## MARK SCHEME for the October/November 2011 question paper

## for the guidance of teachers

# 9701 CHEMISTRY

9701/42

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

This mark scheme is published as an aid to teachers and candidates, to indicate the requirements of the examination. It shows the basis on which Examiners were instructed to award marks. It does not indicate the details of the discussions that took place at an Examiners' meeting before marking began, which would have considered the acceptability of alternative answers.

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1	(a) (i) (ii)	red/	-	our of bi	n mixture of H <sub>2</sub> + C <i>l</i> <sub>2</sub> <i>but</i> No romine decolourises/disapp		[1
			ainer gets warm/h				[2
	(iii)	H-H	= 436		C <i>l</i> -C <i>l</i> = 244	H-C $l$ = 431	
		ΔH	= 436 + 244 - 2(4	431)	= -182 kJ mol <sup>-1</sup>		[2
		H-H	= 436		Br-Br = 193	H-Br = 366	
		ΔH	= 436 + 193 – 2(3	366)	$= -103 \text{ kJ mol}^{-1}$		[2
	(iv)	H-B	r bond is weaker tl	han the	H-C <i>l</i> bond – allow convers	se.	[1 <b>[8</b>
	(b) (i)	light					[1
	(ii)		ds made = C-I &		= 410 + 151 = 561 = 240 + 299 = 539 = 551 - 539 = +22 kJ m	ol <sup>-1</sup>	[2
	(iii)				othermic <i>or</i> no strong bo		_
	(c) (i)		olytic fission is th electron species	e breał	king of a bond to form (two	o) radicals/neutral	species/ [1

- [1] (ii) •CH<sub>2</sub>Cl [1] **[3]** the C-Br bond is the weakest or needs least energy to break/breaks most easily
- (d) CI

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- 4 structures: [2] 2 or 3 structures: [1]
  - [1] **[3]**

[Total: 18]

Correct chiral atom identified

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2	Ord		er w.r.t. [CH₃CHO] = 1 er w.r.t. [CH₃OH] = 1 er w.r.t. [H <sup>+</sup> ] = 1		[1] [1] [1]
	(ii)	rate	= k[CH₃CHO][CH₃OH][H <sup>+</sup> ]		[1]
	(iii)	units	$s = mol^{-2} dm^6 s^{-1}$		[1]
	(iv)	rate	will be $2 \times 4 = 8$ times as fast as reaction 1 (relative ratio	ate = 8)	[1] <b>[6]</b>

(b)

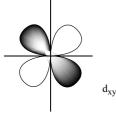
	[CH <sub>3</sub> CHO] /mol dm <sup>-3</sup>	[CH <sub>3</sub> OH] /mol dm <sup>-3</sup>	[H <sup>+</sup> ] /mol dm <sup>-3</sup>	[acetal <b>A</b> ] /mol dm <sup>-3</sup>	[H <sub>2</sub> O] /mol dm <sup>-3</sup>
at start	0.20	0.10	0.05	0.00	0.00
at equilibrium	(0.20 – x)	(0.10 – 2x)	0.05	x	x
at equilibrium	0.175	0.05	0.05	0.025	0.025

(iv)	$K_c = 0.025^2/(0.175 \times 0.05^2) = 1.4(3) \text{ (mol}^{-1} \text{ dm}^3)$	[1] [max 9]
(iii)	$K_c = \{[acetal A][H_2O]\}/\{[CH_3CHO][CH_3OH]^2\}$ units = mol <sup>-1</sup> dm <sup>3</sup>	[1] [1]
(ii)	4 values in third row	4 x [1]
(i)	3 values in second row	3 x [1]

[Total: 15]

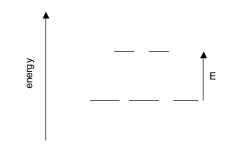
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3 (a) for example.... also allow d<sub>z2</sub>



shape (4 lobes) [1] correct label e.g. d<sub>xy</sub> [1] [2]

(b) (i)

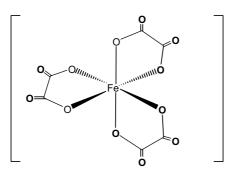


Marks are for 5 degenerate orbitals [1] and 3:2 split [1]

(ii) colour due to the absorption of light NOT emitted light[1] $E = hf or photon's energy = E in above diagram[1]electron promoted from lower to higher orbital[1]size of <math>\Delta E$  depends on the ligand[1]as  $\Delta E$  changes, so does f in E = hf[1][7]

(c) (i) 
$$O.N.(carbon) = +3$$
 (4 × (-2) + 2x = -2, thus 2x = +6) [1]

(iii)



[2]

(iv)  $\underline{2} K_3 Fe(C_2O_4)_3 \rightarrow \underline{3} K_2C_2O_4 + \underline{2} FeC_2O_4 + \underline{2} CO_2$  $Or K_3 Fe(C_2O_4)_3 \rightarrow \underline{3/2} K_2C_2O_4 + FeC_2O_4 + CO_2$ [2]

[max 5]

[Total: 14]

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4	(a) (i)	Allow ⇒ instead of arrow				[1]	
	(ii)					_	
		r	nost basic		least basic		
		е	thylamine	ammonia	phenylamine		
							[1]
	(iii)		ethylamine > NH <sub>3</sub> due to electron-donating ethyl/alkyl group phenylamine < NH <sub>3</sub> due to delocalisation of lone pair over ring				[1] [1] <b>[4]</b>
	(b) (i)	C <sub>6</sub> H	$_{5}\text{OH} + \text{OH}^{-} \rightarrow$	$C_6H_5O^- + H_2O$ (or	with Na <sup>+</sup> /H <sub>2</sub> O/A <sup>-</sup> )		[1]
	(ii)	pKa of nitrophenol is smaller/K <sub>a</sub> is larger because it's a stronger acid/dissociates more than phenol stronger because the anionic charge is spread out moreover the NO <sub>2</sub> group <i>or</i> NO <sub>2</sub> is electron-withdrawing					[1]
	(iii)	pKa	oKa = 1.0				
	(iv)	Nitro	group increas	es acidity / electro	n-withdrawing groups	increase acidity	( [1] [ <b>5]</b>

### (c) (i) **B** is phenyldiazonium cation, $C_6H_5-N^+\equiv N$

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l	l	•	)	F

/			
	reaction	reagent(s)	conditions
	Step 1	NaNO <sub>2</sub> + HC <i>1</i> or HNO <sub>2</sub> [1]	T < 10°C [1]
	Step 2	H₂O / aq	heat/boil/T > 10° (both) [1]
	Step 3	HNO₃ NB HNO₃(aq) OK for both	dilute (both) [1]
			٢ <i>٨</i>

[4] **[5]** 

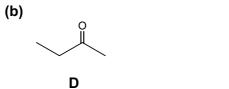
[1]

[Total: 14]

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- 5 (a) (i) C=C double bonds / alkenes
  - (ii) -OH groups / accept alcohols or acids
  - (iii)  $CH_3CO-$  or  $CH_3CH(OH)-$  groups
  - (iv) carbonyl, >C=O, groups / accept aldehydes and ketones 4 × [1]



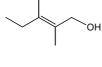






(c) isomers of C





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trans

correct structure (excl. stereochemistry)	[1]
cis and trans drawn correctly	[1]
type of isomerism is cis-trans or geometrical isomerism	[1]
	[3]

[Total: 9]

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6	(a) (i)	2H <sub>2</sub> 1	$NCH_2CO_2H \rightarrow H_2NCH_2CONHCH_2CO_2H + H_2O$		[1]
	(ii)	Skel	etal formula required		[1] <b>[2]</b>
	(b) (i)	α he β ple	lix eated sheet		[1] [1]
	(ii)	For Nee with	dents should choose one of the structures belowα helix:For β pleated sheetd to show a helixNeed to show twoC=O H-Nstrands with C=Oveen turnsthem	parallel 'zig-zag'	
		Whie	chever is chosen, overall structure [1] position of H bo	onds [1]	

[4]

(c)

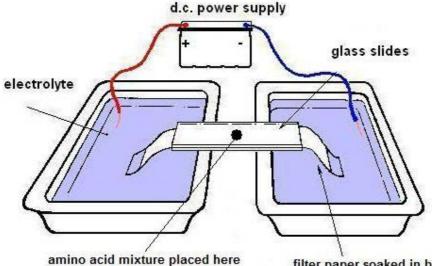
ς,			
	amino acid residue 1	amino acid residue 2	type of bonding
	-HNCH(CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> NH <sub>2</sub> )CO-	$H_2CH_2CH_2CH_2NH_2)CO-$ HNCH(CH <sub>2</sub> CH <sub>2</sub> CO <sub>2</sub> H)CO-	
	-HNCH(CH <sub>3</sub> )CO-	–HNCH(CH <sub>3</sub> )CO–	van der Waals'
	–HNCH(CH₂SH)CO–	-HNCH(CH <sub>2</sub> SH)CO-	Disulfide bonds
	-HNCH(CH <sub>2</sub> OH)CO-	-HNCH(CH <sub>2</sub> CO <sub>2</sub> H)CO-	Hydrogen bonds

[4]

[Total: 10]

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7 (a) Sketch and label the apparatus used to carry out electrophoresis e.g



e filter paper soaked in buffer solution

Marks: power supply / electrolyte + filter paper / buffer / acid mixture central 4 × [1]
[4]

(b)	(i)	pH of the buffer Charge on the amino acid species	[1] [1]
	(ii)	Size of the amino acid species / M <sub>r</sub> Voltage applied Magnitude of the charge (on the amino acid species) Temperature	[1] [1] [1] (max 3) <b>[max 3]</b>
(c)	(i)	They have insufficient electron density / only one electron	[1]

(ii) Sulfur [1] because it has the greatest atomic number / number of electrons [1] [3]

[Total: 10]

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#### 8 (a)

traditional material	modern polymer used
Paper/cardboard/wood/leaves hessian/hemp/jute steel/aluminium	PVC in packaging
Cotton/wool/linen	Terylene in fabrics
Glass/china/porcelain/earthenware metal/leather	Polycarbonate bottle

 $3 \rightarrow 2$  marks,  $2 \rightarrow 1$  mark [2]

(b)	Rea	asons: Plastics/polymers pollute the environment for a long time do not decor biodegrade quickly They are mainly produced from oil Produce toxic gases on burning	mpose/ [1] [1] [1] max two
		ategy 1: Recycle polymer waste / use renewable resources ategy 2: Develop biodegradable polymers	[1] [1] [max 3]
(c)	PVC Combustion would produce HC1/ dioxins as a pollutant or nylon/acrylic Combustion would produce HCN		[1] [1] [1] <b>[2]</b>
(d)	(i)	Polythene (or other addition polymer)	[1]
	(ii)	Addition polymerisation	[1]
		The polymer chains don't have strong bonds between them – easy to melt Could be answered with a suitable diagram	[1] <b>[3]</b>
			[Total: 10]