GCE Advanced Level

MARK SCHEME for the November 2004 question paper

9701 CHEMISTRY

9701/04

Paper 4 (Structured Questions A2 Core), maximum raw mark 60

This mark scheme is published as an aid to teachers and students, to indicate the requirements of the examination. It shows the basis on which Examiners were initially instructed to award marks. It does not indicate the details of the discussions that took place at an Examiners' meeting before marking began. Any substantial changes to the mark scheme that arose from these discussions will be recorded in the published *Report on the Examination*.

All Examiners are instructed that alternative correct answers and unexpected approaches in candidates' scripts must be given marks that fairly reflect the relevant knowledge and skills demonstrated.

Mark schemes must be read in conjunction with the question papers and the *Report on the Examination*.

• CIE will not enter into discussion or correspondence in connection with these mark schemes.

CIE is publishing the mark schemes for the November 2004 question papers for most IGCSE and GCE Advanced Level syllabuses.



Grade thresholds taken for Syllabus 9701 (Chemistry) in the November 2004 examination.

| | maximum | minimum mark required for grade | | | | |
|-------------|-------------------|---------------------------------|----|----|--|--|
| | mark available | А | В | Е | | |
| Component 4 | 60 | 44 | 39 | 22 | | |

The thresholds (minimum marks) for Grades C and D are normally set by dividing the mark range between the B and the E thresholds into three. For example, if the difference between the B and the E threshold is 24 marks, the C threshold is set 8 marks below the B threshold and the D threshold is set another 8 marks down. If dividing the interval by three results in a fraction of a mark, then the threshold is normally rounded down.



November 2004

GCE A LEVEL

MARK SCHEME

MAXIMUM MARK: 60

SYLLABUS/COMPONENT: 9701/04

CHEMISTRY Paper 4 (Structured Questions A2 Core)



| Page 1 | | Mark Scheme | | | | | | Pap | ber | |
|--------|---------------------------|-------------------------|--|---|------------------------------|----------------------|---------------|--------------------|------|----|
| | | | A LEVEL - | - NOVEMBER 2004 | | | 9701 | 4 | ŀ | |
| 1 | (a) (i) | strong, bec | cause final pH is | about 14 | | | | | [1] | |
| | (ii) | (pH = 0.70) | $) \Rightarrow [H^+] = 10^{-0.7}$ | = 0.20 (mol dm ⁻³) |) | | | | [1] | |
| | | | \therefore [H ₂ SO ₄] | = (0.10 mol dm ⁻³) |) | | | ecf | [1] | |
| | (iii) | (end point i | is at 34.0 cm ³ (: | ± 0.5 cm³), so) | | | | | | |
| | | amount | t of H⁺ used | = 0.2 x 25/1000 | = 0.0050 | mol | ecf fror | n (ii) | [1] | |
| | | moles o | of guanidine | = moles of H^+ | = 0.0050 | mol | | | | |
| | | [guanid | line] | = 0.005 x 1000/3 | 4.0 = 0.14 | 1 7 (mo | l dm⁻³) | | [1] | |
| | | | | allow range: 0.14 | ¹ 5 – 0.149 | ecf i | n 0.005 or | 34.0 | | |
| | (iv) |) M _r | = 8.68/0.147 | = 59 (allow range | ə 58 — 60) | | ecf from | 1 (iii) | [1] | 6 |
| | (b) (i) | | → 7 CaSO ₄ | + 3 Ca(H ₂ PO ₄) ₂ + | 2 HF | | | | [1] | |
| | (ii) | M _r values: | Ca(H ₂ PO ₄) ₂ | = 234.1, | $H_2SO_4 =$ | 98.0 | | | [1] | |
| | | | 234.1 x 3 | = 702.3 | 98 x 7 = | 686 | | both | [1] | |
| | | | | ecf from r | atios in equ | ation, | and from I | M _r val | ues | |
| | | | \therefore mass of H ₂ S | SO_4 needed = 1.0 : | x 686/702.3 | 8 = 0.9 | 8 kg | | [1] | |
| | | | (correct answ Allow ecf fron | er = [3] marks. acc n incorrect M _r or inc | curate value correct mult | e is: 0. tipliers | 977 kg. s) | | | 4 |
| | (c) (i) | A solution t | that resists cha | nges in pH [NO | T: results ir | n no p | H change] | | [1] | |
| | | when sma l | II amounts of H | ⁺ or OH ⁻ are addeo | b | | | | [1] | |
| | (ii) | pH = -log ₁₀ | (6.3 x 10 ⁻⁸) + log | g ₁₀ (0.1/0.2) = 6.9 | | | | | [1] | |
| | | or [H⁺] | = (6.3 x 10 ⁻⁸) x | 0.2/0.1 = 1.26 x 1 | 0-7 | | | | | |
| | | ∴ pH | = -log ₁₀ (1.26 x ⁻ | 10 ⁻⁷) = 6.9 | | | | | | 3 |
| | | | | | | | | т | otal | 13 |
| 2 | (a) O ₂ | + 4H⁺ + 4e⁻ | | 2H ₂ O (or equation | on ÷ 2) | | | | [1] | 1 |
| | (b) ⊕ | | | | | | | | [1] | 1 |
| | (c) 1.2 | 3 (V) (ignor | e sign) | | | | | | [1] | 1 |
| | (d) a b | etter/larger | salt bridge <i>or</i> a | diaphragm or large | er (area of) | electr | odes | | | |
| | or i | ncrease cor | ncentrations/pre | essure | | | | | [1] | 1 |
| | | | | | | | | | | |

| | Page 2 | | Mark Scheme A LEVEL – NOVEMBER 2004 | Syllabus 9701 | P | Paper 4 | \neg |
|------|---------------------------|-----------------------|---|------------------|----------|--------------|--------|
| L | (e) time = 400 | 0 x 24 x | 60 x 60 = 34 560 000 seconds | | <u> </u> | [1] |] |
| | charge = o | current | x time = 0.01 x 34 560 000 = 345 600 C | | ecf | [1] | |
| | moles of H | H = 345 | 600/96 500 = 3.6 mol ∴ mass of H = 3.6 | g | ecf | [1] | 3 |
| | (f) advantage | es: | less pollution/CO ₂ /NO _x etc. <i>or</i> cleaner by-product | (S | | | |
| | | | less dependence on fossil fuels/finite resources | any (| one | [1] | |
| | disadvanta | ages: | more expensive (to develop <i>or</i> to run) | | | | |
| | | | takes up more space | | | | |
| | | | poor power-to-volume ratio | | | | |
| | | | hydrogen is difficult to store or to transport | any d | one | [1] | |
| | | | NOT hydrogen is explosive/flammable | | | | 2 |
| | | | | | | Tota | 19 |
| 3 | solubilities de | crease | down the group | | | [1] | |
| | hydration ene | ergy of t | ne cation decreases | | | [1] | |
| | lattice energy | stays t | ne same, or decreases less than H.E. | | | [1] | |
| | making ΔH_{solut} | _{ition} more | e endothermic or H.E. no longer able to overcome | -L.E. | | [1] | 4 |
| | | | | | | Tota | 4 |
| 4 | (a) an elemer | nt formi | ng one or more ions with a partially filled/incomplet | e d-shell | | [1] | 1 |
| | (b) (i) almos | t no cha | ange (allow <i>slight</i> increase or <i>slight</i> decrease) | | | [1] | |
| | (ii) densit | y shoul | d increase | | | [1] | |
| | becau | ise A _r is | increasing but size/volume/radius stays the same | | | [1] | - |
| | (-) | -19 | (allow pa | rtial ect fro | om b | (i)) | 3 |
| (-1) | (c)30 | | | £ | | [1] | 1 |
| (a) | (I) an ion form | mea wn | en a ligand (datively) bonds to a (central metal) ca | tion | | [1] | |
| | (11) | | $H_2 O H_2 H_2 O H_2 H_2 O H_2 H_2 O H_2$ | | | | |

| | Page 3 | | Mark Schen | Mark Scheme Sy | | | |
|---|--|--|--------------------------------------|--|---------------|----------------------|----|
| | | <u> </u> | | | 5701 | | |
| | (e) (i) dark/d | eep/navy/royal | /Oxford blue <i>or</i> pu | rple [NOT Prussian blue | or lilac or r | nauve] [1] | |
| | (ii) 4NH ₃ | + [Cu(H ₂ O) ₆] ²⁺ | | [Cu(NH ₃) ₄ (H ₂ O) ₂] ²⁺ + 4 | H₂O | | |
| | | or | > | [Cu(NH ₃) ₄] ²⁺ + 6H ₂ O | | [1] | 2 |
| | (f) CuCl ₄ ²⁻ is | produced | | | | [1] | |
| | the equilit | prium is revers | ible or ⇒ in equation | on | | [1] | |
| | Cl⁻ ligand | s replace/exc ł | nange with H ₂ O ligation | ands (in words) | | [1] | |
| | (the follow | ing equation is | s worth the first two | marks) | | | |
| | [Cu(H ₂ O)6 | δ] ²⁺ + 4C <i>l</i> ⁻ ⇒ [C | uC1₄]²- + 6H₂O | | | | 3 |
| | | | | | | Total | 12 |
| 5 | (a) (i) A <i>l</i> Cl ₃ / | FeC <i>l</i> ₃/A <i>l</i> /Fe/l₂ (| + heat) [aq negate | s] (N.B. NOT A <i>t</i> Br ₃ etc.) | 1 | [1] | |
| | (or na | mes) | | | | | |
| | (ii) (sun)li | ght/hf/UV (aq r | negates) | | | [1] | 2 |
| | (b) SOC <i>l</i> ₂ /PC | ; <i>l</i> ₃/PC <i>l</i> ₅ [aq neg | jates] | | | [1] | |
| | (or names | \$) | | | | | 1 |
| | (c) (i) C > B | > A (i.e. a mar | k in the penultimate | e box) | | [1] | |
| | (ii) (acyl o | chloride fastest |) highly δ + carbor | atom joined to 2 electro | onegative a | toms | |
| | | or | addition-eliminati | on mechanism is possib | le | [1] | |
| | (aryl c | hloride slowes | t) delocalisation of | lone pair over ring \Rightarrow str | onger C-Cl | bond | |
| | | or | impossibility of 'b | ackside' attack on the C | -Cl bond | [1] | 3 |

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|-----|--------------------|-------------------------------------|---------------------------------------|------------------------------------|---------------------|----------------------|
| (d) | C ₆ H | $_{5}$ -CO $_{2}$ C $_{6}$ H $_{5}$ | C ₆ H₅-CONHCH ₃ | C ₆ H ₅ -C | O ₂ H | |
| | | [1] | [1] | | [1] | |
| | OR | | | | | |
| | | j.O | NHCH3 | Ô | ОН | |
| | | | | | | 3 |
| | | | | | | Total 9 |
| 6 | (a) (i) E | | | | | [1] |
| | (ii) CH₃Cł | H₂CH₂CO₂ ⁻ (Na⁺) | [NOT C ₃ H ₇ CC | 00-Na <i>or</i> C | C₃H⁊COOŀ | -] [1] |
| | | | [but allow CH | $_{3}CH_{2}CH_{2}C$ | O ₂ Na] | |
| | CHl ₃ c | or name | | | | [1] 3 |
| | (b) the alcoho | ol from E has four diffe | erent groups around a carb | on atom | | [1] |
| | ∴ it is chira | al/asymmetric <i>or</i> it is p | produced as a 50:50 mixtur | e of mirror | [·] images | [1] |
| | or its mirr | or images are non-su | perimposable | | | |
| | | CH3 C3H7 OH | Нишини | 3 C ₃ H ₇ | | |
| | formulae: | | i | | | [1] |
| | the alcoho | ol from D has 2 identio | cal groups on its central car | bon atom | | [1] |
| | | | | | | 4 max 3 |
| | | | | | | Total 6 |
| 7 | (a) orange co | lour disappears/brom | ine is decolourised (NOT d | iscoloured | , <i>or</i> goes (| clear) [1] |
| | (white) pre | ecipitate/solid/crystals | is formed | | | [1] 2 |

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|--------|--------|-------------|---|-------------------------------|---|---------------------------|------|----|
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| (b) | e.g. | add | neutral FeCl ₃ | (aq) – viole | et colour with phenol | | | |
| | or | add | universal indi | cator – red | /orange colour with phenol | | | |
| | or | add | Na metal – fiz | | | | | |
| | or | add | NaOH(aq) to | the pure co | ompound – phenol would dissolve | e | | |
| | or | add | H^+ (aq) to the | pure comp | oound – phenylamine would disso | olve | | |
| | or | add | HNO ₂ at room | n temperati | ure – phenylamine would produce | e gaseous N ₂₋ | | |
| | or | add woul | HNO₂ at 5 °C ld produce a c | , followed k coloured (o | by an alkaline solution of phenol - range) dye | - phenylamine | [1] | 1 |
| | (c) IV | KMn | ıO₄ + heat | | | | [1] | |
| | V | HNC |) ₃ + H ₂ SO ₄ | [1] | (both) conc ^d and at 50 $^{\circ}\text{C}$ \cdot | < T < 60 °C | [1] | |
| | VI | Sn + | HC1 (NOT L | _iA <i>1</i> H ₄) | | | [1] | 4 |
| | | | | | | | Tota | 17 |