

# Mark Scheme (Results)

GCE Chemistry June 2017

Paper 9CH0/02 Advanced Organic and Physical Chemistry



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#### **General Marking Guidance**

- All candidates must receive the same treatment. Examiners must mark the first candidate in exactly the same way as they mark the last.
- Mark schemes should be applied positively. Candidates must be rewarded for what they have shown they can do rather than penalised for omissions.
- Examiners should mark according to the mark scheme not according to their perception of where the grade boundaries may lie.
- There is no ceiling on achievement. All marks on the mark scheme should be used appropriately.
- All the marks on the mark scheme are designed to be awarded. Examiners should always award full marks if deserved, i.e. if the answer matches the mark scheme. Examiners should also be prepared to award zero marks if the candidate's response is not worthy of credit according to the mark scheme.
- Where some judgement is required, mark schemes will provide the principles by which marks will be awarded and exemplification may be limited.
- When examiners are in doubt regarding the application of the mark scheme to a candidate's response, the team leader must be consulted.
- Crossed out work should be marked UNLESS the candidate has replaced it with an alternative response.
- Mark schemes will indicate within the table where, and which strands of QWC, are being assessed. The strands are as follows:

i) ensure that text is legible and that spelling, punctuation and grammar are accurate so that meaning is clear

ii) select and use a form and style of writing appropriate to purpose and to complex subject matter

iii) organise information clearly and coherently, using specialist vocabulary when appropriate

#### **Using the Mark Scheme**

Examiners should look for qualities to reward rather than faults to penalise. This does NOT mean giving credit for incorrect or inadequate answers, but it does mean allowing candidates to be rewarded for answers showing correct application of principles and knowledge. Examiners should therefore read carefully and consider every response: even if it is not what is expected it may be worthy of credit.

The mark scheme gives examiners:

- an idea of the types of response expected
- how individual marks are to be awarded
- the total mark for each question
- examples of responses that should NOT receive credit.

/ means that the responses are alternatives and either answer should receive full credit.

() means that a phrase/word is not essential for the award of the mark, but helps the examiner to get the sense of the expected answer.

Phrases/words in **bold** indicate that the <u>meaning</u> of the phrase or the actual word is **essential** to the answer.

ecf/TE/cq (error carried forward) means that a wrong answer given in an earlier part of a question is used correctly in answer to a later part of the same question.

Candidates must make their meaning clear to the examiner to gain the mark. Make sure that the answer makes sense. Do not give credit for correct words/phrases which are put together in a meaningless manner. Answers must be in the correct context.

### **Quality of Written Communication**

Questions which involve the writing of continuous prose will expect candidates to:

- write legibly, with accurate use of spelling, grammar and punctuation in order to make the meaning clear
- select and use a form and style of writing appropriate to purpose and to complex subject matter
- organise information clearly and coherently, using specialist vocabulary when appropriate.

Full marks will be awarded if the candidate has demonstrated the above abilities.

Questions where QWC is likely to be particularly important are indicated (QWC) in the mark scheme, but this does not preclude others.

Question Number	Answer	Mark
1(a)	The only correct answer is D (free radical substitution)	(1)
	<b>A</b> is not correct because electrophilic addition would be the reaction mechanism between alkenes, such as ethene, and chlorine but the substance in the question is the alkane, ethane	
	<b>B</b> is not correct because electrophilic substitution involves the reactions of aromatic systems but this question refers to the reaction of the alkane, ethane	
	<b><i>C</i></b> is not correct because the mechanism of reaction between an alkane such as ethane and chlorine involves substitution and not addition	

Question Number	Answer	Mark
1(b)	The only correct answer is B (4-ethyl-3-methylheptane)	(1)
	<b>A</b> is not correct because the longest consecutive carbon chain involves the seven carbon atoms from the top left of the molecular drawing down to the bottom right which means that the stem of the name is not pentane but heptane, with the consequential effect on the numbering and length of side chains	
	C is not correct because the longest consecutive carbon chain involves the seven carbon atoms from the top left of the molecular drawing down to the bottom right which means that the stem of the name is not hexane but heptane, with the consequential effect on the numbering and length of side chains	
	<b>D</b> is not correct because the longest consecutive carbon chain involves the seven carbon atoms from the top left of the molecular drawing down to the bottom right which means that the stem of the name is not hexane but heptane, with the consequential effect on the numbering and length of side chains	

Question Number	Answer	Additional Guidance	Mark
1(c)(i)	An explanation that makes reference to the following points:		(2)
	<ul> <li>different alkanes have different boiling temperatures/points (1)</li> </ul>	Allow Volatility for boiling temperature Allow Different alkanes condense at different temperatures Ignore melting temperatures if given with boiling temperatures Ignore densities	
	<ul> <li>because of (different) chain length/molar mass /strength of intermolecular forces/ number of electrons</li> <li>(1)</li> </ul>	Accept London /dispersion /van der Waals forces Allow reference to size A comparison such as 'longer alkanes have higher boiling points' scores 2 Ignore any reference to surface area Do not award references to cracking Do not award reference to just weight/mass Do not award incorrect trend	

Question Number	Answer	Additional Guidance	Mark
1(c)(ii)			(1)
	Correct equation	Example of equation:	
		$C_8H_{18} \rightarrow C_2H_4 + C_6H_{14}$ OR	
		$\begin{array}{rcl} C_8 H_{18} \ \to \ 2 C_2 H_4 \ + \ C_4 H_{10} \\ OR \end{array}$	
		$C_8H_{18} \rightarrow 3C_2H_4 + C_2H_6$	
		Allow	
		$CH_2=CH_2$ for $C_2H_4$	
		Products can be given in either order	
		Do not award equations forming $H_2$	

Question Number	Answer	Additional Guidance	Mark
1(c)(iii)	Correct equation	Example of equation:	(1)
		$\begin{array}{c cccccc} H & H & H & H & H & H \\ H & H & H & H &$	
		Accept bonds to hydrogen atoms inside the ring, e.g. $H \rightarrow C \rightarrow H \rightarrow C \rightarrow H$ $H \rightarrow C \rightarrow H \rightarrow C \rightarrow H$ $H \rightarrow C \rightarrow H \rightarrow C \rightarrow H$	
		Allow skeletal or structural formulae for hexane and for cyclohexane	
		or CH <sub>3</sub> (CH <sub>2</sub> ) <sub>4</sub> CH <sub>3</sub> or (CH <sub>2</sub> ) <sub>6</sub>	
		Ignore	
		$C_6H_{14} \rightarrow C_6H_{12} + H_2$	

(Total for Question 1 = 6 marks)

Question Number	Answer	Mark
2(a)	The only correct answer is C (hexagonal rings within a layer)	(1)
	<b>A</b> is not correct because in the layers of graphite and graphene the carbon atoms are bonded to three other carbon atoms and not four	
	<b>B</b> is not correct because graphite and graphene do not have pentagonal rings within their layers	
	<b>D</b> is not correct because graphene is a two-dimensional structure consisting of a single layer	

Question Number	Answer	Mark
2(b)	The only correct answer is D (all 120°)	(1)
	<b>A</b> is not correct because the angles within a layer of graphite and graphene are neither 90° nor 109.5° but are all 120°	
	<b>B</b> is not correct because the angles within a layer of graphite and graphene are not 109.5° but are all 120°	
	<b>C</b> is not correct because there are no angles within a layer of graphite and graphene that are 109.5° but they are all 120°	

Question Number	Answer	Mark
2(c)	The only correct answer is C (poor electrical conductivity)	(1)
	<b>A</b> is not correct because both graphene and graphite are similar to diamond in having a high melting temperature	
	<b>B</b> is not correct because neither graphene nor graphite nor diamond have a precise molecular formula since they are giant molecular structures	
	<b>C</b> is not correct because graphene, graphite and diamond are all giant molecular structures	

(Total for Question 2 = 3 marks)

Question Number	Answer	Additional Guidance	Mark
3(a)	An explanation that makes reference to the following points:	Accept reverse argument	(2)
		References to halogen reactivity scores (0)	
	• reactivity increases down Group (7) (1)	Do not award references to ions/halides	
	<ul> <li>because (C-X) bond enthalpy decreases / because (C-X) bond gets weaker down Group 7 (1)</li> </ul>	Do not award explanation in terms of just electronegativity or C-X dipoles	
		Ignore references to atom size, shielding etc and references to intermolecular forces	
		No TE on incorrect reactivity trend	

Question Number	Answer	Additional Guidance	Mark
3(b)(i)	<ul> <li>dipole on C—Br bond</li> <li>and</li> <li>curly arrow from C-Br bond to Br or just</li> <li>beyond</li> <li>(1)</li> </ul>	$\begin{array}{c} H \\ H_{3} \\ \hline \\ H \\ H \\ \hline \\ H \\$	(4)
	<ul> <li>curly arrow from <b>lone pair</b> on oxygen of hydroxide ion to carbon bonded to Br (1)</li> </ul>	Allow curly arrow to C <sup>+</sup> of carbocation	
	<ul> <li>formula of transition state with correct charge, partial bonding</li> <li>(1)</li> </ul>	Do not award if carbocation formed as intermediate Square brackets are not essential Allow charge on Br or OH of transition state Allow longer bonds for partial bonding Ignore geometry of transition state	
	• correct final products (1)	Allow NaBr product if mechanism starts with NaOH	
		Only penalise horizontal bond from the H of OH to C in the product e.g. $OH-CH_2CH_3$	
		Use of incorrect halogenoalkane loses this mark	
		One mark max deducted for omission of charge on ions, including transition state	
		$S_N1$ mechanism can score M1, M2 and M4 but not M3. M2 can be awarded for curly arrow from the lone pair on the oxygen of the hydroxide ion to the C+ of the carbocation intermediate	

Question Number	Answer	Additional Guidance	Mark	
3(b)(ii)	Reagents: nitric acid / HNO <sub>3</sub> and silver nitrate (solution) /AgNO <sub>3</sub> (1)	Use of hydrochloric acid/HCl OR sulfuric acid/H <sub>2</sub> SO <sub>4</sub> scores <b>(0)</b> Do not award acidified silver nitrate If name and formula given then both must be correct	(2)	
	(Result) cream/off-white precipitate (1)	Allow (very) <b>pale</b> yellow Do not award just white or just yellow Ignore subsequent additions of ammonia even if incorrect Result mark dependent on reagents mark or 'near miss' such as omitting to add nitric acid, using ethanolic silver nitrate, incorrect formulae		

Question Number		Answer		Additional Guidance	Mark
3(c)	(1)	(1)	(1)	Accept formulae in any order Award 2 if 3 correct displayed/structural formulae given Award 1 if 2 correct displayed/structural formulae given If more than 3 skeletal formulae drawn then deduct one mark for each additional formula 2-methylpropene negates a correct formula only if four formulae given View any formulae given with skeletal formula as working and ignore Ignore names even if incorrect Penalise any other alkenes such as pentenes, once only	(3)

Question Number	Answer	Additional Guidance	Mark
3(d)	An explanation that makes reference to the following points:		(3)
	<ul> <li>(only) ethanol has hydrogen bonding (and dipole-dipole and London forces)</li> <li>(1)</li> </ul>	Ignore references to ethanol having stronger London forces	
	<ul> <li>ethene (only) has (weaker) London/ instantaneous dipole –induced dipole forces (1)</li> </ul>	Accept dispersion /van der Waals forces	
	<ul> <li>more energy required to break the (stronger) intermolecular forces/hydrogen bonds in alcohols (1)</li> </ul>	A comparison is needed Allow overcome for break Allow 'heat' for energy Accept reverse argument	
		Do not award if the more energy required is given in response to just breaking stronger London forces for ethanol	
		Do not award M3 for covalent bonds breaking	

(Total for Question 3 = 14 marks)

Question Number	Answer	Additional Guidance	Mark
4(a)(i)	<ul> <li>conversion of pressure, volume and temperature to correct units (1)</li> </ul>	Example of calculation: $207kPa = 207\ 000\ Pa$ $8.98\ dm^3 = 0.00898\ m^3,$ $20^{\circ}C = 293\ K$	(3)
	<ul> <li>rearrangement of ideal gas equation so n=PV ÷ RT and calculation of n</li> <li>(1)</li> </ul>	n= <u>207 000 x 0.00898</u> = 8.31 x 293 = 0.7634	
	• conversion of answer into mass to 2/3 SF (1)	<ul> <li>= 0.7634 x 28 = 21.37647</li> <li>= 21.4 / 21 (g)</li> <li>Correct answer with no working scores 3</li> <li>TE on both parts of the calculation</li> </ul>	

Question Number		Answer	Additional Guidance	Mark
4(a)(ii)	•	The temperature increase will result in an increase in pressure because p is (directly) proportional to T (at constant volume and moles of gas)	Allow $p \propto T$ Reference to $p=nRT/V$	(1)

Question Number	Answer	Additional Guidance	Mark
4(b)	An explanation that makes reference to the following points:	Reference to molecule scores (0)	(2)
	• fewer protons (in nitrogen) (1)	Accept reverse arguments in terms of oxygen Allow weaker (effective) nuclear charge Allow smaller atomic number	
	<ul> <li>result in a weaker <b>nuclear</b> attraction because shielding is the same/electrons are in the same (sub)shell (in oxygen)/same number of electron shells</li> <li>(1)</li> </ul>	Do not award if incorrect numbers of protons stated or if ions referred to Do not award 'charge density' Ignore references to electron repulsion and electronegativity	

(Total for Question 4 = 6 marks)

Question Number	Answer	Additional Guidance	Mark
5(a)(i)	Correct equation	$\begin{array}{l} 2NO + 2CO \ \rightarrow \ N_2 + 2CO_2 \\ Accept \ multiples \end{array}$	(1)
		Ignore catalysts and conditions if stated	

Question Number	Answer	Additional Guidance	Mark
5(a)(ii)	A description that makes reference to the following points:	Absence of reference to the catalytic surface results in a deduction of one mark	(3)
	• <b>adsorption</b> of gases to catalytic surface (1)	Do not award absorption or "stick"	
	<ul> <li>weakening of bonds (and chemical reaction) on catalytic surface (1)</li> </ul>	Allow bonds break (and reaction occurs) on catalytic surface Ignore the type of interaction referred to between the reactants and the catalytic surface	
	• desorption of products from catalytic surface (1)	Allow 'release' of products from catalytic surface Allow de-adsorbed	

Question Number	Answer	Mark
5(b)	The only correct answer is A (area A)	(1)
	<b>B</b> is not correct because this is the area representing the number of particles with sufficient energy to react in the absence of a catalyst	
	<b>C</b> is not correct because this area subtraction does not represent the increase in the number of particles with sufficient energy to react	
	<b>D</b> is not correct because this sum of areas represents the total number of particles with sufficient energy to react in the presence of a catalyst	

Question Number	Answer	Mark
5(c)	The only correct answer is B (2.15 x 10 <sup>22</sup> )	(1)
	<b>A</b> is not correct because the molar mass of carbon dioxide has been used in the calculation instead of that of carbon monoxide	
	<b>C</b> is not correct because this is the number of molecules that are in one mole and not one gram of carbon monoxide	
	<b>D</b> is not correct because this is the result of incorrectly using the molar mass of carbon monoxide rather than the number of moles of carbon monoxide	

# (Total for Question 5 = 6 marks)

Question Number		Answer		Additional Guidance	Mark
6(a)	(CH <sub>2</sub> =CHCH <sub>2</sub> CH <sub>3</sub> $\rightarrow$ )	$ \begin{array}{c ccccccccccc} H & H & H & H \\ \hline                                  $	(1)	Accept skeletal, structural or displayed formulae or combination of which is clear, e.g. $-C_2H_5$ Brackets are not essential	(2)
		CH <sub>3</sub> CH <sub>3</sub>		Ignore 'n' Ignore orientation of side chains Ignore bond length Ignore where bond goes to for the ethyl groups	
	( →)	-[]-	(1)	Penalise lack of 'end-bonds' once only Award 1 mark max if only one repeat unit given for each polymer	
				Ignore more than 2 repeat units	

Question Number		Answer	Additional Guidance	Mark
6(b)	•	$HOOC - (CH_2)_4 - COOH$ or $CIOC - (CH_2)_4 - COCI$ (1)	Accept skeletal, structural or displayed formulae	(2)
	•	$H_2N - (CH_2)_4 - NH_2$ (1)	Penalise use of $C_4H_8$ once only Penalise missing H's once only	
			The monomers can be in either order Do not award monofunctionality	

Question Number	Answer	Additional Guidance	Mark
6(c)	$H_2N - (CH_2)_4 - COOH$	Accept skeletal, structural or displayed formulae	(1)
		Allow $H_2N - (CH_2)_4 - COCI$	
		Ignore connectivity	
		Allow	
		N H	
		Allow use of $C_4H_8$ here only	
		Penalise missing hydrogens	

Question Number	Answer	Additional Guidance	Mark
6(d)(i)	Allow H OOC $H_3C$ H H H H H H H H	Diagram must be 3-dimensional with either wedges or dashes to score 2 marks Ignore orientation of group at the top Ignore vertical bond to H of OH group	(2)

Question Number	Answer	Additional Guidance	Mark
6(d)(ii)	<ul> <li>They rotate the plane of plane-polarised light (equally) and in opposite/different directions OR Determine in which direction they rotate the plane of plane-polarised light</li> </ul>	Allow one plane	(1)

Question Number		Answer	Additional Guidance	Mark
6(d)(iii)	•	Does not accumulate in the environment/does not occupy landfill	Accept answers that outline the benefit of avoiding other means of disposal such as incineration, use of toxic chemicals Ignore just less harm to environment/less harm to animal life/less pollution/less of an "eyesore"/less energy to break it down	(1)

(Total for Question 6 = 9 marks)

Question Number	Answer		Additional Guidance	Mark
Number 7(a)			Example of synthetic pathway $ \bigcirc & \overset{(H_2S)_+}{R_{eflux}} & \bigcirc & & & & & \\ & & & & & & & & \\ & & & &$	(5)
	<ul> <li>A synthetic pathway that consists of:</li> <li>(reagents and conditions for the nitration of benzen conc. Nitric (HNO<sub>3</sub>) and sulfuric acids (H<sub>2</sub>SO<sub>4</sub>) and</li> </ul>	ie)	unbalanced, incorrect equations or reaction mechanisms Allow any single value or range between 50-60°C/warm/<55°C	
	55°C/heat/reflux	(1) (1)	Intermediate marks are standalone	
	<ul> <li>(reduction of nitrobenzene) tin and conc. hydrochloric acid and heat/reflux</li> </ul>	(1)	Allow iron & c.HCl Do not award dilute Ignore subsequent addition of NaOH Penalise lack of heat once only in M1 and M3	
	structure of phenylamine	(1)	Penalise just the names of intermediates once only	
	(reaction of phenylamine with) ethanoyl chloride	(1)	Ignore heat Do not award use of AICI <sub>3</sub>	

Question Number	Answer	Mark
7(b)	The only correct answer is B (6)	(1)
	<ul> <li>A is not correct because four carbon atoms in the aromatic ring are non-equivalent and not just three, so the correct total of non-equivalent carbon atoms and therefore peaks is six</li> <li>C is not correct because there are two sets of equivalent carbon atoms in the aromatic ring and not just one which means that the correct total of non-equivalent carbon atoms and therefore peaks is six</li> </ul>	
	<b>D</b> is not correct because this is the total number of carbon atoms in antifebrin but carbon atoms 2 and 6 in the aromatic ring are equivalent, as are 3 and 5, which gives a correct total of six non-equivalent carbon atoms and therefore six peaks	

Question Number	Answer	Additional Guidance	Mark
7(c)	An explanation that makes reference to the following points:		(2)
	<ul> <li>lone pair (of electrons) from the oxygen and will interact with the delocalised ring of electrons / increase the (pi/π) electron density of the benzene ring (1)</li> </ul>	Allow reference to the <b>lone pair</b> (of electrons) from the <b>nitrogen</b> Ignore activation of ring	
		Do not award charge density	
	<ul> <li>which increases the reactivity toward electrophiles (such as bromine)/ which means that the bromine is more easily polarised (1)</li> </ul>	Allow $Br^+/Br^{\delta+}$ for electrophile	
		Allow reference to benzene as being a stronger nucleophile	
		Do not award references to electrophilic addition	

Question Number	Answer	Additional Guidance	Mark
7(d)		Example of calculation:	(2)
	• conversion of moles to mass of paracetamol (1)	(mass of paracetamol = $3.10 \times 10^{-3} \times 151$ = 0.4681 (g)	
	• conversion of answer into percentage to 2/3 SF (1)	% = (0.4681÷ 0.500) x 100 = 93.62% ) =94 (%)/93.6 (%)	
		Allow TE for second mark from incorrect molar mass as long as value derived from dividing by 0.500/500mg and percentage is less than 100%	
		Correct answer without working scores 2	

(Total for Question 7 = 10 marks)

Question Number	Answer	Additional Guidance	Mark
8(a)(i)	$CH_3COCI + AICI_3 \rightarrow CH_3CO^+ + AICI_4^-$	Accept use of FeCl <sub>3</sub> /Fe +Cl <sub>2</sub> Allow displayed formulae Do not award C <sub>2</sub> H <sub>3</sub> OCl	(1)
		Ignore state symbols even if incorrect	

Question Number	Answer		Additional Guidance	Mark
8(a)(ii)			$CH_{3}$ $C$	(4)
	<ul> <li>electron pair movement from ring to electrophile</li> </ul>	(1)	$H^+ + AIG_+ \rightarrow AIG_3 + HG$ Do not award curly arrow that ends at the CH <sub>3</sub> Allow arrow starting anywhere within the hexagon	
	<ul> <li>formula of intermediate ion</li> </ul>	(1)	'Horseshoe' to cover at least three carbon atoms and face the tetrahedral carbon and with some part of the plus sign inside 'horseshoe' Allow Kekulé diagrams Do not award dotted bonds unless part of a 3D structure	
	<ul> <li>curly arrow from C–H bond to reform delocalised ring</li> </ul>	(1)		
	<ul> <li>correct product and equation to show regeneration of catalyst and HCI</li> </ul>	(1)	Could be shown in reaction mechanism Ignore curly arrows	

Question Number	Answer	Mark
8(b)(i)	The only correct answer is B (alkaline iodine solution)	(1)
	<b>A</b> is not correct because this oxidising agent would react with phenylethanal and not with phenylethanone which is the wrong way round	
	<b>C</b> is not correct because test is for aldehydes and so would react with phenylethanal and not with phenylethanone which is the wrong way round	
	<b>D</b> is not correct because test is for aldehydes and so would react with phenylethanal and not with phenylethanone which is the wrong way round	

Question Number	Answer	Additional Guidance	Mark
8(b)(ii)	An answer that makes reference to the following points:		(4)
	<ul> <li>formation of yellow/orange/red (crystalline) precipitate (1)</li> </ul>	Colour and state are both required Allow solid for ppt Ignore any conditions given with the use of 2,4-DNPH	
	• (Filter then) recrystallisation of products (1)		
	• determination of melting temperature (1)	Penalise M3 if any reference to boiling temperature	
	<ul> <li>comparison (and hence identification) from use of database/known values (1)</li> </ul>	Award only in the context of melting temperature of the hydrazones or as a TE of boiling temperature	
		Max 3 out of 4 if test is only carried out with one of the carbonyls	

Question Number	Ansv	ver	Additional Guidance	Mark	
*8(b)(iii)		red answer with linkages ve content and for how the vs lines of reasoning. v the marks should be	Guidance on how the mark scheme should be applied: The mark for indicative content should be added to the mark for lines of reasoning. For example, a response with four indicative marking points that is partially structured with some linkages and lines of reasoning scores 4 marks (3 marks for indicative content and 1 mark for partial structure and some linkages and lines of reasoning).		
	6 5-4 3-2 1 0	4 3 2 1 0	If there were no linkages between the points, then the same indicative marking points would yield and overall score of 3 marks (3 marks for indicative content and zero marks for linkages). In general it would be expected that 5 or 6 indicative points would get 2 reasoning marks, and 3 or 4 indicative points would get 1 mark for reasoning, and 0, 1 or 2 indicative points would score zero marks for reasoning. If there is any incorrect chemistry, deduct mark(s)		
	The following table shows how awarded for structure and line Answer shows a coherent logical structure with linkages and fully sustained lines of				
	reasoning demonstrated throughout Answer is partially structured with some linkages and lines of	1	from the reasoning. If no reasoning mark(s) awarded do not deduct mark(s).		
	reasoning Answer has no linkages between points and is unstructured	0	If there is no mention of protons/hydrogens in the response then deduct one structure and reasoning mark		

	ndicative content imilarities	Ignore references to C <sup>13</sup> nmr Accept annotations on a structure towards crediting the following IPs Allow either a single chemical shift value or a range within the stated values Penalise incorrect chemical shifts	
•	<b>IP1</b> : aromatic hydrogens will give similar/same peaks	Both have peaks in the range 6.5-8.4 (ppm) Ignore any splitting description	
•	<b>IP2</b> : both have a peak in the range $1.7-3.0$ (ppm) (due to the hydrogen of the H–C–C=O type)	Ignore any splitting pattern given for this peak to award this mark	
Di •	ifferences IP3 (Hydrogen environments): Phenylethanone has one less peak/hydrogen environment than phenylethanal	Allow any difference of one in the number of peaks stated	
•	<b>IP4</b> (Splitting patterns): a singlet for phenylethanone but a doublet and a triplet in phenylethanal	All these splitting patterns required for this IP	
•	<b>IP5</b> (Peak area ratios): relative peak (area) ratio in phenylethanone is 3 but in phenylethanal the peak (area) ratio is 2 to 1	Ignore the splitting pattern for this IP and ignore any peak areas given for the aryl hydrogens	
•	<b>IP6</b> (Chemical shifts): (Only) phenylethanal has an aldehyde (hydrogen) peak in the range 9 – 10.1 (ppm)	Ignore the splitting pattern for this IP	

Question Number	Answer	Additional Guidance	Mark
8(c)	• Reagent: lithium tetrahydridoaluminate((III)) /LiAlH <sub>4</sub> (1)	Allow lithium aluminium hydride	(2)
		Accept sodium tetrahydridoborate /sodium borohydride/ NaBH <sub>4</sub> In water/alcohol for 2 marks	
		Ignore reference to addition of acid after use of LiAlH $_4$ in dry ether	
		Do not award with additional reagents	
	• Conditions: (dry) ether/ethoxyethane (1)	Ignore heat	
		The mark for conditions is dependent on correct reagent or near miss such as incorrect formula LiAIH <sub>3</sub> /LiAIH/LiAI	

(Total Question 8 = 18 marks)

Question Number	Answer	Additional Guidance	Mark
9(a)	Correct answer to 2 SF	Example of calculation: (Four half-lives to decrease 600 g to 37.5 g so 4 x 14 mins) = 56 (mins)	(1)
		Penalise wrong units, e.g. "m"	

Question Number	Answe	r	Additional Guidance	Mark
9(b)(i)	Reaction Orders:			(3)
	<ul> <li>X First/1</li> <li>Y Second/2</li> </ul>	(1)		
	<ul> <li>Y Second/2</li> <li>Z Zero/0</li> </ul>	(1) (1)	Allow "none"/"no order"	

Question Number	Answer	Additional Guidance	Mark
9(b)(ii)	Marking point 1 • Rate = $k[X][Y]^2[Z]^0$ (1)	Reactants can be in any order Z does not have to be included in the rate equation TE from (b)(i) which will apply for all four marking points	(4)
	Marking point 2(1)• rearrangement of rate expression(1)Marking point 3(1)• calculation of value for k to 2/3 SF(1)	$k = rate / [X][Y]^2$	
	Marking point 4 • units dm <sup>6</sup> mol <sup>-2</sup> s <sup>-1</sup> (1)	Any 'run' can be used No TE on incorrect rearrangement Allow units in any order Correct answer without working and with correct units to 2/3 SF scores marking points 2, 3 and 4	

Question Number	Answer	
	The only correct answer is D (Fourth)	(1)
	<b>A</b> is not correct because this is the individual reaction order with respect to $bromate(V)$ ions and with respect to bromide ions but is not the overall reaction order	
9(c)(i)	<b>B</b> is not correct because this is the reaction order with respect to hydrogen ions but is not the overall reaction order	
	<b>C</b> is not correct because this is the number of species in the rate equation but is not the overall reaction order	

Question Number	Answer		Additional Guidance	Mark
9(c)(ii)			Example of calculation:	(2)
	<ul> <li>rearrange rate equation so [Br<sup>-</sup>] =</li> </ul>	(1)	$[Br^{-}] = \underline{rate} \\ k [BrO_{3}^{-}][H^{+}]^{2}$	
	<ul> <li>calculation of value to 2/3 SF</li> </ul>	(1)	= 0.255/0.26 (mol dm <sup>-3</sup> )	
			Correct answer without working to 2/3 SF scores 2 marks	
			If units given then must be correct	
			No TE on incorrect rearrangement	

Question Number			Answer Additional Guidance		ce	Mark	
9(d)	Example of suitable graph			1/7 1/0-3/16-1			(7)
<b>J</b> (u)		2.90 3:00	3.10	1/T ×10-3/K-1 3 20 3:30	3:40 3.	50	
		X					
		-510-					
		-6.0-					
		-710-	$\sim$				
				$\left<$			
		6.R		$\sim$			
				X			
		- 9:0-					
		-100			X		
		-110 - Gradient	= (-10.50-	-5/10) = -5/10 -2.95×103) 5/3×10	= (-)10,189	$\mathbf{X}$	
			(3.48×10 <sup>-3</sup> -	-2.95×153) 5'3×15	4		

(3.30, 3.19, 3.10, (3.00)) (5), -8.70, -7.55, -6.60, -3) pomission of end zero
3)
se more than 3SF once only
l points must cover at least ½ aph paper on each axis award 1/t
±½ square
$nt = -10200$ Allow $\pm 500$ this mark if the value is seen in calculation
0200 x 8.31 / 1000 +) 84.8 (kJ mol <sup>-1</sup> ) nswer must be positive and in nge (+) 80.6 – 88.9 (kJ mol <sup>-1</sup> ) value given in J mol <sup>-1</sup> but then units are essential SF for gradient and activation
ν ι

# **TOTAL FOR PAPER = 90 MARKS**

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