

MARKSCHEME

November 2001

CHEMISTRY

Standard Level

Paper 2

SECTION A

| 1. | (a) | $\Delta H_{\rm rxn} = \sum {\rm BE}_{\rm breaking} - \sum {\rm BE}_{\rm making}$ | |
|----|-----|---|-------------|
| | | $= (BE_{C=C} + BE_{H-H}) - (BE_{C-C} + 2BE_{C-H})$ | [1] |
| | | =(612+436)-[348+2(412)] | [1] |
| | | =1048-1172 (OR 2696-2820 if all bonds are broken and made) | |
| | | $= -124 \text{ kJ mol}^{-1} (\text{accept} -124 \text{ kJ});$ | [1] |
| | | $(+124 kJ mol^{-1} scores [2])$ | |
| | (b) | 2(-124) = -248 kJ (allow ECF from (a)) | [1] |
| | | Has $2 \text{ C} = \text{C}$, needs 2 H_2 / forms twice the bonds | [1] |
| | | 40 00 6 72 53 28 | |
| 2. | (a) | $n_{\rm C} = \frac{10000}{12.01}; n_{\rm H} = \frac{0002}{1.01} =; n_{\rm O} = \frac{00020}{16.0}$ | |
| | | =3.333 = 6.65 = 3.333 | [1] |
| | | Empirical formula: CH ₂ O | [1] |
| | (b) | $(CH_2O): (12+2+16) = 30;$ molar mass = 2× empirical mass / OWTTE | [1] |
| | | (Award only [1] if reasoning not given) | [1] |
| | (c) | СН.СООН | Ш |
| | (0) | HCOOCH. | [1] |
| | | (Accept other formulas e.g. CH_2OHCHO and $CHOH=CHOH$) | |
| | (d) | pH: CH ₃ COOH pH < 7; HCOOCH ₃ pH = 7 / ester will be higher / acid will be lower | |
| | | Smell: CH ₃ COOH pungent/vinegar smell; HCOOCH ₃ sweet smell | |
| | | Boiling point: CH ₃ COOH higher; HCOOCH ₃ lower | [4] |
| | | ([1] for each test, [1] for results of each test, [4 max]) | |
| 3. | (a) | Mg : 0 Cu^{2+} : + 2 (need both for mark) | [1] |
| | (b) | Cu^{2+} | [1] |
| | (-) | (Accept copper / Cu) | |
| | (c) | Ti (reacts with Ni^{2+} and Cu^{2+}) is a stronger reducing agent than Ni and Cu | [1] |
| | | But weaker reducing agent than Mg Therefore, Mg, Ti, Ni, Cu | [1] [1] |
| | | | Total [20] |
| | | | - vini [#v] |

SECTION B

| 4. | (a) | (i) | Acid: proton / H ⁺ donor | [1] |
|----|-----|--------------|---|------------|
| | | | Base: proton / H ⁺ acceptor | [1] |
| | | (ii) | A pair that differs by a proton / H^+ | [1] |
| | | | H_3O^+/H_2O OR H_2O/OH^- (accept other valid answers) | [1] |
| | | (iii) | Both | [1] |
| | | | Acid: $HCO_3^- + H_2O \rightleftharpoons CO_3^{2-} + H_3O^+$ | [1] |
| | | | Base: $HCO_3^- + H_2O \rightleftharpoons H_2CO_3 + OH^-$ | [1] |
| | | | (States not required; accept single arrow) | |
| | | (iv) | $\mathrm{CO}_3^{2-} + \mathrm{H}_2\mathrm{O} \rightleftharpoons \mathrm{HCO}_3^- + \mathrm{OH}^-$ | [1] |
| | | | Undergoes (base) hydrolysis / reacts with water to produce OH ⁻ OR Products of hydrolysis are a weak acid and strong base | [1] |
| | (b) | Mea | sure pH / use pH meter / use universal indicator | [1] |
| | | Stro | ng acid has lower pH / weak acid has higher pH | [1] |
| | | Mea Stroi | sure conductivity | [1] |
| | | Read | et with magnesium / calcium carbonate | [1] |
| | | Stro | ng acid reacts faster / weak acid reacts slower | [1] |
| | | Mea | sure heat change / temperature rise on adding NaOH | |
| | | Stro | ng acid has greater value / weak acid has lower value | [1] |
| | | (acc | ept any suitable pair of method and difference for [2] each) | |
| | | Stro | ng acid fully ionized / dissociated | [1] |
| | | Wea | k acid partly ionized / dissociated | [1] |
| | (c) | (i) | Urea first in list and NaOH last | [1] |
| | | | Ammonia before caffeine | [1] |
| | | (ii) | Each pH unit represents a tenfold change in acidity | [1] |
| | | | 5 pH units $\Rightarrow 10 \times 10 \times 10 \times 10 \times 10 = 100000$ times different | [1] |
| | | | pH = 12 - 5 = 7 | [1] |
| | | | | Total [20] |

| (a) | (i) | $H_2O(l) \rightleftharpoons H^+(aq) + OH^-(aq) / 2H_2O(l) \rightleftharpoons H_3O^+(aq) + OH^-(aq)$ (States and equilibrium sign needed for second mark) | [2] |
|-----|----------------------|---|--------------------------|
| | (ii) | Endothermic Bond breaking needs energy | [1] [1] |
| | (iii) | Forward reaction K increases, thus more H^+/OH^- formed OR Temperature increase favours forward reaction to use up some of the heat supplied (<i>second mark not awarded for only saying 'due to Le Chatelier's principle'</i>). (Allow ECF from (ii)) | [1] [1] |
| (b) | (i) | All substances are in the same phase / state (Accept all are gases) | [1] |
| | (ii) | No further change in temperature / colour (of iodine) / concentrations | [1] |
| | (iii) | $K_c > 1$: [products] exceed [reactants] at equilibrium / OWTTE $K_c \ll 1$: reaction hardly proceeds / does not proceed / [reactants] \gg [products] | [1] [1] |
| | (iv) | Decreasing volume increases pressure / concentration However no change in equilibrium position Since number of moles of gases the same in reactants and products | [1] [1] [1] |
| | (v) | No effect on position of equilibrium Speeds up both forward and reverse reaction No effect on K_c Value of K_c is affected only by T | [1] [1] [1] [1] |
| (c) | (One 0.25 Ther | e H ₂ reacts with one I ₂ to form 2HI) mol H ₂ and 0.25 mol H ₂ react to form 0.50 mol HI refore, $[H_2] = 0.15 + 0.25 = 0.40$ mol dm ⁻³ $[I_2] = 0.05 + 0.25 = 0.30$ mol dm ⁻³ | [1] [1] [1] |

Total [20]

5.

6. (a) (i) Hydrogen bonding [1] A hydrogen atom bonded to a highly electronegative oxygen atom [1] Strong / special type of dipole-dipole interaction [1] (ii) Van der Waals' / London / dispersion forces [1] Temporary distortion of electron cloud in the non-polar atoms / OWTTE [1] Weak / temporary / instantaneous dipole-dipole interaction [1] [1] (iii) Electrostatic attraction between Na⁺ and Cl⁻/ ions of opposite charge Ions formed due to electron transfer [1]

(b) (i)



[1] [1] [1] [1] [1] (Penalise only once if charge on ion or lone e⁻ pairs on terminal atoms are missing)

| (ii) | PCl ₃ : | trigonal pyramidal | [1] |
|------|---------------------|---|-----|
| | | $<109^{\circ}$ / $\simeq 109^{\circ}$ (but not $109\frac{1}{2}^{\circ}$ or tetrahedral angle) | [1] |
| | NH_2^{-1} : | bent / angular / v-shaped | [1] |
| | _ | $<109^{\circ}/\simeq 109^{\circ}$ (but not $109\frac{1}{2}^{\circ}$) | [1] |
| | NH_4^+ : | tetrahedral | [1] |
| | · | $109\frac{1}{2}^{\circ}$ | [1] |
| | | 2 | |

Total [20]