Cambridge International Advanced Level

MARK SCHEME for the October/November 2014 series

9701 CHEMISTRY

9701/42

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

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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 ×2, rate doubles	1	
(ii)	(ii) (0.00408×27) rate = <u>0.11</u> (mol dm ⁻³ s ⁻¹) to 2sf		
(iii)	(Rate =) $k [O_2][NO]^2$	1	
(iv)	(iv) $k = 332(.03125)$ mol ⁻² dm ⁶ s ⁻¹		
(b) (i)	(i) labelled axes <i>x</i> -axis: energy (KE) and <i>y</i> -axis: molecules or particles two curves: starts origin; not touching <i>x</i> -axis again; no levelling out; curves only intersecting once curves labelled and T2 is to the right and lower max than T1		
(ii)	rate increases and energy of the particles increases	1	
	more particles have <i>E</i> _a	1	[5]
(c)	1 mole of F_2 and 1 mole NO reacting in the slow step	1	
	a balanced mechanism consistent with overall equation		
	e.g. $F_2 + NO \rightarrow NOF + F$ OR $F_2 + NO \rightarrow NOF_2$ NO + F \rightarrow NOF NOF NOF ₂ \rightarrow 2NOF		[2]
Total			[13]

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2 (a)	3d4s	1	
	(Ni) $\uparrow \downarrow$ $\uparrow \downarrow$ $\uparrow \downarrow$ \uparrow \uparrow	1	[2]
	(Ni ²⁺) $\land \downarrow$ $\land \downarrow$ $\land \downarrow$ \land		
(b) (i)	degenerate	1	
(ii)	2 upper orbitals and 3 lower orbitals	1	
(iii)	correct upper orbital diagram	1	[4]
(c)	electron(s) move from lower to upper level	1	
	absorb (red/blue) light/photon	1	
	complementary colour (green) is seen OR green light is transmitted	1	[3]

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(d)	Ni ²⁺ + 2 OR [Ni(H OR [Ni(H	$\begin{array}{l} H_{2} \textbf{OR} Ni(OH)_{2}(H_{2}O)_{4} \\ H_{3})_{6}]^{2^{+}} \textbf{OR} [Ni(NH_{3})_{n}(H_{2}O)_{6^{-n}}]^{2^{+}} \textbf{OR} [Ni(NH_{3})_{n}(H_{2}O)_{4^{-n}}]^{2^{+}} \\ POH^{-} \rightarrow Ni(OH)_{2} \\ POH^{-} \rightarrow Ni(OH)_{2} \\ POH^{-} \rightarrow Ni(OH)_{2} + 6H_{2}O \\ H_{2}O)_{6}]^{2^{+}} + 2OH^{-} \rightarrow Ni(OH)_{2} + 4H_{2}O + 2NH_{4}^{+} \\ POH^{-} \rightarrow Ni(OH)_{2} + 2OH^{-} \rightarrow Ni(OH)_{2} + 2H_{2}O \end{array}$			1 1 1	
	Ni(OH) ₂ - OR Ni(H ₂	+ $6NH_3 \rightarrow [Ni(NH_3)_6]^{2+-} + 2OH^-$ $_2O)_6]^{2+} + 6NH_3 \rightarrow [Ni(NH_3)_6]^{2+-} + 6H_2O$			1	[4]
Total						[13]

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3 (a) (i)	$101 = P^{35}Cl^{35}Cl$ $103 = P^{35}Cl^{37}Cl$ $105 = P^{37}Cl^{37}Cl$	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)	$P_{4}O_{6} \text{ structure where each P has three P-O bonds and each O has two P-O bonds e.g.}$ $O = P - O - P - O - P = O$	1	
(ii)	(molecule/ion/species) that donates a lone pair of electrons (to a central transition metal atom or ion)	1	[2]
(d) (i)	$K_{\rm sp} = [Ca^{2+}]^3 [PO_4^{3-}]^2$	1	

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

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(ii)	 int cu pr 	e of irly arrow from inside the benzene ring to NO_2^+ group termediate – penalise NO_2 connectivity or missing methyl group (ond irly arrow from C-H bond into ring oduct + H ⁺ (or as diagram –H ⁺) and 3-substituted nitromethylbenzene) $\downarrow -H^+$ $H^- NO_2$ NO_2^+	e)		3	[4]
(b) (i)	Cl	$C_{1}CH_{2}CO_{2}H > CH_{3}CO_{2}H$ AND ($C_{1}CH_{2}CO_{2}H$) as an electronegative/e		-	1	
(ii)	OR benze	phenol > CH ₃ CH ₂ OH AND electrons on oxygen (on phenol) delocal ene ring withdraws electrons from oxygen	sed into ring		1	
	stronger a	acid linked to weakening O-H bond/anion being stabilised			1	[3]

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

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Total				13
5	(a)	$CH_{3}CH_{2}COCl > CH_{3}CH_{2}CH_{2}Cl > C_{6}H_{5}Cl$		
		 any two of: C-Cl bond strength is weakest in CH₃CH₂COCl ora In C₆H₅Cl (no hydrolysis) C-Cl bond is part of delocalised system OR p-orbital on Cl overlaps with π system OR electrons from Cl overlap with π system CH₃CH₂COCl carbon in C-Cl bond is more electron deficient since it is also attached to an oxygen atom ora 	1+1	[3]
	(b) ketone, amine, carboxylic acid two correct 1 mark, all three 2		2	[2]
	(C) (i)	dipole on C-Br curly arrow breaking C-Br bond curly arrow from lone pair on N to carbon in C-Br bond H_{2N} H_{2N} H_{3C} H_{3C} H_{3C} H_{2} H_{3C} H_{2} H_{3C} H_{2} H_{3C} H_{2} H_{3C} H_{2} H_{3C} H_{2} H_{2} H_{3C} H_{2} H_{2} H_{2} H_{3C} H_{2} H_{2} H_{3C} H_{2} H_{2} H_{2} H_{3C} H_{2} H_{2} H_{3C} H_{2} H_{2} H_{3C} H_{2}	1 1 1	
	(ii)	nucleophilic substitution	1	
	(iii) HBr or hydrogen bromide		1	[5]

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(d)		3	[3]
	$Y = H_2N OH OH$		
	$W = \bigcup_{\substack{H_3N^+ \\ (CI^-) \\ O}} OH \qquad X = \bigcup_{\substack{H_3C^- \\ H_3C^- \\ OH}} OH$		
	each structure 1 mark		
(e)	$ \begin{array}{ c c c c } & \bigcirc & $		
	correct displayed amide formula correct polyamide with two repeat units	1 1	[2]
Total			15
6 (a)	(move in different directions)	1	
	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	[3]
(b) (i)	mobile – solvent or water stationary – alumina/silica (supported on glass/plastic/Al)		
(ii)	by adsorption		

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	(c)	any three	e of: (all can be awarded from a clear, labelled diagram)				
		 H- tw an (go 	ase pairing) A to T OR C to G bonds between bases o/double stranded/chains iti-parallel strands eneral structure) sugar-phosphate backbone OR BASE-SUGAR-PHC a diagram	OSPHATE bo	onded	3	[3]
	(d)	van der Waals' forces lost (in val) H-bonding gained (in ser)					[2]
Total							11
7	(a)	amide group circled OR indicated as diagram ester group circled OR indicated as diagram $H_{3}C \longrightarrow CH_{3}$ $H_{3}C \longrightarrow CH_{3}$		1 1	[2]		
	(b)	OR impro	es of the drug required ved activity of the drug ed side effects			1	[1]

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				1			
	(c) decreases enzyme activity OR decreases rate at which product is formed						
		binds with the enzyme's active site OR has a complementary shape to active site OR similar shape to substrate					
		(competitive inhibition can be overcome by) increasing [substrate] OR increasing substrate concentration					
	(d)	energy source/carrier OR releases energy when hydrolysed					
То	tal			7			
8	(a)	M:M+1 = 100/(1.1 x n) 20.4/0.9 = 100/(1.1 x n)	1				
		x = 4	1				
	(ii)	C ₄ H ₁₀ O	1	[3]			
	(b) (i)	2-methylpropan-1-ol OR correct structure H_3C OH	1				
	(ii)	$\begin{array}{llllllllllllllllllllllllllllllllllll$	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	
		n is labile or exchanges for D of D ₂ O equation e.g. D ₂ O + OH \rightarrow DOH + OD as a minimum			1	[9]
Total						12
						100