u>	Level 2
<b>11</b>	Level 3
HEAD .	Benefit of doubt not given
SET N	Noted but no credit given
Г.	Ignore

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

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

#### Subject-specific Marking Instructions

#### INTRODUCTION

Your first task as an Examiner is to become thoroughly familiar with the material on which the examination depends. This material includes:

- the specification, especially the assessment objectives
- the question paper
- the mark scheme.

You should ensure that you have copies of these materials.

You should ensure also that you are familiar with the administrative procedures related to the marking process. These are set out in the OCR booklet **Instructions for Examiners**. If you are examining for the first time, please read carefully **Appendix 5 Introduction to Script Marking: Notes for New Examiners**.

Please ask for help or guidance whenever you need it. Your first point of contact is your Team Leader

### Section A

Q	Кеу	Mark	
1	C	1	
2	D	1	
3	С	1	
4	В	1	
5	В	1	
6	С	1	
7	C	1	
8	C	1	
9	D	1	
10	B C	1	
11	C	1	
12	C	1	
13	Α	1	
14	В	1	
15 16	В	1	
16	C	1	
17	В	1	
18	В	1	
19	Α	1	
20	В	1	
21	D	1	
22	C	1	
23 24	В	1	
24	В	1	
25	Α	1	
26	В	1	
27	Α	1	
28	Α	1	
29 30	D	1	
30	C	1	
		30	

Question		on	Answer	Marks	Guidance
31	(a)		FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer = 63(%) or rounds to 63.0(%) award 3 marks Moles of $C_{12}H_{26} = 1.5 \times 10^6/170 \ (=8.824 \times 10^3) \checkmark$ Expected yield of $C_6H_{12} = 8.824 \times 10^3 \times 86 \ (=7.589 \times 10^5 \text{g or } 758.9 \text{ kg}) \checkmark$ % yield = 478 x100/758.9 = 63.0(%) (2 or more sf) ✓	3	<ul> <li>ALLOW alternative method:</li> <li>Moles of hexane =478000/86 = (5.558 x 10<sup>3</sup>) ✓</li> <li>% yield = 5.558 x 10<sup>3</sup>x 100/8.824 x 10<sup>3</sup> = 63.0 ✓</li> <li>A correctly rounded answer to 1sf scores 1</li> <li>If units incorrectly converted ALLOW ECF for second mark</li> </ul>
	(b)	(i)	Set up: burning fuel under a container of water OR measure the temperature <u>increase</u> of water ✓	1	
		(ii)	Find energy transferred to water using Q= mc∆T. AND Find energy that would be transferred per mole of fuel. ✓	1	Must make a comment about how the moles are obtained (i.e. using the mass of fuel burnt)
		(iii)	Any two from:	2	
			Have a lid on the container of water to reduce heat loss/stop water evaporating $\checkmark$		
			Use draught excluders <b>OR</b> insulate sides of calorimeter $\checkmark$		
			Allow enough air/oxygen to reach flame to minimise incomplete combustion <b>OR</b> Move burner closer to calorimeter ✓		ALLOW well ventilated
			Cover the wick of the burner when it is not in use to reduce evaporation of the fuel $\checkmark$		
			Use a bomb calorimeter ✓		
			Use copper calorimeter instead of beaker $\checkmark$		
			Make sure thermometer is not in contact with bottom of beaker $\checkmark$		
			Stir to improve heat distribution ✓		
	(c)		FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer = -4161 (kJ mol <sup>-1</sup> ) award 2 marks	2	<b>ALLOW ECF</b> from incorrect cycle as long as some working is shown

Question	Answer	Marks	Guidance
	$\Delta_c H^{\Theta}$ hexane = (6 x -393) + (7 x -286) – (-199) (expression must be correct) <b>OR</b> shown on an appropriate cycle $\checkmark$ – 4161 (kJ mol <sup>-1</sup> ) $\checkmark$		ALLOW -4160 (3sf based on question data) 2358 + 2002199 = -4161 -480 and a cycle scores 1 (+) 4161 scores 1
(d)	$ \begin{array}{cccccccc} H & H & H \\ H & -C & -C & -C & -H \\ H & O & H \\ H & \checkmark & & & \\ \end{array} $ <sup>13</sup> C spectrum has only 2 peaks so only 2 carbon environments $\checkmark$	2	ALLOW OH
(e)	Acidified potassium/sodium dichromate <b>AND</b> heat/high temperature ✓	1	<b>IGNORE</b> reflux or distil <b>IGNORE</b> dichromate or $Cr_2O_7^{2-}$ alone
(f)	Dipole $\checkmark$ , both curly arrows $\checkmark$ intermediate and curly arrow and product $\checkmark$ Nucleophilic addition $\checkmark$	4	Curly arrow on carbonyl must start at double bond and end on oxygen atom. Other curly arrows must start either at lone pair or negative charge and point either to atom attacked or bond between atoms. <b>ALLOW</b> dipole and movement of electrons to O for 1 mark , then C+ intermediate and attack by CN <sup>-</sup> for the second mark Intermediate and final product must have correct bonds (i.e. not through the N atom)
(g)	Please refer to the marking instructions on page 5 of this mark scheme for guidance on how to mark this question.Level 3 (5–6 marks)Deduces correct structure with detailed evidence referring to all	6	Indicative scientific points may include: Infrared spectrum: C=O as strong absorbance at approx 1750 cm <sup>-1</sup> No O-H from carboxylic acid or alcohol

Question	Answer	Marks	Guidance
	three spectra. There is a well-developed line of reasoning which is clear and logically structured. The information presented is relevant and substantiated. Level 2 (3–4 marks) Deduces correct structure using some evidence. OR		C-H at approx. 2950 cm <sup>-1</sup> possibly ester NMR: 5 proton environments as 5 peaks $\delta = 0.9, 1.1, 1.6$ H-CR. $\delta = 2.3$ HC-C=O $\delta = 4.0$ HC-O Splitting:
	<ul> <li>Deduces compound A is an ester with evidence from at least two spectra.</li> <li>OR</li> <li>Gives detailed analysis of three spectra while failing to determine the structure of compound A.</li> <li>There is a line of reasoning presented with some structure. The information presented is relevant and supported by some evidence.</li> </ul>		<ul> <li>0.9, 1.1 and 4.0 triplets so 2 protons attached to adjacent C/ CH<sub>3</sub>-CH<sub>2</sub></li> <li>2.3 quartet so 3 protons attached to adjacent C/ CH<sub>2</sub>-CH<sub>3</sub></li> <li>1.6 multiplet, several protons attached to adjacent C, possibly CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub></li> <li>Mass Spectrum: Mol mass is 116</li> <li>Extra detail</li> </ul>
	Level 1 (1–2 marks) Gives some evidence from two spectra. There is an attempt at a logical structure with a line of reasoning. The information is in the most part relevant.		Sensible discussion of at least 1 fragment e.g. peak at 87 loss of $CH_3CH_2$ or peak at 73 loss of $CH_3CH_2CH_2$ or peak at 57 due to $CH_3CH_2C=O^+$ <b>OR</b> 116 – 6C = 44 (2O) possibly ester Structure is
	0 marks No response or no response worthy of credit		
	Total	22	

Q	uesti	on	Answer	Marks	Guidance
32	(a)		The 3D shape <b>OR</b> the shape produced by the folding of the protein molecule ✓	1	
	(b)		Any two from: ✓ Instantaneous dipole-induced dipole hydrogen bonds ionic bonds covalent bonds	1	<b>IGNORE</b> specific groups mentioned after bond types.
	(c)	(i)	$\begin{array}{c c} -\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\text{CH}_2-\overrightarrow{C}-\text{NH}-\text{CH}_2-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{CH}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\overrightarrow{C}-\text{NH}-\overrightarrow{C}-\overrightarrow{C}-\overrightarrow{C}-\overrightarrow{C}-\overrightarrow{C}-\overrightarrow{C}-\overrightarrow{C}-C$	1	ALLOW C or CH ringed Extra carbons ringed are CON
		(ii)	$\begin{array}{c} \begin{array}{c} & & & & \\ & +H_{3}N - CH - C - OH \\ & & +H_{3}N - CH_{2} - C - OH \\ & & & \\ & $	4	<ul> <li>ALLOW ECF if all the NH<sub>3</sub><sup>+</sup> groups are not protonated</li> <li>IGNORE Cl<sup>-</sup>ions.</li> <li>IGNORE number of moles of aminoethanoic acid.</li> <li>Structures with deprotonated carboxylate groups score 0 (no ECF)</li> <li>Extra incorrect structures CON a correct one</li> </ul>

Question	Answer	Marks	Guidance
(d)	<ul> <li>Please refer to the marking instructions on page 5 of this mark scheme for guidance on how to mark this question.</li> <li>Level 3 (5–6 marks)</li> <li>Gives a clear and detailed account of all three parts, including most of the points listed.</li> <li>There is a well-developed line of reasoning which is clear and logically structured. The information presented is relevant and substantiated.</li> <li>Level 2 (3–4 marks)</li> <li>Gives an outline account of all three parts OR gives a detailed account of two parts.</li> <li>There is a line of reasoning presented with some structure. The information presented is relevant and supported by some evidence.</li> <li>Level 1 (1–2 marks)</li> <li>Makes some relevant points</li> <li>There is an attempt at a logical structure with a line of reasoning. The information is in the most part relevant.</li> <li>O marks</li> <li>No response or no response worthy of credit</li> </ul>	6	Indicative scientific points may include: Developing • spray with ninhydrin ALLOW UV light • dry (in an oven/ fume cupboard) Chromatogram • Start line • Starting dot of hydrolysate OR Dots of suspected hydrolysis products for reference • (four spots above) • Spots level with suspected hydrolysis products • Mark position of solvent front • Lid • Stop when solvent gets near the top of the paper Analysis • Measure Rf values of spots • Rf = distance moved by spot/distance moved by solvent front • Look up Rf values for the three amino acids • Compare with measured values OR Compare R <sub>f</sub> values with reference amino acids IGNORE use of tlc plate instead of paper
	Total	13	

Q	uestic	on		Answer		Marks	Guidance
33	33 (a)		NaOH(aq) AND (Heat under) Reflux ✓				ALLOW warm for reflux
		(ii)	Acidify (until neutral) ✓			2	ALLOW any dilute named acid
			Filter off C OR evaporate to give C ✓				
		(iii)	0			1	ALLOW any unambiguous representation
							<b>ALLOW</b> C <sub>15</sub> H <sub>22</sub> O <sub>3</sub> N <sub>2</sub>
			HN	Ń. V			if both shown an incorrect formula <b>CON</b> s a correct structure or vice versa
	(b)		Monomer	Repeat unit	Type of polymerisation	2	ALLOW any unambiguous representation.
			CH <sub>2</sub> CHN(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub>	$ \begin{array}{ccc} H & N(C_2H_5)_2 \\C &C \\C \\ H & H \end{array} $	Addition		
			O H <sub>2</sub> N Cl	O HN	Condensation		
			Completely correct – 2 marks	-	<i>ı</i> or column		
			Total			6	

Q	uestic	on	Answer	Marks	Guidance
34	(a)		N S O N S O Diagram AND unpaired electron ✓	1	Incorrect structure scores 0
	(b)	(i)	A radical is used and produced (to continue the reaction) $\checkmark$	1	<b>ALLOW</b> there is a radical on both sides of the equation (AW)
		(ii)	$CO + 2O_2 \rightarrow O_3 + CO_2 \checkmark$	1	IGNORE hv Non reacting species shown on both sides are CON
	(c)		Frequency to break C—C <i>l</i> is 346000/ (6.02 x $10^{23}$ x 6.63 x $10^{-34}$ ) = 8.67 x $10^{14}$ Hz $\checkmark$	3	<b>ALLOW ECF</b> if kJ not turned into J or if Avogadro's constant is omitted.
			Frequency to break C—F is 467000/ (6.02 x $10^{23}$ x 6.63 x $10^{-34}$ ) =11.7 x $10^{14}$ Hz $\checkmark$ C—CI is broken, but UV absorbed is not of a harmful frequency <b>AND</b> C-F is broken and harmful UV absorbed. (AW) $\checkmark$ <b>OR</b> CFC-12 absorbs at both ends of the harmful range of radiation but not in the middle (AW) $\checkmark$		ALLOW a correct calculation of the bond energy needed to absorb 14.0 x $10^{14} =$ 559 kJmol <sup>-1</sup> and 10.1 x $10^{14} = 403$ kJmol <sup>-1</sup> for marks 1 or 2 ALLOW a correct calculation of energy (hv) of UV light and then comparison with energy per bond (J/N <sub>A</sub> ) for C-Cl and C-F for marks 1 and 2. E (10.1 x $10^{14}$ ) = 6.70 x $10^{-19}$ , E (14.0 x $10^{14}$ ) = 9.28 x $10^{-19}$ E (C-Cl) = 5.75 x $10^{-19}$ E (C-F) = 7.76 x $10^{-19}$
					<b>ALLOW</b> 1 mark for a correctly calculated frequency based on the sum of the bond enthalpies
					ALLOW correct comment based on incorrectly calculated frequencies
					<b>ALLOW</b> CFC-12 breaks down (AW) or both bonds break if incorrect calculation supports the statement.
			Total	6	

Ques	stion	Answer	Marks	Guidance	
35 (a)	)	Dissolve bolt in warm sulfuric acid ✓ Transfer to 1 dm <sup>3</sup> volumetric flask, (transfer washings) and make up to the mark (AW) ✓		Conc sulfuric acid is CON	
(b)	)	FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer = 45.6 or rounds to 46 (g) award 3 marks Moles of $MnO_4^-$ not needed by the rusty nail =(0.01792-0.00975) x 0.2 (= 1.634 x 10 <sup>-3</sup> ) $\checkmark$ Moles of Iron rusted in 10cm <sup>3</sup> solution =5 x 1.634 x 10 <sup>-3</sup> (=8.17 x10 <sup>-3</sup> ) $\checkmark$ In 1dm <sup>3</sup> mass = 0.817 x 55.8 = 45.6(g) $\checkmark$	3	ALLOW 2 or more sf ALLOW ECF between steps An answer rounding to 0.46 scores 2 (omission of the factor of 100 from 10 cm <sup>3</sup> to 1000cm <sup>3</sup> )	
(C)	) (i)	$O_2 + 2H_2O + 4e^- \rightarrow 4OH^- \checkmark$ Fe→ Fe <sup>2+</sup> + 2 e <sup>-</sup> <b>OR</b> Fe – 2e <sup>-</sup> → Fe <sup>2+</sup> ✓	2	ALLOW halved ALLOW reversible reactions shown either direction Extra half equations beyond 2 CONs 1 mark each	
	(ii)	Green solid is $Fe(OH)_2$ <b>AND</b> orange solid is $Fe_2O_3(.xH_2O)\checkmark$	1	ALLOW Fe(OH) <sub>3</sub> [Fe(OH) <sub>2</sub> (H <sub>2</sub> O) <sub>4</sub> ],[Fe(OH) <sub>3</sub> (H <sub>2</sub> O) <sub>3</sub> ]	
	(iii)	(Faster in salt water as) more (dissolved) ions (make it a better conductor) ✓ More OH <sup>-</sup> ions is <b>CON</b>	1	ALLOW '(water acts as a) 'salt bridge' and sea water contains a higher concentration of ions' ALLOW (the salt) acts as an electrolyte	
(d)	)	Fe <sup>2+</sup> 1s <sup>2</sup> 2s <sup>2</sup> 2p <sup>6</sup> 3s <sup>2</sup> 3p <sup>6</sup> 3d <sup>6</sup> ✓	1	IGNORE 4s <sup>o</sup> IGNORE working elsewhere. No of electrons in orbitals must be superscripts NOT [Ar]	
(e)	)	Ni(/Ni <sup>2+</sup> ) electrode potential is more negative than $H_2(H^+)$ <b>AND</b> thus $H^+$ can oxidise Ni to Ni <sup>2+</sup> (ORA) $\checkmark$ Cu(/Cu <sup>2+</sup> ) electrode potential is more positive than $H_2(H^+)$ so $H^+$ cannot oxidise Cu to Cu <sup>2+</sup> (ORA) $\checkmark$	2	ALLOW answers in terms of electron flow instead of oxidation We need a comment about each metal in relation to hydrog	
(f)	) (i)	Reaction is ligand substitution/exchange <b>AND</b> new ligand splits the d-orbitals differently ✓	1	ALLOW new complex ion has a different colour ALLOW nucleophilic substitution	
	(ii)	FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer = [Ni(EDTA)] <sup>2</sup> award 2 marks	2	DO NOT ALLOW charges inside brackets	

C	Question		Answer	Marks	Guidance
			Moles of Ni <sup>2+</sup> =0.025 x 0.25 = $6.25 \times 10^{-3}$ <b>AND</b> Moles EDTA = 0.0417 x 0.15 = $6.26 \times 10^{-3} \checkmark$ Ratio is 1:1 so formula is [Ni(EDTA)] <sup>2</sup> $\checkmark$		
			Total	15	

Question		on	Answer	Marks	Guidance	
36	(a)		Triple bond between N atoms requires a lot of energy to break (AW) /has a high bond enthalpy $\checkmark$	1	IGNORE very strong	
	(b)	(i)	$\Delta S = (3 \times 130.6) + 197.6 - (186.2 + 189.0)$	1	Sign must be included	
			Correct Expression evaluated with sign = $+214.2 \checkmark$			
		(ii)	Increase in entropy/positive as there are more molecules of products/gas ✓	1	<b>NOT</b> comments inconsistent with sign of $\Delta S$ calculated	
	(C)		FIRST CHECK THE ANSWER ON THE ANSWER LINE If answer = 962 (K) award 2 marks	2	ALLOW ECF from (b)(i)	
			T = 206000/214.2 ✓		ALLOW 963 (early rounding of 214.2) for 1 mark	
			Evaluated to 3sf =962 (K) ✓			
	(d)		$CO_2$ is used in <b>36.2</b> so it removes a greenhouse gas from the atmosphere, (this is greener) $\checkmark$	3		
			<ul> <li>Plus 2 from: ✓ ✓</li> <li>Both reactions need high T as both are endothermic but become more feasible at higher T as both have + ΔS, so no difference</li> <li>Both reactions give a higher yield at lower T</li> <li>Both need low pressure as 2 moles → 4, so no difference</li> <li>36.2 produces less hydrogen per mole of methane, so less green/ Atom economy is lower in 36.2. (ORA)</li> </ul>		ALLOW 36.2 requires more energy than 36.1, so less green Comments about <b>36.2</b> producing more toxic CO must be qualified (burn off $\rightarrow$ CO <sub>2</sub> or use as fuel) to score. Toxicity alone does not score.	

Question	Answer		Guidance
(e)	FIRST CHECK ANSWER ON ANSWER LINE If answer= 0.13 units dm <sup>6</sup> mol <sup>-2</sup> award 3 marks (0.1 moles of N <sub>2</sub> react so 0.3 moles of H <sub>2</sub> used and) 0.2 moles NH <sub>3</sub> form, 0.7 moles H <sub>2</sub> left $\checkmark$ ([NH <sub>3</sub> ] <sup>2</sup> /[N <sub>2</sub> ][H <sub>2</sub> ] <sup>3</sup> K <sub>c</sub> = 0.2 <sup>2</sup> /0.9 x 0.7 <sup>3</sup> ) evaluated = (0.13) $\checkmark$ units dm <sup>6</sup> mol <sup>-2</sup> $\checkmark$	3	<ul> <li>ALLOW ECF from incorrect concentrations but not from incorrect K<sub>c</sub> expression</li> <li>ALLOW 2 or more sf</li> <li>ALLOW mol<sup>-2</sup> dm<sup>6</sup></li> </ul>
(f) (i)	$4NH_3 + 5O_2 \rightarrow 4NO + 6H_2O \checkmark$	1	ALLOW multiples, halves
(ii)	FIRST CHECK ANSWER ON ANSWER LINE If answer = 12 (tonnes) award 4 marksMoles of NH4NO3 needed= $25 \times 10^{6}/80 \ (=3.125 \times 10^{5}) \checkmark$ Moles of NH3 needed to make nitric acid = $(100 \times 3.125 \times 10^{5})/77 \checkmark$ Total moles of ammonia = $3.125 \times 10^{5} + (100 \times 3.125 \times 10^{5})/77 = 7.18 \times 10^{5}$ Mass = $7.18 \times 10^{5} \times 17 = 1.22 \times 10^{7} \ g$ , 12 (tonnes) $\checkmark$	4	ALLOW 2 or more sf ALLOW ECF between stages MP1 convert to tonnes and then divide by 80 MP2 x100/77 MP3 Total moles ammonia (to make nitric acid + ammonia needed for salt) MP4 X 17 and evaluation and conversion to tonnes
(iii)	Add NaOH and Devarda's alloy or Al powder and warm ✓ Test gas with indicator paper/ red litmus/ rod dipped in HCl turns blue/ dense white fumes (due to ammonia) ✓	2	Reagents and heat needed Test and positive result for ammonia needed <b>ALLOW</b> Brown Ring Test (add $Fe_2SO_4$ solution followed by conc $H_2SO_4$ ) – a brown ring forms at the layer interface
	Total	18	

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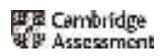
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