

Mark Scheme (Results) Summer 2007

GCE

GCE Chemistry (6244) Paper 01

Edexcel Limited. Registered in England and Wales No. 4496750 Registered Office: One90 High Holborn, London WC1V 7BH



General Guidance on Marking

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.

Crossed out work should be marked UNLESS the candidate has replaced it with an alternative response.

Using the mark scheme

The mark scheme gives you:

- 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.

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.

- 1 / means that the responses are alternatives and either answer should receive full credit.
- 2 () 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.
- 3 [] words inside square brackets are instructions or guidance for examiners.
- 4 Phrases/words in bold indicate that the <u>meaning</u> of the phrase or the actual word is essential to the answer.
- 5 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.

6244/01

	EXPECTED ANSWER		EXPECTED ANSWER	ACCEPT	REJECT	MARK
1.	(a)	(i)	$Mg^{2+}(g)$ (+) $O^{2-}(g)$		if state symbols missing If 2e⁻ included in box	(1)
		(ii)	$\Delta H_1 (\text{Enthalpy of}) \text{ formation (of MgO)} \qquad (1)$ $\Delta H_2 (\text{Enthalpy of}) \text{ atomisation (of Mg)} \qquad (1)$ $\Delta H_3 1^{\text{st}} \text{ plus } 2^{\text{nd}} \text{ electron affinity (of O)}$ $\text{OR } 1^{\text{st}} \text{ and } 2^{\text{nd}} \text{ electron affinity (of O)} \qquad (1)$	Recognisable abbreviation such as "EA" for electron affinity.		(3)
		(iii)	$\Delta H_{f} = -602 = (+150) + (+2186) + (+249) + (+657) + LE$ OR (LE =) -(+657)-(+249)-(+2186)-(+150)+(-602)(1) (LE =) -3844 (kJ mol ⁻¹) (1) Correct answer only with no working (1 max)		Doubling electron affinity and/or atomisation values scores (0) any incorrect sign in algebraic expression (0)	(2)

		EXPECTED ANSWER	ACCEPT	REJECT	MARK
(b)	(i)	The electrons around the iodide ion are drawn towards the magnesium ion	(Mg ²⁺) polarises (I ⁻ ion) "distortion" if clearly linked to the iodide ion "Mg ion" "I ion" OR "iodine ion"	Any reference to atoms or molecules e.g. "Mg polarises" "iodine/I/I ₂ is polarised" Wrong polarisation e.g. "magnesium ion is polarised" "I ⁻ polarises Mg ²⁺ "	(1)
	(ii)	Radius/size (of ions) (1) charge (on ions) (1)	Distance between ions OR Sum of (ionic) radii OR Type of crystal structure OR Madelung constant "Charge density"	"atomic radius"	(2)
	(iii)	Less (exothermic) (1) covalent character (strengthens lattice) (1)	Smaller OR more endothermic OR Less negative OR Lower Theoretical value based on purely/100 % ionic model	Higher/greater Any implication of magnesium iodide being totally covalent	(2)
		Mark each aspect independently			

		EXPECTED ANSWER	ACCEPT	REJECT	MARK
(c)	(i)	Enthalpy change when 1 mol of gaseous ions (1)	Energy or heat	Any implication of an endothermic process e.g. energy required	
		is dissolved such that further dilution causes no further heat change (1)		"1 mol of gaseous atoms"	
		IGNORE "standard conditions" Mark each aspect independently	"Added to water"/"reacts with water" instead of "dissolved"	Just "hydrated" Just "completely hydrated"	
			Is dissolved to form an infinitely dilute solution OR Is dissolved in a large/excess/infinite amount of water		(2)

	EXPECTED ANSWER	ACCEPT	REJECT	MARK
(ii)	EITHER $\Delta H_{SOLN} = (-\Delta H_{LE} + \Delta H_{HYD})$ (1) Expression quoted or correctly used in at least one of the calculations below			
	$\Delta H_{SOLN} MgSO_4 = -(-2874) + (-1920)$ = +954(kJ mol ⁻¹) (1)	Answer only with no working (1)		
	$\Delta H_{SOLN} BaSO_4 = -(-2374) + (-1360)$ = +1014(kJ mol ⁻¹) (1)	Answer only with no working (1)		
	Enthalpy of solution of MgSO ₄ less endothermic/more exothermic/more negative than for BaSO ₄ , so MgSO ₄ more soluble than BaSO ₄ (or reverse argument) (1)		Just "solubility/Δ <i>H</i> _{soln} depends on a balance between lattice and hydration energies"	
	(both) lattice energies and hydration enthalpies decrease from $MgSO_4$ to $BaSO_4$ (or down group) (1)			(4)
	(but) lattice energies change less (1)	"The hydration energies decrease faster"	(-)500 and (-)560 stated without further explanation	
	$\Delta H_{SOLN} = (-\Delta H_{LE} + \Delta H_{HYD})$ (1) stated in words or symbols			
	so ΔH_{soln} less exothermic/more endothermic/more positive for BaSO ₄ so less soluble OR so ΔH_{soln} more exothermic/more negative/less endothermic for MgSO ₄ so MgSO ₄ more soluble (1)		Just "solubility/Δ <i>H</i> _{solution} depends on a balance between lattice and hydration energies"	
			(Tot	al 17 marks)

		EXPECTED ANSWER	ACCEPT	REJECT	MARK
2.	(a)	V: H H H H H H H H H H H H H H H H H H H		Any compressed formulae e.g. C₂H₅Br OR CH₃CH₂Br	
				Any compressed formulae e.g. CH ₃ CONH ₂ OR CH ₃ CNH ₂	(2)

	EXPECTED ANSWER	ACCEPT	REJECT	MARK
(b)	Accept names or formulae, but ignore correct or incorrect conditions Step A: NH ₃ (1) Ignore state or dilution or solvent for ammonia			
	Step B: $K_2Cr_2O_7$ (1) and H_2SO_4 (1)	"Hydrochloric acid"/"HCl" instead of "H2SO4"	Incorrect oxidation number for dichromate(VI)	
	N.B. only award the acid/H ₂ SO ₄ mark if a correct (or a near-miss) oxidising agent given	$Cr_2O_7^{2-}$ and H^+ (2) OR CrO_4^{2-} and H^+ (2) OR Acidified dichromate ions (2) OR Acidified $K_2Cr_2O_7$ (2) OR acidified dichromate((VI)) (1) OR KMnO ₄ (1) and H_2SO_4 (1) OR alkaline KMnO ₄ (1) then acidify (accept any acid) (1)	CrO₄ ²⁻ alone (0) Hydrochloric acid/HCl with KMnO₄ (0) Incorrect oxidation number for manganate(VII)	(6)
	Step C: PCl_5 OR $SOCl_2$ OR PCl_3 (1)			
	Step D: P_2O_5 OR P_4O_{10} (1)			
	Step E: LiAlH₄ (1)	C₂H₅OH and Na OR H₂ and Ni /Pt/Pd (catalyst) OR Lithal	NaBH₄ (O)	

	EXPECTED ANSWER		ACCEPT	REJECT	MARK
(c)	(i)	(substituted) amide OR (N-substituted) amide OR secondary amide OR 2° amide		Polyamide amine	(1)
	(ii)	$\begin{array}{c c} & H & 0 & 0 \\ & & & \\$		– CONH	(2)
(d)	Loss	of smell:	HCl as a reactant and		
	CH ₃	$CH_2NH_2 + H^+ \rightarrow CH_3CH_2NH_3^+$ (1)	$CH_3CH_2NH_3^+Cl^-$ as product		
			(charges are not required).		
	Retu <i>CH</i> ₃ OR CH ₃ C	urn of smell: $CH_2NH_3^+ + OH^- \rightarrow CH_3CH_2NH_2 + H_2O$ $CH_2NH_3Cl + NaOH \rightarrow CH_3CH_2NH_2 + H_2O + NaCl (1)$	$CH_{3}CH_{2}NH_{3}^{+} + NaOH \rightarrow$ $CH_{3}CH_{2}NH_{2} + H_{2}O + Na^{+} (1)$		(2)
					(Total 13 marks)

		EXPECTED ANSWER	ACCEPT	REJECT	MARK
3.	(a)	PH 14 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	ACCEPT	pH range 3 to 5	(4)
		 e.g. pH range 3 to 6 OR 3 to 7 OR 3 to 8 OR 4 to 7 OR 4 to 8 OR 5 to 8 (1) (do not need to start/finish on whole numbers) Middle of vertical pH range between 4 and 6 (1) 		ph range 5 to 5	

	EXPECTED ANSWER	ACCEPT	REJECT	MARK
(b)	Bromocresol green Indicator(s) CQ on graph [check table on question paper]	More than one indicator for extended vertical regions		(1)
(c)	pH change around equivalence point too small OR pH changes over too big a volume (1)	Too small a vertical (region) OR no vertical (region) OR no point of inflexion OR no sudden change in pH OR no straight section		
	for a sharp colour change of indicator (1)	No sharp/clear/precise end point OR very small range over which indicator changes colour	No suitable indicator OR No "easy" colour change	(2)
	[If say ammonia is a strong base or ethanoic acid is a strong acid, or both, (0 out of 2)]			
			(То	tal 7 marks)

			EXPECTED ANSWER	ACCEPT	REJECT	MARK
4.	(a)	Н				(1)
	(b)	este	r			(1)
	(C)	(i)	Moles: C ₂ H ₅ OH: 3.75 (1) Moles: HCOOC ₂ H ₅ : 2.50 and moles H ₂ O : 2.50 (1) for both			(2)
		(ii)	$K_{c} = \frac{[HCOOC_{2}H_{5}][H_{2}O]}{[HCOOH][C_{2}H_{5}OH]}$		Obviously round brackets "()"	(1)
		(iii)	$K_{c} = \frac{2.50}{0.485} \times \frac{2.50}{0.485}$ Must have clearly divided moles of each component by 0.485 for 1 st mark e.g. [HCOOC ₂ H ₅] = [H ₂ O] = 5.16 (mol dm ⁻³) and [HCOOH] = 1.03 (mol dm ⁻³) and [C ₂ H ₅ OH] = 7.73 (mol dm ⁻³) = 3.33 (1) stand alone mark IGNORE sig.figs.	$K_{c} = \frac{(2.50)^{2}}{0.50 \times 3.75} = 3.33 \text{ only scores}$ (2) if it is stated that V cancels either here or in (iv) If $[H_{2}O]$ omitted in (ii), then answer $K_{c} = 0.647 \text{ mol}^{-1}\text{dm}^{3}$ (2) but this will give $K_{c} = 1.33 \text{ mol}^{-1}\text{dm}^{3}$ with V omitted from calculation (1)	1 st mark if 485 used as V in expression	(2)
		(iv)	No, (as) equal numbers of moles on both sides OR volumes cancel OR mol dm ⁻³ cancel OR units cancel OR crossing out units to show they cancel	 "equal powers/moles on both sides" OR " powers cancel" Mark CQ on K_c expression in (ii) 	"concentrations cancel"	(1)

		EXPECTED ANSWER	ACCEPT	REJECT	MARK
(d)	(i)	(as reaction) endothermic (1) K _e decreases (1)	Exothermic in backward direction (or words to that effect)		
			direction, 1 mark only (out of 4) for CQ "increase in K _c "		(4)
		numerator in quotient (has to) decrease OR denominator in quotient (has to) increase OR fraction (has to) decrease (1)			
		yield of $HCOOC_2H_5$ decreases (1)			
	(ii)	no effect as catalysts do not affect (the value of) K OR no effect as catalysts do not affect the position of equilibrium OR no effect as catalysts do not affect the yield OR No effect as catalysts increase the rate of the forward and backward reactions equally/to the same extent OR no effect as catalysts only increase the rate OR no effect as catalysts only alter the rate "no effect" can be stated or implied		Just "catalysts increase rate"	(1)
		IGNORE any references to activation energy			
				Т) (Т	otal 13 marks)

			EXPECTED ANSWER	ACCEPT	REJECT	MARK
5.	IGNO	ORE st	ate symbols throughout this question.			
	(a)	(i)	$Na_2O + H_2O \rightarrow 2NaOH$	$ \rightarrow 2Na^+OH^-$ OR $ \rightarrow 2Na^+ + 2OH^-$ Multiples e.g. $2Na_2O + 2H_2O \rightarrow 4NaOH$		(1)
		(ii)	ionic	Giant ionic OR electrovalent		(1)
	(b)	(i)	$P_4O_{10} + 6H_2O \rightarrow 4H_3PO_4$ OR $P_2O_5 + 3H_2O \rightarrow 2H_3PO_4$	Multiples		(1)
		(ii)	covalent	'molecular covalent' 'simple covalent'	'co <u>n</u> valent' OR 'giant covalent' OR dative covalent	(1)
	(C)	basi IGN(IGN(c (oxides) to acidic (oxides) (1) both words needed DRE references to Al_2O_3 DRE references to amphoteric character of Al_2O_3		The elements change from basic to acidic	(2)
		met.	allic character (of the elements) decreases (1) DRE "across group" if used instead of "across period"	metal to non-metal		

EXPECTED ANSWER		EXPECTED ANSWER	ACCEPT	REJECT	MARK
(d)	(i)	$CO_2 + H_2O \rightleftharpoons H^+ + HCO_3^-$ OR $CO_2 + 2H_2O \rightleftharpoons H_3O + + HCO_3^-$	$CO_{2} + H_{2}O = 2H^{+} + CO_{3}^{2}$ OR $CO_{2} + H_{2}O = H_{2}CO_{3} = H^{+} + HCO_{3}^{-}$ OR $CO_{2} + H_{2}O = H_{2}CO_{3} = 2H^{+} + CO_{3}^{2}$ $"\rightarrow" \text{ instead of "="}$	$\begin{array}{l} JUST\\ CO_2 + H_2O \ \rightleftharpoons \ H_2CO_3 \end{array}$	(1)
	(ii)	$PbO + 2H^{+} \rightarrow Pb^{2+} + H_{2}O$ OR $PbO + 2HNO_{3} \rightarrow Pb(NO_{3})_{2} + H_{2}O$ (1) $PbO + 2OH^{-} + H_{2}O \rightarrow Pb(OH)_{4}^{2^{-}}$ OR $PbO + 2NaOH + H_{2}O \rightarrow Na_{2}Pb(OH)_{4}$ OR $PbO + 2NaOH \rightarrow Na_{2}PbO_{2} + H_{2}O$ (1)	formation of $Pb(OH)_6^{4^-}$ OR $Na_4Pb(OH)_6$	PbO with other acids Any equations with PbO ₂	(2)

EXPECTED ANSWER		EXPECTED ANSWER	ACCEPT	REJECT	MARK
(e)	Incre	eases as		Just "increases" on its own (0) OR "increases" followed by incorrect justification (0)	
	<u>EITH</u> oxid ampl [amp	I <u>ER</u> es acidic at the top of the group (1) hoteric at bottom of group (1) photeric must be stated for the 2 nd mark]	Oxides more basic down group 1 (out of 2) OR oxides less acidic down group 1 (out of 2)		
	<u>OR</u> atom	ns become larger (1)	atoms have more shielding/shells/energy levels OR "(outer) electrons further from nucleus" [no need to refer to atoms in this case]	Just "more electrons" OR "the elements become larger"	(2)
	(so) OR h OR m [allo made	more easily lose electrons have lower ionisation energies hore easily form positive ions (in compounds) (1) w 2 nd mark even if no specific reference has been e to atoms]			
	[N.B. EITH	. Marking cannot allow points taken from both the ER and OR arguments together]			
				(Tota	i i i marks)

	EXPECTED ANSWER		EXPECTED ANSWER	ACCEPT	REJECT	MARK
6.	(a)	(i)	$CH_3CH_2CH_2OH$ (1)	$C_2H_5CH_2OH$ OR full structural formulae.g. H H H H	C ₃ H ₇ OH	
				$ \begin{array}{cccccccccccccccccccccccccccccccccccc$		(2)
			Reduction OR nucleophilic addition (1) IGNORE heterolytic	Redox "Nucleophilic reduction"		



EXPECTED ANSWER		EXPECTED ANSWER	ACCEPT	REJECT	MARK
	(iii)	$CH_{3}CH_{2}CH(OH)C \equiv N$ (1) nucleophilic addition (1) both words needed IGNORE "heterolytic"	$C_{2}H_{5}CH(OH)CN$ OR $CH_{3}CH_{2}CH(OH)CN$ OR full structural formula e.g. $H H OH$ $H-C-C-C-C-CN$ $H H H$ OR $H H O$ $H - C-C-C-C=N$ $H H H$		(2)
(b)	(i)	$Mg + C_2H_5Br \rightarrow C_2H_5MgBr$ IGNORE charges		C ₂ H ₅ BrMg C ₂ H ₅ MgI	(1)
	(ii)	Dry ethoxyethane OR dry ether IGNORE references to I ₂ OR heat			(1)

	EXPECTED ANSWER		ACCEPT	REJECT	MARK
(c	z) (i)	With propanone: Structure: (CH ₃) ₂ C(OH)CH ₂ CH ₃ (1) Name: 2-methylbutan-2-ol OR 2-hydroxy-2-methylbutane (1) IGNORE punctuation	C ₂ H ₅ in lieu of CH ₂ CH ₃ OR the full structural formula e.g. $\begin{array}{c} H_{3}C & H & H \\ H_{3}C & C & C & C \\ H_{3}C & C$	С ₅ H ₁₁ OH	(2)

EXPECTED ANSWER		ACCEPT	REJECT	MARK
	With butanal: Structure: $CH_3CH_2CH_2CH(OH)CH_2CH_3$ (1) Name: hexan-3-ol (1)	the full structural formula e.g. H H H H H H H-C-C-C-C-C-C-H H H H H H H OR H H H H H H H H-C-C-C-C-C-C-H H H H H H H H-C-C-C-C-C-C-C-H H H H O H H H CQ structure provided it is a secondary alcohol hexane-3-ol CQ name provided that it is a secondary alcohol	С ₆ H ₁₃ OH	(2)
(ii)	H (group in butanal) is planar (1) attacked (with equal probability) from two directions (1) Mark each aspect independently unless reference made to carbocations etc		Butanal/the molecule/it is planar OR Butanal/the molecule/it is linear References to carbocations OR carbonium ions OR Planar intermediates OR S _N 1 mechanisms scores (0 out of 2)	(2)