

CAMBRIDGE INTERNATIONAL EXAMINATIONS

Cambridge International Advanced Level

MARK SCHEME for the October/November 2014 series

9701 CHEMISTRY

9701/41

Paper 4 (A2 Structured Questions), maximum raw mark 100

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Mark schemes should be read in conjunction with the question paper and the Principal Examiner Report for Teachers.

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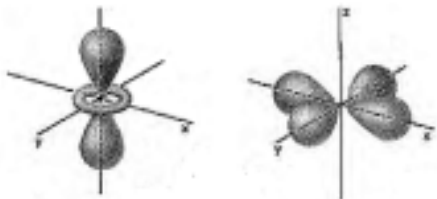
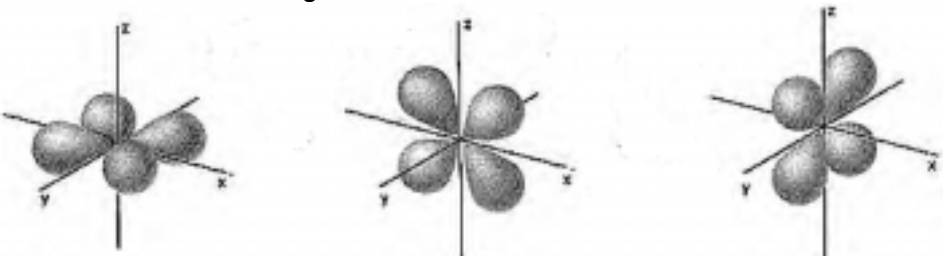
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Question	Marking point	Marks	Marks total
1 (a) (i)	[NO] 2 nd order and the concentration is $\times 2$, rate $\times 4$	1	
	[O ₂] 1 st order and evidence of using expt 1 & 2 when the concentration is $\times 2$, rate doubles	1	
(ii)	(0.00408 $\times$ 27) rate = 0.11 (mol dm ⁻³ s ⁻¹) to 2sf	1	
(iii)	(Rate =) $k [\text{O}_2][\text{NO}]^2$	1	
(iv)	$k = 332(.03125)$ mol ⁻² dm ⁶ s ⁻¹	1 1	[6]
(b) (i)	labelled axes x-axis: energy (KE) and y-axis: molecules or particles two curves: starts origin; not touching x-axis again; no levelling out; curves only intersecting once curves labelled and T2 is to the right and lower max than T1	1 1 1	
	(ii) rate increases and energy of the particles increases more particles have E_a	1 1	
(c)	1 mole of F ₂ and 1 mole NO reacting in the slow step a balanced mechanism consistent with overall equation e.g. $\text{F}_2 + \text{NO} \rightarrow \text{NOF} + \text{F}$ OR $\text{F}_2 + \text{NO} \rightarrow \text{NOF}_2$ $\text{NO} + \text{F} \rightarrow \text{NOF}$ $\text{NO} + \text{NOF}_2 \rightarrow 2\text{NOF}$	1 1	[2]
Total			[13]

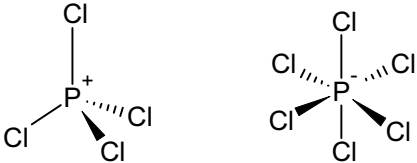
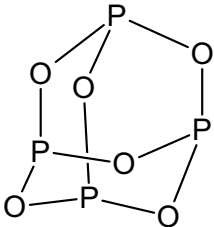
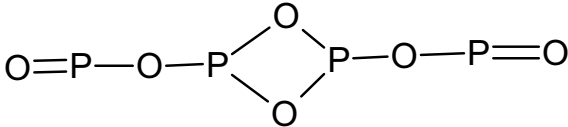
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2 (a)	3d <table><tr><td>(Ni)</td><td>↑↓</td><td>↑↓</td><td>↑↓</td><td>↑</td><td>↑</td></tr><tr><td>(Ni²⁺)</td><td>↑↓</td><td>↑↓</td><td>↑↓</td><td>↑</td><td>↑</td></tr></table> 4s <table><tr><td>↑↓</td></tr><tr><td></td></tr></table>	(Ni)	↑↓	↑↓	↑↓	↑	↑	(Ni ²⁺)	↑↓	↑↓	↑↓	↑	↑	↑↓		1 1	[2]
(Ni)	↑↓	↑↓	↑↓	↑	↑												
(Ni ²⁺)	↑↓	↑↓	↑↓	↑	↑												
↑↓																	
(b) (i)	degenerate	1															
(ii)	2 upper orbitals and 3 lower orbitals	1															
(iii)	correct upper orbital diagram  correct lower orbital diagram 	1 1	[4]														
(c)	electron(s) move from lower to upper level absorb (red/blue) light/photon complementary colour (green) is seen OR green light is transmitted	1 1 1	[3]														

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(d)	A $\text{Ni}(\text{OH})_2$ OR $\text{Ni}(\text{OH})_2(\text{H}_2\text{O})_4$	1	
	B $[\text{Ni}(\text{NH}_3)_6]^{2+}$ OR $[\text{Ni}(\text{NH}_3)_n(\text{H}_2\text{O})_{6-n}]^{2+}$ OR $[\text{Ni}(\text{NH}_3)_n(\text{H}_2\text{O})_{4-n}]^{2+}$	1	
	$\text{Ni}^{2+} + 2\text{OH}^- \rightarrow \text{Ni}(\text{OH})_2$ OR $[\text{Ni}(\text{H}_2\text{O})_6]^{2+} + 2\text{OH}^- \rightarrow \text{Ni}(\text{OH})_2 + 6\text{H}_2\text{O}$ OR $[\text{Ni}(\text{H}_2\text{O})_6]^{2+} + 2\text{NH}_3 \rightarrow \text{Ni}(\text{OH})_2 + 4\text{H}_2\text{O} + 2\text{NH}_4^+$ OR $[\text{Ni}(\text{H}_2\text{O})_6]^{2+} + 2\text{OH}^- \rightarrow \text{Ni}(\text{OH})_2(\text{H}_2\text{O})_4 + 2\text{H}_2\text{O}$	1	
	$\text{Ni}(\text{OH})_2 + 6\text{NH}_3 \rightarrow [\text{Ni}(\text{NH}_3)_6]^{2+} + 2\text{OH}^-$ OR $[\text{Ni}(\text{H}_2\text{O})_6]^{2+} + 6\text{NH}_3 \rightarrow [\text{Ni}(\text{NH}_3)_6]^{2+} + 6\text{H}_2\text{O}$	1	[4]
Total			[13]

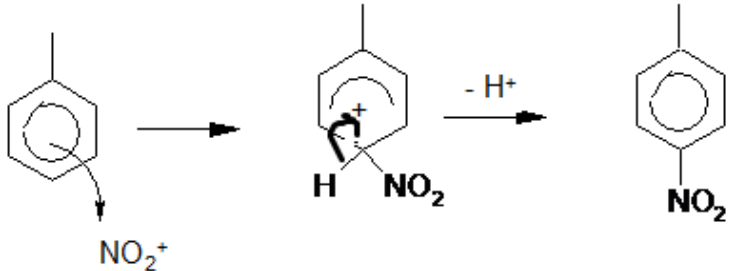
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3	(a) (i)	$101 = \text{P}^{35}\text{Cl}^{35}\text{Cl}$ $103 = \text{P}^{35}\text{Cl}^{37}\text{Cl}$ $105 = \text{P}^{37}\text{Cl}^{37}\text{Cl}$	1 1 1	
	(ii)	9:6:1	1	[4]
	(b) (i)	PCl_5 5 bonding pairs around P	1	
	(ii)		1 1	[3]
	(c) (i)	 P_4O_6 structure where each P has three P-O bonds and each O has two P-O bonds e.g. 	1	
	(ii)	(molecule/ion/species) that donates a lone pair of electrons (to a central transition metal atom or ion)	1	[2]
(d)	(i)	$K_{\text{sp}} = [\text{Ca}^{2+}]^3[\text{PO}_4^{3-}]^2$	1	

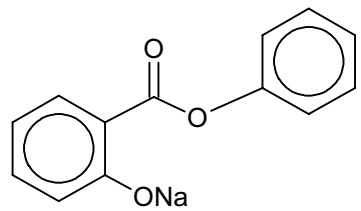
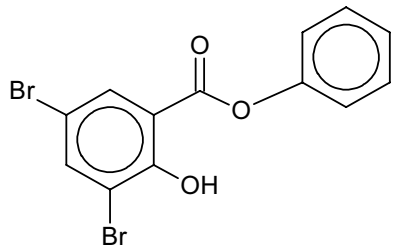
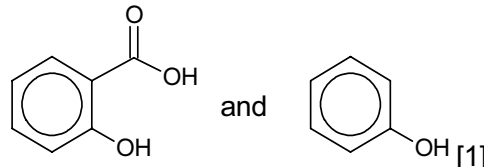
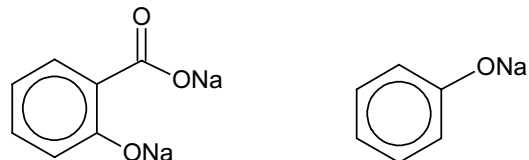
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(ii)	$[Ca^{2+}] = 3 \times 2.50 \times 10^{-6} = 7.50 \times 10^{-6} \text{ mol dm}^{-3}$ $[PO_4^{3-}] = 2 \times 2.50 \times 10^{-6} = 5.00 \times 10^{-6} \text{ mol dm}^{-3}$ $= (7.50 \times 10^{-6})^3 (5.00 \times 10^{-6})^2$ $= 1.05(1.1) \times 10^{-26}$ $\text{mol}^5 \text{dm}^{-15}$	1 1 1	[4]
(e) (i)	(enthalpy change) when 1 mole of an ionic compound is formed from its gaseous ions	1 1	
(ii)	Mg ²⁺ has a smaller (ionic) radii than Ca ²⁺ OR Mg ²⁺ is smaller than Ca ²⁺	1	[3]
Total			[16]
4 (a) (i)	$2H_2SO_4 + HNO_3 \rightarrow 2HSO_4^- + NO_2^+ + H_3O^+$ OR $H_2SO_4 + HNO_3 \rightarrow HSO_4^- + NO_2^+ + H_2O$	1	

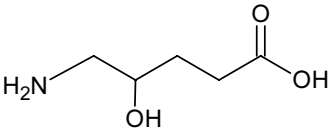
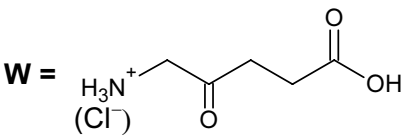
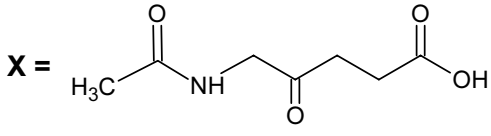
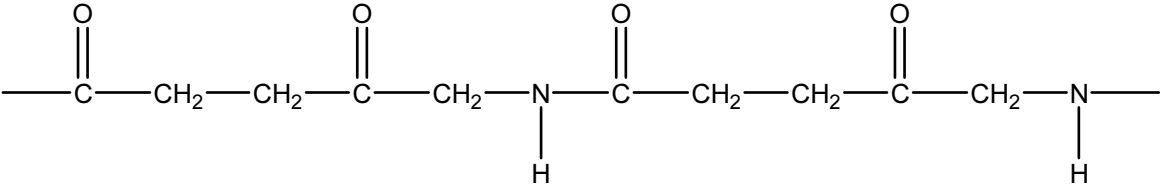
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(ii)	any three of • curly arrow from inside the benzene ring to NO_2^+ group • intermediate – penalise NO_2 connectivity or missing methyl group (once) • curly arrow from C-H bond into ring • product + H^+ (or as diagram $-\text{H}^+$) allow 2- and 3-substituted nitromethylbenzene) 	3	[4]
(b) (i)	acidity of $\text{ClCH}_2\text{CO}_2\text{H} > \text{CH}_3\text{CO}_2\text{H}$ AND ($\text{ClCH}_2\text{CO}_2\text{H}$) as an electronegative/electron withdrawing Cl	1	[3]
(ii)	acidity of phenol $> \text{CH}_3\text{CH}_2\text{OH}$ AND electrons on oxygen (on phenol) delocalised into ring	1	
	OR benzene ring withdraws electrons from oxygen stronger acid linked to weakening O-H bond/anion being stabilised	1	

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(c)	Na	 (or ionic)	redox/reduction				
	Br ₂		(electrophilic) substitution				
	NaOH	 and  or ionic	hydrolysis/ acid-base/				
1 mark for each correct structure for reaction types, 2 correct = 1 mark, 3 correct = 2 marks						4 2	[6]

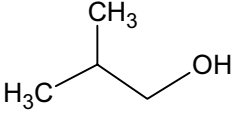
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(d)	Y =  W =  X =  each structure 1 mark	3	[3]
(e)	 correct displayed amide formula correct polyamide with two repeat units	1 1	[2]
Total			15
6 (a)	• (move in different directions) some amino acids have a different charge • (move at different speeds) some amino acids have a different size/different charge • (some amino acids do not move at all) some amino acids exist as a zwitterions/have no net(overall) charge/neutral/both NH₂/COOH are charged in amino acids	1 1 1	[3]
(b) (i)	mobile – solvent or water stationary – alumina/silica (supported on glass/plastic/Al)	1 1	
(ii)	by adsorption	1	[3]

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(c)	any three of: (all can be awarded from a clear, labelled diagram) • (base pairing) A to T OR C to G • H-bonds between bases • two/double stranded/chains • anti-parallel strands • (general structure) sugar-phosphate backbone OR BASE-SUGAR-PHOSPHATE bonded in a diagram	3	[3]
(d)	van der Waals' forces lost (in val) H-bonding gained (in ser)	1 1	[2]
Total			11
7 (a)	amide group circled OR indicated as diagram ester group circled OR indicated as diagram	1 1	[2]
(b)	lower doses of the drug required OR improved activity of the drug OR reduced side effects	1	[1]

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(c)	decreases enzyme activity OR decreases rate at which product is formed	1	
	binds with the enzyme's active site OR has a complementary shape to active site OR similar shape to substrate	1	
	(competitive inhibition can be overcome by) increasing [substrate] OR increasing substrate concentration	1	[3]
(d)	energy source/carrier OR releases energy when hydrolysed	1	[1]
Total			7
8 (a)	M:M+1 = 100/(1.1 x n)	1	
	20.4/0.9 = 100/(1.1 x n) x = 4	1	
(ii)	C ₄ H ₁₀ O	1	[3]
(b) (i)	2-methylpropan-1-ol OR correct structure 	1	
(ii)	0.9-1.0 is (2 x)CH ₃ R/CH ₃ /RCH multiplet/1.8 is CHR/R ₃ CH singlet/2.5 is OH 3.4 is CH ₂ O/CH ₃ O	1 1 1 1	
(iii)	doublet 1H/one proton on adjacent carbon	1 1	

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(iv)	OH peak or one peak disappears	1	[9]
	OH proton is labile or exchanges for D of D ₂ O or as an equation e.g. $\text{D}_2\text{O} + \text{OH} \rightarrow \text{DOH} + \text{OD}$ as a minimum	1	
Total			12
			100