



JABchem



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Past Papers Higher Chemistry

2014 Marking Scheme

| Grade Awarded | Mark Required (/100) | % candidates achieving grade |
|---------------|----------------------|------------------------------|
| A | 73+ | 30.1% |
| B | 60+ | 23.2% |
| C | 47+ | 23.0% |
| D | 40+ | 10.1% |
| No award | <40 | 13.6% |

| Section: | Multiple Choice | Extended Answer |
|---------------|-----------------|-----------------|
| Average Mark: | 25.9 /40 | 35 /60 |

2014 Higher Chemistry Marking Scheme

| MC Qu | Answer | % Pupils Correct | Reasoning | | | | | | | | | | | | | | | | | | | | | |
|--|--------|------------------|---|---------|-----|-------------------------------------|---|----|----|---------------------------|--------------------------------------|-----|----|----|----|----|--------------------------------|--|-----|-----|-----|-----|-----|-------------------------------------|
| 1 | A | 67 | <input checked="" type="checkbox"/> A Electron Arrangements: $\text{Ca}^{2+} = 2,8,8$ and $\text{Cl}^- = 2,8,8$ <input checked="" type="checkbox"/> B Electron Arrangements: $\text{K}^+ = 2,8,8$ and $\text{Br}^- = 2,8,18,8$ <input checked="" type="checkbox"/> C Electron Arrangements: $\text{Mg}^{2+} = 2,8$ and $\text{Cl}^- = 2,8,8$ <input checked="" type="checkbox"/> D Electron Arrangements: $\text{Na}^+ = 2,8$ and $\text{S}^{2-} = 2,8,8$ | | | | | | | | | | | | | | | | | | | | | |
| 2 | D | 57 | 1mol of NaCl formula units \leftrightarrow 1mol Na^+ ions 1mol of Na_2SO_4 formula units \leftrightarrow 2mol Na^+ ions \therefore 0.6mol NaCl f.u. \leftrightarrow 0.6mol Na^+ ions \therefore 0.2mol Na_2SO_4 f.u. \leftrightarrow 0.4mol Na^+ ions Total number of Na^+ ions = 0.6mol + 0.4mol = 1mol | | | | | | | | | | | | | | | | | | | | | |
| 3 | D | 75 | <input checked="" type="checkbox"/> A Mass number is unchanged as nucleus is unchanged as $\text{X} \rightarrow \text{X}^+ + \text{e}^-$ <input checked="" type="checkbox"/> B The charge on the nucleus is unchanged as the number of electrons decreases <input checked="" type="checkbox"/> C The nucleus is unchanged during $\text{X} \rightarrow \text{X}^+ + \text{e}^-$ \therefore atomic number is unchanged <input checked="" type="checkbox"/> D A group 1 element has 1 outer electron which is lost during $\text{X} \rightarrow \text{X}^+ + \text{e}^-$ | | | | | | | | | | | | | | | | | | | | | |
| 4 | B | 49 | <input checked="" type="checkbox"/> A displacement: higher up metals displaces a lower down metal from its ion <input checked="" type="checkbox"/> B neutralisation: Hydrogen H^+ ions react to become water H_2O <input checked="" type="checkbox"/> C oxidation: increase in the oxygen : hydrogen ratio with electrons being lost <input checked="" type="checkbox"/> D reduction: decrease in the oxygen : hydrogen ratio with electrons being gained | | | | | | | | | | | | | | | | | | | | | |
| 5 | C | 76 | $\begin{array}{rcl} \text{activation energy} & = & \text{activation energy} + \text{enthalpy change} \\ \text{(reverse catalysed reaction)} & = & \text{(catalysed forward reaction)} + (190-160) \\ & = & 35 + 30 \\ & = & 65\text{kJ mol}^{-1} \end{array}$ | | | | | | | | | | | | | | | | | | | | | |
| 6 | A | 78 | <input checked="" type="checkbox"/> A enthalpy change $\Delta H = 100-200 = -100\text{kJ mol}^{-1}$ <input checked="" type="checkbox"/> B enthalpy change $\Delta H = 100-150 = -50\text{kJ mol}^{-1}$ <input checked="" type="checkbox"/> C endothermic reaction as products are higher in potential energy than reactants <input checked="" type="checkbox"/> D endothermic reaction as products are higher in potential energy than reactants | | | | | | | | | | | | | | | | | | | | | |
| 7 | C | 82 | $\text{Rate} = \frac{\Delta \text{quantity}}{\Delta \text{time}} = \frac{0.0062-0.0020}{5-0} = \frac{0.0042}{5} = 0.00084 \text{ mol l}^{-1} \text{ min}^{-1}$ | | | | | | | | | | | | | | | | | | | | | |
| 8 | C | 57 | <input checked="" type="checkbox"/> A Rate decreases as reaction proceeds <input checked="" type="checkbox"/> B Reaction comes to a slow stop not a sudden stop as reaction proceeds <input checked="" type="checkbox"/> C Reaction rate slows as reaction proceeds and comes a gradual stop <input checked="" type="checkbox"/> D Rate decreases as reaction proceeds | | | | | | | | | | | | | | | | | | | | | |
| 9 | A | 51 | <input checked="" type="checkbox"/> A volume of gas given off is less (0.5g chalk) and lump means less steep gradient <input checked="" type="checkbox"/> B lump would give a much slower initial reaction rate and gradient would be less <input checked="" type="checkbox"/> C reduction in mass of chalk to 0.5g would half the volume of gas given off <input checked="" type="checkbox"/> D Reaction rate would be approximately same so gradient would be similar | | | | | | | | | | | | | | | | | | | | | |
| 10 | A | 39 | <input checked="" type="checkbox"/> A Curve R: higher temperature & greater number of particles (area under curve) <input checked="" type="checkbox"/> B Curve R must have a higher temperature than curve Q <input checked="" type="checkbox"/> C Curve R has greater area due to greater number of particles <input checked="" type="checkbox"/> D Curve R must have a higher temperature than curve Q | | | | | | | | | | | | | | | | | | | | | |
| 11 | C | 52 | no. of mol = volume x concentration = 0.1litre x $1\text{mol l}^{-1} = 0.1\text{mol}$ 0.1mol releases -3.1kJ \therefore 1mol releases -31kJ | | | | | | | | | | | | | | | | | | | | | |
| 12 | D | 83 | <input checked="" type="checkbox"/> A boiling points increase due to greater London dispersion forces down group 7 <input checked="" type="checkbox"/> B covalent radius increases as extra shell of electrons is added <input checked="" type="checkbox"/> C 1 st ionisation energy decreases (electrons easier to remove further from nucleus) <input checked="" type="checkbox"/> D Van der Waals' forces are greater as bigger atoms more likely to temp dipole | | | | | | | | | | | | | | | | | | | | | |
| 13 | B | 61 | <table border="1"> <thead> <tr> <th>Element</th> <th>Li</th> <th>Na</th> <th>K</th> <th>Rb</th> <th>Cs</th> <th>Deductions about Francium</th> </tr> </thead> <tbody> <tr> <td>Melting Point ($^{\circ}\text{C}$)</td> <td>181</td> <td>98</td> <td>63</td> <td>39</td> <td>28</td> <td>Less than 28°C</td> </tr> <tr> <td>1st Ionisation Energy (kJ mol^{-1})</td> <td>520</td> <td>496</td> <td>419</td> <td>403</td> <td>376</td> <td>Less than 282 kJ mol^{-1}</td> </tr> </tbody> </table> | Element | Li | Na | K | Rb | Cs | Deductions about Francium | Melting Point ($^{\circ}\text{C}$) | 181 | 98 | 63 | 39 | 28 | Less than 28°C | 1 st Ionisation Energy (kJ mol^{-1}) | 520 | 496 | 419 | 403 | 376 | Less than 282 kJ mol^{-1} |
| Element | Li | Na | K | Rb | Cs | Deductions about Francium | | | | | | | | | | | | | | | | | | |
| Melting Point ($^{\circ}\text{C}$) | 181 | 98 | 63 | 39 | 28 | Less than 28°C | | | | | | | | | | | | | | | | | | |
| 1 st Ionisation Energy (kJ mol^{-1}) | 520 | 496 | 419 | 403 | 376 | Less than 282 kJ mol^{-1} | | | | | | | | | | | | | | | | | | |

| | | | |
|----|---|----|---|
| 14 | A | 56 | <input checked="" type="checkbox"/> A Electronegativities: Be=1.5 & Cl=3.0 ∴ difference = 1.5 ∴ least ionic character <input checked="" type="checkbox"/> B Electronegativities: Ca=1.3 & Cl=3.0 ∴ difference = 1.7 <input checked="" type="checkbox"/> C Electronegativities: Li=1.0 & Cl=3.0 ∴ difference = 2.0 <input checked="" type="checkbox"/> D Electronegativities: Cs=0.8 & Cl=3.0 ∴ difference = 2.2 ∴ most ionic character |
| 15 | B | 72 | <input checked="" type="checkbox"/> A CO ₂ is linear and the molecule is non-polar due to shape <input checked="" type="checkbox"/> B NH ₃ is trigonal pyramidal. Difference in the electronegativities means it is polar <input checked="" type="checkbox"/> C CCl ₄ is tetrahedral and the molecule is non-polar due to shape <input checked="" type="checkbox"/> D CH ₄ is non-polar due to the similar electronegativities of carbon and hydrogen |
| 16 | C | 58 | <input checked="" type="checkbox"/> A Metallic substances (metal elements and alloys) are never compounds <input checked="" type="checkbox"/> B Covalent bonds are found in diatomic elements e.g. H ₂ , N ₂ , O ₂ , F ₂ , Cl ₂ , Br ₂ , I ₂ <input checked="" type="checkbox"/> C Hydrogen bonds are only found in covalent compounds with -OH, -NH or H-F bonds <input checked="" type="checkbox"/> D London dispersion forces are found in every substance |
| 17 | C | 45 | <input checked="" type="checkbox"/> A 1mol C atoms = 12g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{24}{12} = 2\text{mol}$ carbon atoms <input checked="" type="checkbox"/> B 1mol O ₂ molecules = 32g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{16}{32} = 0.5\text{mol}$ O ₂ molecules <input checked="" type="checkbox"/> C 1mol H ₂ molecules = 2g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{2}{2} = 1\text{mol}$ H atoms ∴ 1mol e ⁻ <input checked="" type="checkbox"/> D no. of mol = v x c = 1litre x 1mol l ⁻¹ = 1mol ∴ 1mol NaCl f.u. = 2mol ions |
| 18 | B | 67 | gfm of SO ₂ =64.1g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{128.2}{64.1} = 2\text{mol}$ of SO ₂ <input checked="" type="checkbox"/> A gfm H ₂ =2g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{2.0}{2} = 1\text{mol}$ of H ₂ <input checked="" type="checkbox"/> B gfm He=4g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{8.0}{4} = 2\text{mol}$ of He <input checked="" type="checkbox"/> C gfm O ₂ =32g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{32}{32} = 1\text{mol}$ of O ₂ <input checked="" type="checkbox"/> D gfm Ne=20.2g ∴ no. of mol = $\frac{\text{mass}}{\text{gfm}} = \frac{80.8}{20.2} = 4\text{mol}$ of Ne |
| 19 | B | 36 | <input checked="" type="checkbox"/> A Cracking produces smaller unsaturated hydrocarbons e.g. ethene and propene <input checked="" type="checkbox"/> B Cracking naphtha does not produce petrol directly (bits might be added to petrol) <input checked="" type="checkbox"/> C Reforming can produce benzene-based aromatic hydrocarbons <input checked="" type="checkbox"/> D Reforming can produce ringed cycloalkane hydrocarbons |
| 20 | A | 78 | <input checked="" type="checkbox"/> A Straight chained hydrocarbons are more likely to auto-ignite <input checked="" type="checkbox"/> B Branched-chain hydrocarbons are less likely to auto-ignite <input checked="" type="checkbox"/> C Ringed cycloalkane hydrocarbons are less likely to auto-ignite <input checked="" type="checkbox"/> D Ringed aromatic hydrocarbons are less likely to auto-ignite |
| 21 | B | 79 | <input checked="" type="checkbox"/> A ethanol has a higher relative cost ∴ petrol would cost more <input checked="" type="checkbox"/> B more ethanol required due to lower octane number and higher costs <input checked="" type="checkbox"/> C ethanol has a lower octane rating ∴ more ethanol would be required <input checked="" type="checkbox"/> D ethanol has a higher relative cost ∴ petrol would cost more |
| 22 | D | 69 | Comparison of heptane and 2-methylhexane: 2-methyl branch raises octane number from 0 to 43 Adding 2-methyl branch to 2,4-dimethylhexane to form 2,2,4-trimethylpentane: ∴ adding 2-methyl branch to 2,4-dimethylhexane raises octane number from 66 by similar amount ∴ Octane number of 2,2,4-trimethylpentane = 66 + 43 = 99 |
| 23 | C | 70 | <input checked="" type="checkbox"/> A esters are insoluble in water and are used as solvents <input checked="" type="checkbox"/> B esters are sweet smelling and are used in perfumes <input checked="" type="checkbox"/> C Esters are not used in toothpastes <input checked="" type="checkbox"/> D esters are sweet smelling and are used in flavourings |
| 24 | D | 74 | <input checked="" type="checkbox"/> A Polyester textile fibres are linear and not cross-linked <input checked="" type="checkbox"/> B Polyester fibres are long chain molecules <input checked="" type="checkbox"/> C polyesters are formed by condensation polymerisation <input checked="" type="checkbox"/> D cured polyester resins are strong due to cross-linking of the polyester chains |
| 25 | B | 63 | <input checked="" type="checkbox"/> A ethene is not one of the two monomers used as ethene has only 2 carbons <input checked="" type="checkbox"/> B but-2-ene gives -CH ₃ groups on adjacent carbons and propene has 1x -CH ₃ group <input checked="" type="checkbox"/> C ethene is not one of the two monomers used as ethene has only 2 carbons <input checked="" type="checkbox"/> D but-1-ene is not a monomer used as there are no -C ₂ H ₅ side groups in polymer |

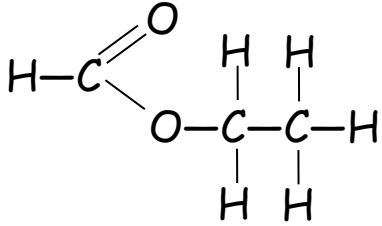
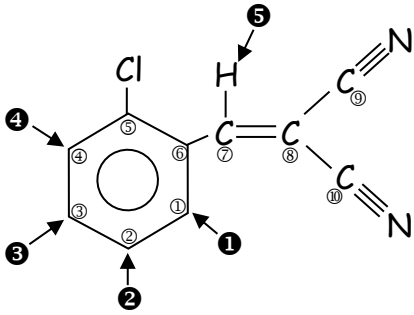
| 26 | D | 65 | <input checked="" type="checkbox"/> A isoprene monomer will form a polymer by addition due to C=C double bonds <input checked="" type="checkbox"/> B No -OH, -NH or H-F groups to form hydrogen bonds <input checked="" type="checkbox"/> C isoprene monomer will form a polymer by addition due to C=C double bonds <input checked="" type="checkbox"/> D addition polymer is formed due to C=C bonds and London dispersion between | | | | | | | | | | | | | | |
|-------------------------------|----|---------------------|--|-------------------------------|----|----------------|--------------|--------------------------|------------------|----|---------------------|--------|-------------|-------------------|----|-------------------|-------|
| 27 | D | 72 | <input checked="" type="checkbox"/> A fats and oils do not cross-link between chains during hardening <input checked="" type="checkbox"/> B hardening of oils into fats is caused by adding hydrogen across C=C bonds <input checked="" type="checkbox"/> C chain length does not change during hardening <input checked="" type="checkbox"/> D number of C=C double bonds decreases during hardening/hydrogenation | | | | | | | | | | | | | | |
| 28 | A | 79 | <input checked="" type="checkbox"/> A Acidified Dichromate gives a orange → green colour change on alcohol oxidation <input checked="" type="checkbox"/> B Benedict's solution does not react with primary alcohols like ethanol <input checked="" type="checkbox"/> C Fehling's solution does not react with primary alcohols like ethanol <input checked="" type="checkbox"/> D Tollen's reagent does not react with primary alcohols like ethanol | | | | | | | | | | | | | | |
| 29 | A | 75 | <input checked="" type="checkbox"/> A Benzene is not a raw material as it has to be extracted from crude oil, etc <input checked="" type="checkbox"/> B Iron ore is iron oxide which is extracted from the ground ∴ raw material <input checked="" type="checkbox"/> C sodium chloride is salt and is extracted from the ground ∴ raw material <input checked="" type="checkbox"/> D water is widely available on this planet ∴ raw material | | | | | | | | | | | | | | |
| 30 | B | 70 | <input checked="" type="checkbox"/> A due to the large quantities involved, iron production is a continuous process <input checked="" type="checkbox"/> B Medicines like aspirin are made by batch process as smaller quantities are required <input checked="" type="checkbox"/> C due to the large quantities involved, ammonia production is a continuous process <input checked="" type="checkbox"/> D due to the large quantities involved, making sulphuric acid is a continuous process | | | | | | | | | | | | | | |
| 31 | C | 65 | $\begin{array}{l} \textcircled{1} \quad C_{(\text{graphite})} + O_2 \rightarrow CO_2 \quad \Delta H = -394 \text{ kJ mol}^{-1} \\ \textcircled{2} \quad C_{(\text{diamond})} + O_2 \rightarrow CO_2 \quad \Delta H = -395 \text{ kJ mol}^{-1} \\ \\ \textcircled{1} \quad C_{(\text{graphite})} + O_2 \rightarrow CO_2 \quad \Delta H = -394 \text{ kJ mol}^{-1} \\ \textcircled{2} \times -1 \quad CO_2 \rightarrow C_{(\text{diamond})} + O_2 \quad \Delta H = +395 \text{ kJ mol}^{-1} \\ \\ \text{Add } \textcircled{1} + \textcircled{2}' \quad C_{(\text{graphite})} \rightarrow C_{(\text{diamond})} \quad \Delta H = +1 \text{ kJ mol}^{-1} \end{array}$ | | | | | | | | | | | | | | |
| 32 | B | 58 | $\begin{array}{l} \textcircled{1} \quad C_2H_4 + 3O_2 \rightarrow 2CO_2 + 2H_2O \quad \Delta H_1 \\ \textcircled{2} \quad CH_3CHO + 2\frac{1}{2}O_2 \rightarrow 2CO_2 + 2H_2O \quad \Delta H_2 \\ \textcircled{3} \quad 2O_3 \rightarrow 3O_2 \quad \Delta H_3 \\ \\ \textcircled{1} \quad C_2H_4 + 3O_2 \rightarrow 2CO_2 + 2H_2O \quad \Delta H_1 \\ \textcircled{2} \times -1 \quad 2CO_2 + 2H_2O \rightarrow CH_3CHO + 2\frac{1}{2}O_2 \quad -\Delta H_2 \\ \textcircled{3} \times \frac{1}{2} \quad O_3 \rightarrow 1\frac{1}{2}O_2 \quad \frac{1}{2}\Delta H_3 \\ \\ \text{Add } \textcircled{1} + \textcircled{2}' + \textcircled{3}' \quad C_2H_4 + O_3 \rightarrow CH_3CHO + O_2 \quad \Delta H_1 - \Delta H_2 + \frac{1}{2}\Delta H_3 \end{array}$ | | | | | | | | | | | | | | |
| 33 | A | 67 | <input checked="" type="checkbox"/> A At equilibrium, rates of forward & reverse reactions are equal At equilibrium, concentrations of reactants and products are constant <input checked="" type="checkbox"/> B At equilibrium, rate of forward reaction = rate of reverse reaction <input checked="" type="checkbox"/> C At equilibrium, concentrations of reactants and products are constant <input checked="" type="checkbox"/> D At equilibrium, concentrations of reactants and products are constant | | | | | | | | | | | | | | |
| 34 | C | 59 | <input checked="" type="checkbox"/> A forward reaction is endothermic ∴ high temperature favours forward reaction <input checked="" type="checkbox"/> B forward reaction is endothermic ∴ high temperature favours forward reaction <input checked="" type="checkbox"/> C low pressure and high temperature favour the production of carbon dioxide <input checked="" type="checkbox"/> D forward reaction increases pressure ∴ low pressure favours forward reaction | | | | | | | | | | | | | | |
| 35 | D | 68 | <table border="1"> <thead> <tr> <th>0.1mol l⁻¹ Alkali</th> <th>pH</th> <th>Mass of Solute</th> <th>Conductivity</th> <th>Moles of HCl neutralised</th> </tr> </thead> <tbody> <tr> <td>sodium hydroxide</td> <td>13</td> <td>higher gfm ∴ higher</td> <td>Higher</td> <td rowspan="2">Same volume</td> </tr> <tr> <td>ammonium solution</td> <td>10</td> <td>lower gfm ∴ lower</td> <td>Lower</td> </tr> </tbody> </table> | 0.1mol l ⁻¹ Alkali | pH | Mass of Solute | Conductivity | Moles of HCl neutralised | sodium hydroxide | 13 | higher gfm ∴ higher | Higher | Same volume | ammonium solution | 10 | lower gfm ∴ lower | Lower |
| 0.1mol l ⁻¹ Alkali | pH | Mass of Solute | Conductivity | Moles of HCl neutralised | | | | | | | | | | | | | |
| sodium hydroxide | 13 | higher gfm ∴ higher | Higher | Same volume | | | | | | | | | | | | | |
| ammonium solution | 10 | lower gfm ∴ lower | Lower | | | | | | | | | | | | | | |
| 36 | B | 77 | <input checked="" type="checkbox"/> A catalyst will result in same final concentration of reactant <input checked="" type="checkbox"/> B catalyst gives steeper initial gradient on graph and same final concentration <input checked="" type="checkbox"/> C catalyst will result in same final concentration of reactant <input checked="" type="checkbox"/> D catalyst will speed up reaction and give steeper initial gradient on graph | | | | | | | | | | | | | | |

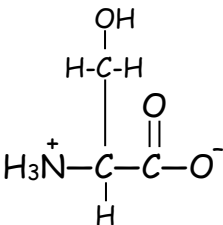
| | | | |
|----|---|----|---|
| 37 | A | 47 | <input checked="" type="checkbox"/> A $Al \rightarrow Al^{3+} + 3e^-$: aluminium metal is oxidised \therefore aluminium metal is reducing agent <input checked="" type="checkbox"/> B $Ag^+ + e^- \rightarrow Ag$: silver ions are reduced \therefore silver ions acting as oxidising agent <input checked="" type="checkbox"/> C $Ag^+ + e^- \rightarrow Ag$: silver ions are reduced \therefore silver ions are electron acceptors <input checked="" type="checkbox"/> D sulphide ions are unchanged and are spectator ions |
| 38 | B | 58 | <input checked="" type="checkbox"/> A the low conductivity of water is due to the low concentrations of ions in water <input checked="" type="checkbox"/> B there are very few ions in pure water and this results in low conductivity <input checked="" type="checkbox"/> C hydrogen bonding between water molecules does not contribute to conductivity <input checked="" type="checkbox"/> D the low conductivity is due to number of ions not the ratio of H^+ and OH^- ions |
| 39 | D | 62 | <input checked="" type="checkbox"/> A concentration of H^+ ions decreases during dilution from pH=4 to pH=6 <input checked="" type="checkbox"/> B concentration of H^+ ions decreases during dilution from pH=4 to pH=6 <input checked="" type="checkbox"/> C pH=4 to pH=6 is a difference in H^+ concentration by a factor of 100 not 2 <input checked="" type="checkbox"/> D Dilution from pH=4 to pH=6 is a decrease in H^+ concentration by a factor of 100 |
| 40 | D | 76 | <input checked="" type="checkbox"/> A beta radiation is negative and is attracted to the positive plate <input checked="" type="checkbox"/> B beta radiation is negative and is attracted to the positive plate <input checked="" type="checkbox"/> C beta radiation is negative and is attracted to the positive plate <input checked="" type="checkbox"/> D alpha bends to negative, beta bends to positive and gamma does not bend |

2014 Higher Chemistry Marking Scheme

| Long Qu | Answer | Reasoning | | | | | | | | | | | | | | | | | | | | |
|--|--|---|--|--|--|--|---------|--|-----------|--|---------|---------------------------------|----------------|--|---------------|---------------------|------------------|---|---------------------------------|--|-----------------------------------|---|
| 1a | <table border="1"> <tr> <td>metallic</td> <td></td> <td>covalent</td> <td></td> </tr> <tr> <td></td> <td>network</td> <td></td> <td>molecular</td> </tr> </table> | metallic | | covalent | | | network | | molecular | <table border="1"> <tr> <th>Bonding</th> <th>1st Twenty Elements</th> </tr> <tr> <td>Metallic solid</td> <td>lithium, beryllium, sodium, magnesium, aluminium, potassium, calcium</td> </tr> <tr> <td>Monatomic gas</td> <td>helium, neon, argon</td> </tr> <tr> <td>Covalent network</td> <td>boron, carbon (diamond), carbon (graphite), silicon</td> </tr> <tr> <td>Discrete covalent molecular gas</td> <td>hydrogen, nitrogen, oxygen, fluorine, chlorine</td> </tr> <tr> <td>Discrete covalent molecular solid</td> <td>sulphur, phosphorus, carbon (fullerene)</td> </tr> </table> | Bonding | 1 st Twenty Elements | Metallic solid | lithium, beryllium, sodium, magnesium, aluminium, potassium, calcium | Monatomic gas | helium, neon, argon | Covalent network | boron, carbon (diamond), carbon (graphite), silicon | Discrete covalent molecular gas | hydrogen, nitrogen, oxygen, fluorine, chlorine | Discrete covalent molecular solid | sulphur, phosphorus, carbon (fullerene) |
| | | metallic | | covalent | | | | | | | | | | | | | | | | | | |
| | | | network | | molecular | | | | | | | | | | | | | | | | | |
| | | Bonding | 1 st Twenty Elements | | | | | | | | | | | | | | | | | | | |
| | | Metallic solid | lithium, beryllium, sodium, magnesium, aluminium, potassium, calcium | | | | | | | | | | | | | | | | | | | |
| | | Monatomic gas | helium, neon, argon | | | | | | | | | | | | | | | | | | | |
| Covalent network | boron, carbon (diamond), carbon (graphite), silicon | | | | | | | | | | | | | | | | | | | | | |
| Discrete covalent molecular gas | hydrogen, nitrogen, oxygen, fluorine, chlorine | | | | | | | | | | | | | | | | | | | | | |
| Discrete covalent molecular solid | sulphur, phosphorus, carbon (fullerene) | | | | | | | | | | | | | | | | | | | | | |
| 1b | Delocalised electrons jump from atom to atom | The electrons in the outer shells of metals are delocalised as they are able to jump from atom to atom allowing electrical conduction through the metal. | | | | | | | | | | | | | | | | | | | | |
| 1c | Same shell filling up and more positive nucleus pulls in outer shell more | Elements in same period have same number of occupied electron shells meaning the element does not increase in size across period. The nucleus becomes increasing positive across a period and this increased charge is attracted to the outer shell more and decreases the size of the atom. | | | | | | | | | | | | | | | | | | | | |
| 2a | naphtha | Petrol is made by reforming the naphtha fraction. Diesel is made by blending the gas oil fraction. | | | | | | | | | | | | | | | | | | | | |
| 2b(i) | diagram showing: | $ \begin{array}{ccccccc} & & \text{H} & & \text{H} & & \\ & & & & & & \\ & & \text{H}-\text{C}-\text{H} & - & \text{H}-\text{C}-\text{H} & & \\ & & & & & & \\ \text{H} & - & \text{C} & - & \text{C} & - & \text{C} & - & \text{C} & - & \text{C} & - & \text{H} \\ & & & & & & & & & & & & \\ & & \text{H} & & \text{H} & & \text{H}-\text{C}-\text{H} & & \text{H} & & \text{H} & & \\ & & & & & & & & & & & & \\ & & & & & & \text{H} & & & & & & \end{array} $ | | | | | | | | | | | | | | | | | | | | |
| 2b(ii) | cycloalkanes or aromatic hydrocarbons | Petrol will auto-ignite before the spark if there is a high a degree of straight chained hydrocarbons. By adding hydrocarbon molecules with branches or rings then the likelihood of auto-ignition before the spark is reduced. | | | | | | | | | | | | | | | | | | | | |
| 2c(i) | Answer to include: | <table border="1"> <tr> <td style="text-align: center;">½ mark</td> <td style="text-align: center;">½ mark</td> </tr> <tr> <td style="text-align: center;">Temperature not too high to denature the enzyme.</td> <td style="text-align: center;">Temperature high enough to give fast reaction.</td> </tr> </table> | ½ mark | ½ mark | Temperature not too high to denature the enzyme. | Temperature high enough to give fast reaction. | | | | | | | | | | | | | | | | |
| ½ mark | ½ mark | | | | | | | | | | | | | | | | | | | | | |
| Temperature not too high to denature the enzyme. | Temperature high enough to give fast reaction. | | | | | | | | | | | | | | | | | | | | | |
| 2c(ii) | Decrease in oxygen : hydrogen ratio | <p>Reduction is defined as a reduction in the oxygen : hydrogen ratio. It can also be viewed as the opposite of oxidation.</p> <table border="1"> <tr> <td style="width: 50%; vertical-align: top;"> <p>Oxidation:</p> <p>ethanol → ethanal</p> <p style="text-align: center;">primary alcohol aldehyde</p> $\begin{array}{ccc} \begin{array}{c} \text{H} \quad \text{H} \\ \quad \\ \text{H}-\text{C}-\text{C}-\text{OH} \\ \quad \\ \text{H} \quad \text{H} \\ \text{C}_2\text{H}_6\text{O} \end{array} & \rightarrow & \begin{array}{c} \text{H} \quad \quad \text{O} \\ \quad \quad // \\ \text{H}-\text{C}-\text{C} \\ \quad \quad \backslash \\ \text{H} \quad \quad \text{H} \\ \text{C}_2\text{H}_4\text{O} \end{array} \end{array}$ </td> <td style="width: 50%; vertical-align: top;"> <p>Reduction:</p> <p>ethanal → ethanol</p> <p style="text-align: center;">aldehyde primary alcohol</p> $\begin{array}{ccc} \begin{array}{c} \text{H} \quad \quad \text{O} \\ \quad \quad // \\ \text{H}-\text{C}-\text{C} \\ \quad \quad \backslash \\ \text{H} \quad \quad \text{H} \\ \text{C}_2\text{H}_4\text{O} \end{array} & \rightarrow & \begin{array}{c} \text{H} \quad \text{H} \\ \quad \\ \text{H}-\text{C}-\text{C}-\text{OH} \\ \quad \\ \text{H} \quad \text{H} \\ \text{C}_2\text{H}_6\text{O} \end{array} \end{array}$ </td> </tr> </table> | <p>Oxidation:</p> <p>ethanol → ethanal</p> <p style="text-align: center;">primary alcohol aldehyde</p> $ \begin{array}{ccc} \begin{array}{c} \text{H} \quad \text{H} \\ \quad \\ \text{H}-\text{C}-\text{C}-\text{OH} \\ \quad \\ \text{H} \quad \text{H} \\ \text{C}_2\text{H}_6\text{O} \end{array} & \rightarrow & \begin{array}{c} \text{H} \quad \quad \text{O} \\ \quad \quad // \\ \text{H}-\text{C}-\text{C} \\ \quad \quad \backslash \\ \text{H} \quad \quad \text{H} \\ \text{C}_2\text{H}_4\text{O} \end{array} \end{array} $ | <p>Reduction:</p> <p>ethanal → ethanol</p> <p style="text-align: center;">aldehyde primary alcohol</p> $ \begin{array}{ccc} \begin{array}{c} \text{H} \quad \quad \text{O} \\ \quad \quad // \\ \text{H}-\text{C}-\text{C} \\ \quad \quad \backslash \\ \text{H} \quad \quad \text{H} \\ \text{C}_2\text{H}_4\text{O} \end{array} & \rightarrow & \begin{array}{c} \text{H} \quad \text{H} \\ \quad \\ \text{H}-\text{C}-\text{C}-\text{OH} \\ \quad \\ \text{H} \quad \text{H} \\ \text{C}_2\text{H}_6\text{O} \end{array} \end{array} $ | | | | | | | | | | | | | | | | | | |
| <p>Oxidation:</p> <p>ethanol → ethanal</p> <p style="text-align: center;">primary alcohol aldehyde</p> $ \begin{array}{ccc} \begin{array}{c} \text{H} \quad \text{H} \\ \quad \\ \text{H}-\text{C}-\text{C}-\text{OH} \\ \quad \\ \text{H} \quad \text{H} \\ \text{C}_2\text{H}_6\text{O} \end{array} & \rightarrow & \begin{array}{c} \text{H} \quad \quad \text{O} \\ \quad \quad // \\ \text{H}-\text{C}-\text{C} \\ \quad \quad \backslash \\ \text{H} \quad \quad \text{H} \\ \text{C}_2\text{H}_4\text{O} \end{array} \end{array} $ | <p>Reduction:</p> <p>ethanal → ethanol</p> <p style="text-align: center;">aldehyde primary alcohol</p> $ \begin{array}{ccc} \begin{array}{c} \text{H} \quad \quad \text{O} \\ \quad \quad // \\ \text{H}-\text{C}-\text{C} \\ \quad \quad \backslash \\ \text{H} \quad \quad \text{H} \\ \text{C}_2\text{H}_4\text{O} \end{array} & \rightarrow & \begin{array}{c} \text{H} \quad \text{H} \\ \quad \\ \text{H}-\text{C}-\text{C}-\text{OH} \\ \quad \\ \text{H} \quad \text{H} \\ \text{C}_2\text{H}_6\text{O} \end{array} \end{array} $ | | | | | | | | | | | | | | | | | | | | | |
| 2c(iii) | 87.1% | $ \begin{array}{ccc} \text{C}_6\text{H}_{12}\text{O}_6 & \longrightarrow & 2\text{C}_2\text{H}_5\text{OH} + 2\text{CO}_2 \\ 1\text{mol} & & 2\text{mol} \\ 180\text{g} & & 92\text{g} \\ 1000\text{g} & & 92\text{g} \times \frac{1000}{180} \\ & & = 511.1\text{g} \end{array} $ $ \% \text{Yield} = \frac{\text{Actual}}{\text{Theoretical}} \times 100 = \frac{445}{511.1} \times 100 = 87.1\% $ | | | | | | | | | | | | | | | | | | | | |
| 3a(i) | Answer to include: | Decrease the volume of potassium iodide solution and increase the volume of water to ensure the total volume of the two solutions is the same. | | | | | | | | | | | | | | | | | | | | |

| | | | | | | | | | | | | | | |
|----------|--|---|---|--|---------------|--------------|--------|----------|---------------|---------|-----------|-------------|--|--|
| 3a(ii) | 2 answers from: | Use a smaller measuring cylinder or syringe for volumes | Use a white tile for observing colour changes | Ensure the beakers are dry before use so concentrations are accurate | | | | | | | | | | |
| 3b | 1 answer from: | Collision angle must be correct | Collision energy must be in excess of the minimum energy to overcome the activation energy. | | | | | | | | | | | |
| 4a | Answer to include: | H-F has hydrogen bonding which raises the b.pt. by bringing the molecules closer together. F-F is non-polar covalent molecular. B.pt. is lower as there is only weak London dispersion forces between molecules. | | | | | | | | | | | | |
| 4b | pH > 7 e.g. 8-11 | The salt of a weak acid and a strong alkali neutralisation reaction gives an alkaline pH when dissolved in water. | | | | | | | | | | | | |
| 5a(i) | condensation | Nylon is a polyamide polymer made by condensation polymerisation of a diacid monomer and a diamine monomer. Other condensation polymers include Kevlar and polyester | | | | | | | | | | | | |
| 5a(ii) | one functional group on monomer | Condensation polymers require monomers with 2 functional groups per monomer so that the polymer chain keeps on going in both directions. When molecules containing only one functional group are added, the polymer chains are prevented from becoming very long as the one-functional group monomer prevents the polymer from extending any further. | | | | | | | | | | | | |
| 5b | conductor | <table border="1"> <tr> <td>Polymer</td> <td>Biopol</td> <td>Poly(ethanol)</td> <td>Poly(ethyne)</td> <td>kevlar</td> </tr> <tr> <td>Property</td> <td>Biodegradable</td> <td>Soluble</td> <td>Conductor</td> <td>Very strong</td> </tr> </table> | Polymer | Biopol | Poly(ethanol) | Poly(ethyne) | kevlar | Property | Biodegradable | Soluble | Conductor | Very strong | | |
| Polymer | Biopol | Poly(ethanol) | Poly(ethyne) | kevlar | | | | | | | | | | |
| Property | Biodegradable | Soluble | Conductor | Very strong | | | | | | | | | | |
| 6 | Answer to include: | OH ⁻ ions in alkali react with H ⁺ ion in equilibrium by neutralisation. Removal of H ⁺ ions (product) from equilibrium moves equilibrium to right to replace H ⁺ ∴ colour becomes more yellow as equilibrium moves to right. | | | | | | | | | | | | |
| 7a | ${}^1_1\text{p}$ | ${}^{14}_7\text{N} + {}^1_0\text{n} \rightarrow {}^{14}_6\text{C} + {}^1_1\text{p}$ <p style="text-align: center;">nitrogen nucleus neutron captured carbon nucleus proton ejected</p> | | | | | | | | | | | | |
| 7b | Neutron splits into proton and electron | <p>During Beta-emission, a neutron splits into a proton and an electron.</p> ${}^1_0\text{n} \rightarrow {}^1_1\text{p} + {}^0_{-1}\text{e}$ <p style="text-align: center;">Neutron in nucleus Proton stays in nucleus β particle ejected</p> | | | | | | | | | | | | |
| 7c(i) | 24225 years | <p>From graph: 5% of carbon-14 content at 4.25 half lives 1 half-life = 5700 years 4.25 half-lives = 5700 years × 4.25/1 = 24225 years</p> | | | | | | | | | | | | |
| 7c(ii) | Too many half-lives have passed to measure | When too many half-lives have passed, it is not possible to measure half-life with any accuracy as there is too little ¹⁴ C left in the sample. | | | | | | | | | | | | |
| 8a | 37.8 | $E_h = c \times m \times \Delta T$ <p>Energy = specific heat capacity × mass × change in temperature</p> <p>Energy = 4.18 × 0.21 × 50</p> <p>Energy = -43.89kJ</p> <p>1mol CaO = (1×40)+(1×16) = 40+16 = 56g</p> <p style="text-align: center;">-65kJ ↔ 56g -43.89kJ ↔ 56g × ^{-43.89}/₋₆₅ = 37.8g</p> | | | | | | | | | | | | |
| 8b | -147 | $\begin{array}{l} \textcircled{1} \quad \text{Ca(s)} + \frac{1}{2}\text{O}_2\text{(g)} \rightarrow \text{CaO(s)} \quad \Delta H = +635\text{kJ mol}^{-1} \\ \textcircled{2} \quad \text{H}_2\text{(g)} + \frac{1}{2}\text{O}_2\text{(g)} \rightarrow \text{H}_2\text{O(l)} \quad \Delta H = -286\text{kJ mol}^{-1} \\ \textcircled{3} \quad \text{Ca(s)} + \text{O}_2\text{(g)} + \text{H}_2\text{(g)} \rightarrow \text{Ca(OH)}_2\text{(s)} \quad \Delta H = +986\text{kJ mol}^{-1} \\ \textcircled{4} \quad \text{Ca(OH)}_2\text{(s)} \rightarrow \text{Ca(OH)}_2\text{(aq)} \quad \Delta H = -82\text{kJ mol}^{-1} \end{array}$ $\begin{array}{l} \textcircled{1} \times -1 \quad \text{CaO(s)} \rightarrow \text{Ca(s)} + \frac{1}{2}\text{O}_2\text{(g)} \quad \Delta H = -635\text{kJ} \\ \textcircled{2} \times -1 \quad \text{H}_2\text{O(l)} \rightarrow \text{H}_2\text{(g)} + \frac{1}{2}\text{O}_2\text{(g)} \quad \Delta H = +286\text{kJ} \\ \textcircled{3} \quad \text{Ca(s)} + \text{O}_2\text{(g)} + \text{H}_2\text{(g)} \rightarrow \text{Ca(OH)}_2\text{(s)} \quad \Delta H = +986\text{kJ} \\ \textcircled{4} \quad \text{Ca(OH)}_2\text{(s)} \rightarrow \text{Ca(OH)}_2\text{(aq)} \quad \Delta H = -82\text{kJ} \end{array}$ <p>Add</p> $\text{CaO(s)} + \text{H}_2\text{O(l)} \rightarrow \text{Ca(OH)}_2\text{(aq)} \quad \Delta H = -147\text{kJ mol}^{-1}$ | | | | | | | | | | | | |

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|---------------|---|--|-------------|---|---------------|--|---------|-----------|---------|-----------------------------|--|--------------|
| 9a(i) | conc sulphuric acid | PPA Technique Question | | | | | | | | | | |
| 9a(ii) | Reactants flammable | PPA Technique Question | | | | | | | | | | |
| 9a(iii) | Diagram Showing: |  | | | | | | | | | | |
| 9b | $\text{CHCl}_3 + 4\text{NaOH} \rightarrow \text{HCOONa} + 3\text{NaCl} + 2\text{H}_2\text{O}$ | | | | | | | | | | | |
| 10a | heterogeneous | <table border="1"> <tbody> <tr> <td>Homogeneous</td> <td>Catalyst in same state as the reactants</td> </tr> <tr> <td>Heterogeneous</td> <td>Catalyst in different state from the reactants</td> </tr> </tbody> </table> | Homogeneous | Catalyst in same state as the reactants | Heterogeneous | Catalyst in different state from the reactants | | | | | | |
| Homogeneous | Catalyst in same state as the reactants | | | | | | | | | | | |
| Heterogeneous | Catalyst in different state from the reactants | | | | | | | | | | | |
| 10b | $\text{CH}_4(\text{g}) + 2\text{H}_2\text{O}(\text{g}) \rightarrow \text{CO}_2(\text{g}) + 4\text{H}_2(\text{g})$ | $\begin{aligned} \textcircled{1} \quad & \text{CH}_4(\text{g}) + \text{H}_2\text{O}(\text{g}) \rightarrow \text{CO}(\text{g}) + 3\text{H}_2(\text{g}) \\ \textcircled{2} \quad & \text{CO}(\text{g}) + \text{H}_2\text{O}(\text{g}) \rightarrow \text{CO}_2(\text{g}) + \text{H}_2(\text{g}) \\ \text{Add } \textcircled{1} + \textcircled{2} \quad & \text{CH}_4(\text{g}) + 2\text{H}_2\text{O}(\text{g}) \rightarrow \text{CO}_2(\text{g}) + 4\text{H}_2(\text{g}) \end{aligned}$ | | | | | | | | | | |
| 10c(i) | 44.8 | $Q = I \times t = 200 \times (30 \times 60) = 360000\text{C}$ $2\text{H}^+ + 2\text{e}^- \longrightarrow \text{H}_2$ <table style="margin-left: auto; margin-right: auto;"> <tbody> <tr> <td style="text-align: center;">2mol</td> <td style="text-align: center;">1mol</td> </tr> <tr> <td style="text-align: center;">2x96500C</td> <td style="text-align: center;">24 litres</td> </tr> <tr> <td style="text-align: center;">193000C</td> <td style="text-align: center;">24 litres</td> </tr> <tr> <td style="text-align: center;">360000C</td> <td style="text-align: center;">24 litres x $360000/193000$</td> </tr> <tr> <td></td> <td style="text-align: center;">= 44.8litres</td> </tr> </tbody> </table> | 2mol | 1mol | 2x96500C | 24 litres | 193000C | 24 litres | 360000C | 24 litres x $360000/193000$ | | = 44.8litres |
| 2mol | 1mol | | | | | | | | | | | |
| 2x96500C | 24 litres | | | | | | | | | | | |
| 193000C | 24 litres | | | | | | | | | | | |
| 360000C | 24 litres x $360000/193000$ | | | | | | | | | | | |
| | = 44.8litres | | | | | | | | | | | |
| 10c(ii) | Water is a renewable source | The hydrogen formed by the electrolysis of water is burned and turns back into water again. This means water will not run out. | | | | | | | | | | |
| 11a | w=10 x=5 y=2 z=1 |  | | | | | | | | | | |
| 11b | 4-methylpentan-2-one | <ol style="list-style-type: none"> 1. Identify functional group: -one 2. Identify longest chain with functional group: -pentanone 3. Give lowest numbering system to functional group: -pentan-2-one 4. Identify side groups: -methylpentan-2-one 5. Give side groups number from existing numbering system: 4-methylpentan-2-one | | | | | | | | | | |
| 11c | CFCs destroy ozone | Chlorofluorocarbons break down ozone. Ozone is required to absorb harmful ultraviolet light from sun. U.V. light can cause sunburn and skin cancer. | | | | | | | | | | |
| 11d(i) | hydrogenation | Addition of hydrogen across C=C double bond is also known as hydrogenation | | | | | | | | | | |

| | | | | | | | | |
|-----|---|---|------------------------------|------------------------------|-------------------------|---|---|---|
| 14b | $C_5H_6N_6O_6$ \downarrow $2CO+4H_2O+3CO_2+2N_2$ | Rule | Quantity | Reaction | Total Moles of Products | | | |
| | | 1 | 5xC | $5C \rightarrow 5CO$ | 5 | | | |
| | | 2 | 4O of total 7O react with 8H | $8H + 4O \rightarrow 4H_2O$ | 5 | 4 | | |
| | | 3 | 3O remaining react with 3CO | $3CO + 3O \rightarrow 3CO_2$ | 2 | 4 | 3 | |
| | | 4 | 4N join to form N_2 | $4N \rightarrow 2N_2$ | 2 | 4 | 3 | 2 |
| 15a |  | It is important to redraw the serine molecule into the same format as the glycine molecule first. Once the $-CH_2OH$ group in serine has been identified, the zwitterion is easy to draw using the template in the example. | | | | | | |
| 15b | Answer to include: | $1 \times 10^{-5} \text{ mol l}^{-1}$ sodium hydroxide solution $\therefore [OH^-] = 1 \times 10^{-5}$ $\therefore [H^+] = 1 \times 10^{-9} \therefore \text{pH} = 9$ As pH is less than IEP pH=9.7 \therefore positive ion is formed \therefore positive ions attracted to negative electrode | | | | | | |