# MARK SCHEME for the October/November 2011 question paper

## for the guidance of teachers

# 9701 CHEMISTRY

9701/41

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.

Mark schemes must be read in conjunction with the question papers and the report on the examination.

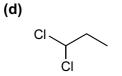
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|                 |        | GCE A LEVEL – O  | ctober/November 2011                | 9701 41                      |                   |
| l (a) (i<br>(ii |        |  | on mixture of $H_2 + Cl_2$ but No   |                              | [1]               |
| ·               | ste    | amy/misty/white fumes pr<br>ntainer gets warm/hot                |                                     |                              | [2]               |
| (iii            | ) H-ŀ  | H = 436  | C <i>l</i> -C <i>l</i> = 244        | H-Cl = 431                   |                   |
|                 | ΔH     | = 436 + 244 - 2(431)   | $= -182 \text{ kJ mol}^{-1}$        |                              | [2]               |
|                 | H-ł    | H = 436  | Br-Br = 193                         | H-Br = 366                   |                   |
|                 | ΔH     | = 436 + 193 - 2(366)   | $= -103 \text{ kJ mol}^{-1}$        |                              | [2]               |
| (iv             | ) H-E  | Br bond is weaker than the                                       | e H-C <i>l</i> bond – allow convers | e.                           | [1]<br><b>[8]</b> |
| (b) (i          | ) ligh | nt   |                                     |                              | [1]               |
| (ii             | ,      | nds broken = C-H & I-I<br>nds made = C-I & H-I<br>$\Delta H$     |                                     | ol <sup>-1</sup>             | [2]               |
| (iii            | ,      | e overall reaction is enc<br>ned <i>or</i> high E <sub>act</sub> | lothermic <i>or</i> no strong bo    | nds/only weak bonds are      | [1]<br><b>[4]</b> |
| (c) (i          | •      | nolytic fission is the brea<br>d-electron species                | 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











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

[Total: 18]

Correct chiral atom identified

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| 2 | (a) (i) | Orde  | er w.r.t. [CH₃CHO] = 1<br>er w.r.t. [CH₃OH] = 1<br>er w.r.t. [H⁺] = 1 |          | [1]<br>[1]<br>[1] |
|   | (ii)    | rate  | = $k[CH_3CHO][CH_3OH][H^+]$   |          | [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 r      | ate = 8) | [1]<br><b>[6]</b> |

(b)

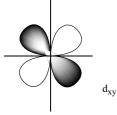
|                | [CH <sub>3</sub> CHO]<br>/mol dm <sup>-3</sup> | [CH <sub>3</sub> OH]<br>/mol dm <sup>-3</sup> | [H <sup>+</sup> ]<br>/mol dm <sup>-3</sup> | [acetal <b>A</b> ]<br>/mol dm <sup>-3</sup> | [H <sub>2</sub> O]<br>/mol dm <sup>-3</sup> |
|----------------|--|---|--|---|---|
| at start       | 0.20   | 0.10  | 0.05                                       | 0.00  | 0.00  |
| at equilibrium | equilibrium (0.20 – x)                         |   | 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]<br>[max 9] |
|-------|---|----------------|
| (iii) | $K_c = \{[acetal A][H_2O]\}/\{[CH_3CHO][CH_3OH]^2\}$<br>units = mol <sup>-1</sup> dm <sup>3</sup> | [1]<br>[1]     |
| (ii)  | 4 values in third row   | 4 x [1]        |
| (i)   | 3 values in second row  | 3 x [1]        |

[Total: 15]

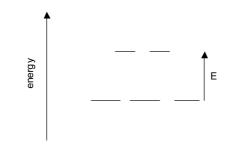
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3 (a) for example.... also allow d<sub>z2</sub>



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

(b) (i)



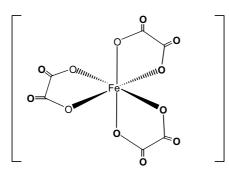
Marks are for 5 degenerate orbitals [1] and 3:2 split [1]

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

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

3-

(iii)



[2]

(iv)  $\underline{2} K_3 \operatorname{Fe}(C_2O_4)_3 \rightarrow \underline{3} K_2C_2O_4 + \underline{2} \operatorname{Fe}C_2O_4 + \underline{2} \operatorname{CO}_2$  $Or K_3 \operatorname{Fe}(C_2O_4)_3 \rightarrow \underline{3/2} K_2C_2O_4 + \operatorname{Fe}C_2O_4 + \operatorname{CO}_2$ [2]

[max 5]

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| 4 | (a) (  |     |                               | $C_2H_5NH_2$ + HA → $C_2H_5NH_3^+$ + A <sup>-</sup> (HA can be H <sub>2</sub> O, HC <i>l</i> etc.)<br>Allow $\rightleftharpoons$ instead of arrow |                        |   |                 |                          |  |
|   | (i     | ii) |                               |   |                        |   |                 |                          |  |
|   |        |     | n                             | nost basic  |                        | least basic                                     |                 |                          |  |
|   |        |     | e                             | thylamine   | ammonia                | phenylamine                                     |                 |                          |  |
|   |        |     |                               |   |                        |   |                 | [1]                      |  |
|   | (ii    |     |                               | -   |                        | ating ethyl/alkyl grou<br>ion of lone pair over |                 | [1]<br>[1]<br><b>[4]</b> |  |
|   | (b) (  | (i) | C <sub>6</sub> H <sub>5</sub> | $_{5}\text{OH} + \text{OH}^{-} \rightarrow$   | $C_6H_5O^- + H_2O$ (or | with Na⁺/H₂O/A⁻)                                |                 | [1]                      |  |
|   | (i     | -   | more<br>stror                 | e than phenol   | he anionic charge      | rger because it's a s<br>is spread out moreov   | -               | [1]                      |  |
|   | (ii    | ii) | рКа                           | = 1.0   |                        |   |                 | [1]                      |  |
|   | (iv    | v)  | Nitro                         | group increas   | es acidity / electro   | n-withdrawing groups                            | increase acidit | y [1]<br>[ <b>5]</b>     |  |
|   |        |     |                               |   |                        |   |                 |                          |  |

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

| , | <br>、 |  |
|---|-------|--|
| ( | )     |  |

| reaction | reagent(s)   | conditions                   |
|----------|--|------------------------------|
| Step 1   | NaNO <sub>2</sub> + HC <i>1</i><br>or HNO <sub>2</sub> [1] | T < 10°C [1]                 |
| Step 2   | H₂O / aq   | heat/boil/T > 10° (both) [1] |
| Step 3   | HNO₃<br>NB HNO₃(aq) OK for both                            | dilute (both) [1]            |
|          |  | 71                           |

[4] **[5]** 

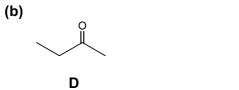
[1]

[Total: 14]

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|        |                                     | •        | -     |

- 5 (a) (i) C=C double bonds / alkenes
  - (ii) -OH groups / accept alcohols or acids
  - (iii) CH<sub>3</sub>CO– or CH<sub>3</sub>CH(OH)– groups
  - (iv) carbonyl, >C=O, groups / accept aldehydes and ketones 4 × [1]



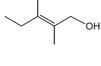






(c) isomers of C





ö

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]<br><b>[2]</b> |
|   | (b)    | (i)  | α he<br>β ple      | lix<br>eated sheet  |                    | [1]<br>[1]        |
|   |        | (ii) | For<br>Nee<br>with | dents should choose one of the structures belowα helix:For β pleated sheed to show a helixNeed to show twoC=O H-Nstrands with C=O -veen turnsthem | parallel 'zig-zag' |                   |
|   |        |      | Whie               | chever is chosen, overall structure [1] position of H bo  | nds [1]            |                   |

| (c) |  |  |                                  |
|-----|--|--|----------------------------------|
|     | 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- | HNCH(CH <sub>2</sub> CH <sub>2</sub> CO <sub>2</sub> H)CO– | lonic bonds or hydrogen<br>bonds |
|     | -HNCH(CH <sub>3</sub> )CO-   | –HNCH(CH₃)CO–  | van der Waals'                   |
|     | -HNCH(CH <sub>2</sub> SH)CO-   | -HNCH(CH <sub>2</sub> SH)CO-                               | Disulfide bonds                  |
|     | –HNCH(CH₂OH)CO–  | -HNCH(CH <sub>2</sub> CO <sub>2</sub> H)CO-                | Hydrogen bonds                   |

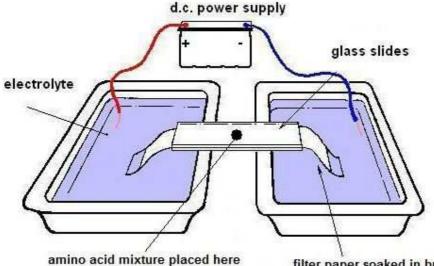
[4]

[4]

[Total: 10]

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



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

because it has the greatest atomic number / number of electrons

[Total: 10]

[1] [3]

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#### 8 (a)

| traditional material  | material modern polymer used |  |
|---|------------------------------|--|
| Paper/cardboard/wood/leaves<br>hessian/hemp/jute<br>steel/aluminium | PVC in packaging             |  |
| Cotton/wool/linen   | Terylene in fabrics          |  |
| Glass/china/porcelain/earthenware<br>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<br>biodegrade quickly<br>They are mainly produced from oil<br>Produce toxic gases on burning | mpose/<br>[1]<br>[1]<br>[1]<br>max two |
|-----|------------------------|--|--|
|     |                        | ategy 1: Recycle polymer waste / use renewable resources<br>ategy 2: Develop biodegradable polymers  | [1]<br>[1]<br>[max 3]                  |
| (c) | PV<br>Coi<br><b>or</b> | C<br>mbustion would produce HC1/ dioxins as a pollutant  | [1]<br>[1]                             |
|     | nyle                   | on/acrylic<br>mbustion would produce HCN   | [1]<br>[1]<br><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]<br><b>[3]</b>                      |
|     |                        |  | [Total: 10]                            |