

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

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)
	A 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 is not correct because electrophilic substitution involves the reactions of aromatic systems but this question refers to the reaction of the alkane, ethane	
	C 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)	
	A 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		
	D 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)
	different alkanes have different boiling temperatures/points (1)	Allow Volatility for boiling temperature Allow Different alkanes condense at different temperatures Ignore melting temperatures if given with boiling temperatures Ignore densities	
	because of (different) chain length/molar mass /strength of intermolecular forces/ number of electrons (1)	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)	Correct equation	$\begin{array}{c} \underline{\text{Example of equation:}} \\ C_8H_{18} \rightarrow C_2H_4 + C_6H_{14} \\ \text{OR} \\ C_8H_{18} \rightarrow 2C_2H_4 + C_4H_{10} \\ \text{OR} \\ C_8H_{18} \rightarrow 3C_2H_4 + C_2H_6 \\ \\ \text{Allow} \\ \text{CH}_2 = \text{CH}_2 \text{ for } C_2H_4 \\ \\ \text{Products can be given in either order} \end{array}$	(1)
		Do not award equations forming H ₂	

Question Number	Answer	Additional Guidance	Mark
1(c)(iii)	Correct equation	Example of equation:	(1)
		$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	
		Accept bonds to hydrogen atoms inside the ring, e.g.	
		Allow skeletal or structural formulae for hexane and for cyclohexane	
		or CH ₃ (CH ₂) ₄ CH ₃ or (CH ₂) ₆	
		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)
	A 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 is not correct because graphite and graphene do not have pentagonal rings within their layers	
	D 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)
	A is not correct because the angles within a layer of graphite and graphene are neither 90° nor 109.5° but are all 120°	
	B is not correct because the angles within a layer of graphite and graphene are not 109.5° but are all 120°	
	${\it C}$ 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)
	A is not correct because both graphene and graphite are similar to diamond in having a high melting temperature	
	B is not correct because neither graphene nor graphite nor diamond have a precise molecular formula since they are giant molecular structures	
	C 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	
	 because (C-X) bond enthalpy decreases / because (C-X) bond gets weaker down Group 7 (1) 	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)	dipole on C—Br bond and curly arrow from C-Br bond to Br or just beyond (1)	$\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array} \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow \begin{array}{c} H_{3}C \longrightarrow \\ H_{3}C \longrightarrow \end{array}$ $\begin{array}{c} H_{3}C \longrightarrow $	(4)
	 curly arrow from lone pair on oxygen of hydroxide ion to carbon bonded to Br (1) 	Allow curly arrow to C ⁺ of carbocation	
	 formula of transition state with correct charge, partial bonding 	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 ₃ and silver nitrate (solution) /AgNO ₃ (1)	Use of hydrochloric acid/HCl OR sulfuric acid/H ₂ SO ₄ scores (0) 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) pale 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	
_	(1)	Answer (1)	(1)	Additional Guidance 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	(3)
				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	

Question Number	Answer	Additional Guidance	Mark
3(d)	An explanation that makes reference to the following points:		(3)
	 (only) ethanol has hydrogen bonding (and dipole-dipole and London forces) 	Ignore references to ethanol having stronger London forces	
	 ethene (only) has (weaker) London/ instantaneous dipole –induced dipole forces (1)	Accept dispersion /van der Waals forces	
	 more energy required to break the (stronger) intermolecular forces/hydrogen bonds in alcohols (1) 	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)	 conversion of pressure, volume and temperature to correct units (1) 	Example of calculation: 207kPa = 207 000 Pa 8.98 dm³ = 0.00898 m³, 20°C = 293 K	(3)
	 rearrangement of ideal gas equation so n=PV ÷ RT and calculation of n 	n= <u>207 000 x 0.00898</u> = 8.31 x 293 = 0.7634	
	• conversion of answer into mass to 2/3 SF (1)	= 0.7634 x 28 = 21.37647 = 21.4 / 21 (g) Correct answer with no working scores 3 TE on both parts of the calculation	

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 ∝ 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	
	result in a weaker nuclear attraction because shielding is the same/electrons are in the same (sub)shell (in oxygen)/same number of electron shells (1)	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	$2NO + 2CO \rightarrow N_2 + 2CO_2$ Accept multiples	(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)
	adsorption of gases to catalytic surface (1)	Do not award absorption or "stick"	
	weakening of bonds (and chemical reaction) on catalytic surface (1)	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 is not correct because this is the area representing the number of particles with sufficient energy to react in the absence of a catalyst	
	C is not correct because this area subtraction does not represent the increase in the number of particles with sufficient energy to react	
	D 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 (2.15×10^{22})	(1)
	A is not correct because the molar mass of carbon dioxide has been used in the calculation instead of that of carbon monoxide	
	C is not correct because this is the number of molecules that are in one mole and not one gram of carbon monoxide	
	D 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)	$(CH_2=CHCH_2CH_3\rightarrow)$		(1)	Accept skeletal, structural or displayed formulae or combination of which is clear, e.g. $-C_2H_5$	(2)
		Н Сн ₂ Н Сн ₂ 	. ,	Ignore 'n' Ignore orientation of side chains Ignore bond length	
				Ignore where bond goes to for the ethyl groups	
				Penalise lack of 'end-bonds' once only	
	(→)		(1)	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 − (CH₂)₄ − COOH 	Accept skeletal, structural or	(2)
	or	displayed formulae	
	CIOC - (CH2)4 - COCI (1		
		Penalise use of C ₄ H ₈ once only	
	• $H_2N - (CH_2)_4 - NH_2$ (1	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 ₄ H ₈ here only	
		Penalise missing hydrogens	

Question Number	Answer	Additional Guidance	Mark
6(d)(i)	Allow H00C H3C H (1) H00C COOH CH3 (1)	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)	 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 	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
7(a)			Example of synthetic pathway	(5)
			C. H250+ C-HNO3 Reflux NO2	
			Sn/cHCl No ₂ Heat	
			CH3(OCI) NH2 CH3(OCI) CH3(OCI) CH3 CH3 CH3 CH3	
			The compounds used can be stated or given within equations. Ignore any unbalanced, incorrect equations or reaction mechanisms	
	A synthetic pathway that consists of:		reaction mechanisms	
	 (reagents and conditions for the nitration of benzer conc. Nitric (HNO₃) and sulfuric acids (H₂SO₄) and 55°C/heat/reflux 	ne) (1)	Allow any single value or range between 50-60°C/warm/<55°C	
	structure of nitrobenzene	(1)	Intermediate marks are standalone	
	 (reduction of nitrobenzene) tin and conc. hydrochloric acid and heat/reflux 	(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 AlCl ₃	

Question Number	Answer	Mark
7(b)	The only correct answer is B (6)	(1)
	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	
	$m{C}$ 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	
	D 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)
	• lone pair (of electrons) from the oxygen and will interact with the delocalised ring of electrons / increase the (pi/п) electron density of the benzene ring (1)	Allow reference to the lone pair (of electrons) from the nitrogen Ignore activation of ring	
		Do not award charge density	
	 which increases the reactivity toward electrophiles (such as bromine)/ which means that the bromine is more easily polarised (1) 	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 ₃ /Fe +Cl ₂	(1)
		Allow displayed formulae Do not award C₂H₃OCl	
		Ignore state symbols even if incorrect	

Question Number	Answer		Additional Guidance	Mark
8(a)(ii)			$\begin{array}{cccccccccccccccccccccccccccccccccccc$	(4)
	 electron pair movement from ring to electrophile 	(1)	$H^+ + AICI_+ \longrightarrow AICI_3 + HCI$ Do not award curly arrow that ends at the CH ₃ Allow arrow starting anywhere within the hexagon	
	formula of intermediate ion	(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	
	 curly arrow from C-H bond to reform delocalised ring 	(1)		
	 correct product and equation to show regeneration of catalyst and HCl 	(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)
	A is not correct because this oxidising agent would react with phenylethanal and not with phenylethanone which is the wrong way round	
	$m{\mathcal{C}}$ is not correct because test is for aldehydes and so would react with phenylethanal and not with phenylethanone which is the wrong way round	
	D 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: • formation of yellow/orange/red (crystalline) precipitate (1)	Colour and state are both required Allow solid for ppt Ignore any conditions given with the use of 2,4-DNPH	(4)
	 (Filter then) recrystallisation of products determination of melting temperature (1) 	Penalise M3 if any reference to boiling temperature	
	 comparison (and hence identification) from use of database/known values (1) 	Award only in the context of melting	
		Max 3 out of 4 if test is only carried out with one of the carbonyls	

Question Number	Ansv	wer	Additional Guidance	Mark
*8(b)(iii)	This question assesses the stucture coherent and logically structure and fully sustained reasoning. Marks are awarded for indicate answer is structured and show the following table shows how	red answer with linkages ive content and for how the vs lines of reasoning.	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	(6)
	marking points seen in fanswer	Number of marks awarded for indicative marking points	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).	
	The following table shows how awarded for structure and line Answer shows a coherent logical structure with 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	
	and fully sustained lines of reasoning demonstrated throughout Answer is partially structured with some linkages and lines of	1	incorrect chemistry, deduct mark(s) 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	

Indicative content

Similarities

• **IP1**: aromatic hydrogens will give similar/same peaks

• **IP2**: both have a peak in the range 1.7-3.0 (ppm) (due to the hydrogen of the H–C–C=O type)

Differences

- **IP3** (Hydrogen environments): Phenylethanone has one less peak/hydrogen environment than phenylethanal
- **IP4** (Splitting patterns): a singlet for phenylethanone but a doublet and a triplet in phenylethanal
- **IP5** (Peak area ratios): relative peak (area) ratio in phenylethanone is 3 but in phenylethanal the peak (area) ratio is 2 to 1
- **IP6** (Chemical shifts): (Only) phenylethanal has an aldehyde (hydrogen) peak in the range 9 10.1 (ppm)

Ignore references to C¹³ 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

Both have peaks in the range 6.5-8.4 (ppm)
Ignore any splitting description

Ignore any splitting pattern given for this peak to award this mark

Allow any difference of one in the number of peaks stated

All these splitting patterns required for this IP

Ignore the splitting pattern for this IP and ignore any peak areas given for the aryl hydrogens

Ignore the splitting pattern for this IP

Question Number	Answer	Additional Guidance	Mark
9(a)	Reagent: lithium tetrahydridoaluminate((III)) /LiAlH ₄ (1)	Allow lithium aluminium hydride	(2)
8(c)		Accept sodium tetrahydridoborate /sodium borohydride/ NaBH ₄ 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 LiAlH ₃ /LiAlH/LiAl	

(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	Answer		Additional Guidance	Mark
	Reaction Orders:			(3)
9(b)(i)				
	X First/1	(1)		
	• Y Second/2	(1)		
	• Z Zero/0	(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 • rearrangement of rate expression (1) Marking point 3 • calculation of value for k to 2/3 SF (1)	Example of calculation: $k = \text{rate} / [X][Y]^2$ $k = \frac{2.17 \times 10^{-6}}{0.00100 \times 0.00300^2}$ = 241.11 = 241/240		
	Marking point 4 • units dm ⁶ mol ⁻² s ⁻¹ (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	Mark
	The only correct answer is D (Fourth)	(1)
	A 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 is not correct because this is the reaction order with respect to hydrogen ions but is not the overall reaction order	
	C 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)
3(c)(ii)	 rearrange rate equation so [Br⁻] = 	(1)	$[Br^{-}] = \frac{\text{rate}}{\text{k } [BrO_3^{-}][H^{+}]^2}$	
	 calculation of value to 2/3 SF 	(1)	= 0.255/0.26 (mol dm ⁻³)	
			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	Answei					Additio	nal Guid	dance	Mark
9(d)	Example of suitable graph	2.40	3.00	3:10	1/T ×10	-3/K-1	3:40	3.50	(7)
		-1:0							
		-510-							
		-610	X						
		-710-		X					
		ln k -8:0							
		-910-							
		-100					X		
		110 -	Gradient=	(-10.50.	-5/10) =.	5140 513×10	=010,189 1		

		<u>'</u>	
• calculation of all three 1/T values x 10 ⁻³	(1)	(3.41), 3.30, 3.19, 3.10, (3.00)	
calculation of all three In k values	(1)	(-9.75), -8.70, -7.55, -6.60, (-5.58) Allow omission of end zero Penalise more than 3SF once only	
axes: correct way round, labelled, suitable scale	e (1)	Plotted points must cover at least ½ the graph paper on each axis Do not award 1/t	
all points plotted correctly, with best-fit straight	t line(1)	Allow ±½ square	
calculation of gradient with sign	(1)	Gradient = -10200 Allow ± 500 Allow this mark if the value is seen in the E_a calculation	
units of gradient	(1)	K	
use of gradient to calculate activation energy	(1)	E _a = 10200 x 8.31 / 1000 = (+) 84.8 (kJ mol ⁻¹) Final answer must be positive and in the range (+) 80.6 – 88.9 (kJ mol ⁻¹) Allow value given in J mol ⁻¹ but then these units are essential Ignore SF for gradient and activation energy values	
		(Total for Question 9 - 18 mark	

(Total for Question 9 = 18 marks)

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