

**Section A**

Questions 1-3 : N/A

Question Number	Correct Answer	Mark
<b>4</b>	C	<b>1</b>

Question Number	Correct Answer	Mark
<b>5</b>	C	<b>1</b>

Question Number	Correct Answer	Mark
<b>6</b>	C	<b>1</b>

Question Number	Correct Answer	Mark
<b>7</b>	A	<b>1</b>

Question Number	Correct Answer	Mark
<b>8</b>	D	<b>1</b>

Question Number	Correct Answer	Mark
<b>9</b>	D	<b>1</b>

Question Number	Correct Answer	Mark
<b>10</b>	D	<b>1</b>

Question Number	Correct Answer	Mark
<b>11</b>	A	<b>1</b>

Question Number	Correct Answer	Mark
<b>12</b>	B	<b>1</b>

Question Number	Correct Answer	Mark
<b>13</b>	B	<b>1</b>

Question Number	Correct Answer	Mark
<b>14</b>	D	<b>1</b>

Question Number	Correct Answer	Mark
<b>15</b>	B	<b>1</b>

Question Number	Correct Answer	Mark
<b>16</b>	B	<b>1</b>

Question Number	Correct Answer	Mark
<b>17</b>	B	<b>1</b>

Question Number	Correct Answer	Mark
<b>18</b>	A	<b>1</b>

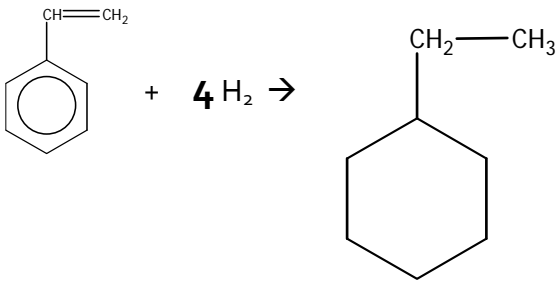
Question Number	Correct Answer	Mark
<b>19</b>	B	<b>1</b>

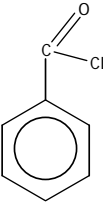
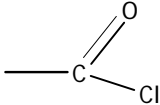
Question Number	Correct Answer	Mark
<b>20</b>	C	<b>1</b>

Question 21 : N/A

Question Number	Acceptable Answers	Reject	Mark
<b>22(a)(i)</b>	Addition / reduction / free-radical addition  IGNORE references to 'hydrogenation'	'redox' 'electrophilic addition' 'nucleophilic addition'	<b>1</b>

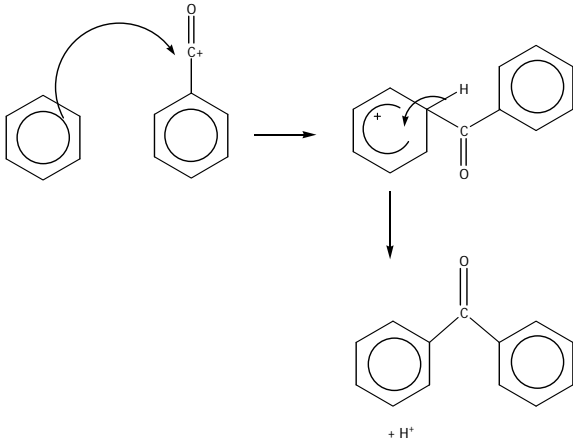
Question Number	Acceptable Answers	Reject	Mark
<b>22(a)(ii)</b>	<p><b>First mark:</b>  <b>Delocalization</b> (of <math>\pi</math>/p electrons in benzene ring) <b>(1)</b></p> <p>IGNORE reference to 'resonance'</p> <p><b>Second mark:</b>  Results in more <b>energy</b> needed to <b>break</b> the <b>bonds</b> in benzene (compared with three separate <math>\pi</math> bonds) <b>(1)</b></p> <p>ALLOW confers <b>stability</b> on the molecule / makes benzene <b>more stable</b> (than expected)</p> <p>IGNORE  Reference to carbon-carbon bond lengths  Values of any enthalpy changes</p> <p>Mark the two points independently</p>		<b>2</b>

Question Number	Acceptable Answers	Reject	Mark
<b>22(a)(iii)</b>	<div style="text-align: center;">  </div> <p style="text-align: center;">( <math>\Delta H =</math> ) <b>— 328</b> (kJ mol<sup>-1</sup>)</p> <p><b>First mark:</b> For “4”</p> <p><b>Second mark:</b> Product as above / correct skeletal formula of product</p> <p>ALLOW Side chain written as —C<sub>2</sub>H<sub>5</sub></p> <p><b>Third mark:</b> —328 (kJ mol<sup>-1</sup>)</p> <p><b>NOTE</b></p> <p><b>One</b> H<sub>2</sub> added showing a CQ correct product with only side chain reduced and cq <math>\Delta H = -120</math> (kJ mol<sup>-1</sup>) scores <b>(2)</b></p> <p><b>Three</b> H<sub>2</sub> added showing a CQ correct product with only the benzene ring reduced and cq <math>\Delta H = -208</math> (kJ mol<sup>-1</sup>) scores <b>(2)</b></p> <p><b>Five</b> H<sub>2</sub> added with fully correct product drawn and <math>\Delta H = -448</math> (kJ mol<sup>-1</sup>) scores <b>(2)</b></p> <p><b>Three and a half</b> H<sub>2</sub> added showing a fully correct product and <math>\Delta H = -268/-293(.3)</math>(kJ mol<sup>-1</sup>) scores <b>(2)</b></p> <p><b>NOTE</b> Mark scoring points independently</p>		<b>3</b>

Question Number	Acceptable Answers	Reject	Mark
<b>22(b)(i)</b>	 <p>Mark awarded for displaying</p> 		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>22(b)(ii)</b>	<p>Electrophilic substitution</p> <p><b>BOTH</b> words needed</p> <p>IGNORE references to 'acylation' and /or 'Friedel-Crafts'</p>		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>22(b)(iii)</b>	<p>Friedel <b>and</b> Crafts</p> <p><b>BOTH</b> names are needed for this mark</p>		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>22(b)(iv)</b>	<p><b>First mark:</b>  <math>\text{C}_6\text{H}_5\text{COCl} + \text{AlCl}_3 \rightarrow \text{C}_6\text{H}_5\text{CO}^+ + \text{AlCl}_4^-</math> <b>(1)</b></p> <p>+ can be anywhere on the <math>\text{C}_6\text{H}_5\text{CO}</math> in the equation for the first mark</p>  <p style="text-align: center;"><math>(\text{AlCl}_4^- + \text{H}^+ \rightarrow \text{HCl} + \text{AlCl}_3)</math></p> <p><b>NOTE:</b>          If ethanoyl chloride or any other acid chloride or the generic <math>\text{RCOCl}</math> is used <b>instead</b> of benzoyl chloride, no first mark can be awarded but the 2nd, 3rd and 4th marks can be awarded consequentially</p> <p><b>Second mark:</b> First curly arrow, as shown, to start from inside the hexagon to the correct <math>\text{C}^+</math> carbon (i.e. not to the benzene ring)          Note the + must be on the C of the <math>\text{C}=\text{O}/\text{CO}</math> for this mark <b>(1)</b></p> <p><b>Third mark:</b> Intermediate correctly drawn <b>(1)</b></p> <p><b>NOTE</b>          + can be shown anywhere in the ring or at the C atom where electrophile is bonded.          The 'horseshoe' in the intermediate to cover at least three carbon atoms</p> <p><b>Fourth mark:</b> Second curly arrow as shown from <math>\text{C}-\text{H}</math> bond to reform the ring, not from the H atom in this bond <b>(1)</b></p> <p><b>NOTE</b>          Products do not have to be shown nor the equation for regeneration of the catalyst given</p>		<p><b>4</b></p> <p style="text-align: right;">6</p>

Question Number	Acceptable Answers	Reject	Mark
<b>22(b)(v)</b>	Absorbs / reflects / blocks / protects from / shields against / <b>uv</b> (light/ radiation) IGNORE 'non-toxic' / references to IR	<b>adsorbs</b> uv light	<b>1</b>

Question Number	Acceptable Answers	Reject	Mark										
22(c)(i)	<p>Any <b>TWO</b> of the following</p> <p>(1) for identifying the bond by formula as shown and (1) for wavenumber in each matching pair</p> <p><b>UNITS</b> are not required</p> <table><tr><th>Bond</th><th>Wavenumber range/wavenumber (cm<sup>-1</sup>)</th></tr><tr><td>C=C</td><td>1600 / 1580 / 1500 / 1450 <b>All four values needed</b></td></tr><tr><td>C=O</td><td>1700 – 1680</td></tr><tr><td>C-H</td><td>3030</td></tr><tr><td>C-H</td><td>750 / 700 <b>Both values needed</b></td></tr></table> <p><b>NOTE</b> ALLOW Correct wavenumber range, or any number within the correct range, <b>for C=O</b></p> <p>Mark identification of the bond and the wavenumber independently (eg a correct bond with a wrong wavenumber, or vice-versa, scores one of the two marks in each case)</p> <p>IGNORE nmr values / chemical shifts</p>	Bond	Wavenumber range/wavenumber (cm <sup>-1</sup> )	C=C	1600 / 1580 / 1500 / 1450 <b>All four values needed</b>	C=O	1700 – 1680	C-H	3030	C-H	750 / 700 <b>Both values needed</b>		4
Bond	Wavenumber range/wavenumber (cm <sup>-1</sup> )												
C=C	1600 / 1580 / 1500 / 1450 <b>All four values needed</b>												
C=O	1700 – 1680												
C-H	3030												
C-H	750 / 700 <b>Both values needed</b>												

Question Number	Acceptable Answers	Reject	Mark
<b>22(c)(ii)</b>	<div data-bbox="363 331 821 646" data-label="Chemical-Block"> </div> <p data-bbox="336 674 507 705"><b>First mark</b></p> <p data-bbox="336 747 459 779"><b>EITHER</b></p> <p data-bbox="336 783 783 852">Identifies correctly the <b>three</b> different proton environments</p> <p data-bbox="336 890 440 921">ALLOW</p> <p data-bbox="336 926 823 1031">If the three different proton environments are only shown on one of the benzene rings</p> <p data-bbox="336 1068 427 1100"><b>NOTE</b></p> <p data-bbox="336 1104 855 1173">On right-hand ring, clockwise from C=O, positions 2, 3 and 4</p> <p data-bbox="336 1178 807 1247">And /or 2,4 and 5 are shown as different environments</p> <p data-bbox="336 1251 443 1283">and /or</p> <p data-bbox="336 1287 823 1356">On left-hand ring, anti-clockwise from C=O, positions 2, 3 and 4</p> <p data-bbox="336 1360 807 1430">And /or 2,4 and 5 are shown as different environments</p> <p data-bbox="336 1467 384 1499"><b>OR</b></p> <p data-bbox="336 1537 874 1606">Identifies proton Z correctly on both benzene rings (1)</p> <p data-bbox="336 1644 552 1675"><b>Second mark</b></p> <p data-bbox="336 1680 815 1749">Fully correct labelling both rings using the letters <b>X, Y</b> and <b>Z</b></p> <p data-bbox="336 1787 427 1818"><b>NOTE</b></p> <p data-bbox="336 1822 868 1892"><b>X</b> and <b>Y</b> labels are interchangeable, <b>Z</b> is not (1)</p>		<b>2</b>

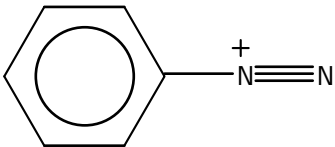
Total for Question 22 = 20 Marks

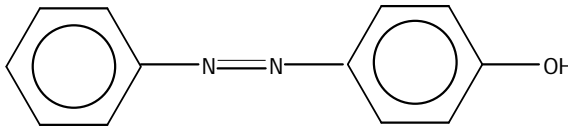


Question Number	Acceptable Answers	Reject	Mark
<b>23(a)(i)</b>	<b>Lone pair</b> (of electrons on the nitrogen atom) ALLOW non-bonded pair (of electrons on the nitrogen atom)	<b>Lone pairs</b> <b>Spare pair</b>	<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23(a)(ii)</b>	<p><b>(with H<sub>2</sub>SO<sub>4</sub>)</b></p> <p><math>(\text{C}_4\text{H}_9\text{NH}_3^+)_2\text{SO}_4^{2-}</math> <b>(1)</b></p> <p>ALLOW</p> <p><math>\text{C}_4\text{H}_9\text{NH}_3^+\text{HSO}_4^-</math></p> <p><b>(with CH<sub>3</sub>COOH)</b></p> <p><math>\text{C}_4\text{H}_9\text{NH}_3^+\text{CH}_3\text{COO}^-</math> <b>(1)</b></p> <p><b>CHARGES</b> not essential</p> <p>Cation and anion can be in either order</p> <p>Max <b>(1)</b> if formula of the amine is incorrect in either case</p> <p>ALLOW (1) if only the correct cation is given in each case (i.e. the anion has been omitted in both cases)</p> <p><b>NOTE</b> The correct ions can be shown separately Eg <math>(\text{C}_4\text{H}_9\text{NH}_3^+)_2 + \text{SO}_4^{2-}</math></p>		<b>2</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23(b)</b>	<p>Tin / Sn ALLOW Iron / Fe <b>(1)</b></p> <p>(concentrated) <b>hydrochloric acid</b></p> <p><b>NOTE</b> If they write 'HCl', there must be some indication of concentrated Eg 'conc HCl' / 'concentrated HCl'</p> <p>ALLOW HCl(aq)</p> <p>(Followed by addition of alkali to liberate the free amine) <b>(1)</b></p> <p>Mark the two points independently</p> <p><b>NOTE</b> Do not allow 2<sup>nd</sup> mark if there is a suggestion that the acid and alkali are added together simultaneously</p>	<p>LiAlH<sub>4</sub></p> <p><b>Just 'HCl'</b></p> <p><b>'dilute'</b> hydrochloric acid / sulfuric acid</p>	<b>2</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23(c)(i)</b>	 <p><b>NOTE</b> If the above structure is drawn, the + charge must be on the N connected directly to the benzene ring</p> <p>ALLOW —N=N<sup>+</sup> on ring</p> <p>IGNORE Cl<sup>-</sup></p>	N <sub>2</sub> <sup>+</sup> on ring	<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23(c)(ii)</b>			<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23(c)(iii)</b>	<p><b>(Conditions)</b>  (Presence of) NaOH / KOH / alkali / OH<sup>-</sup> <b>(1)</b></p> <p>ALLOW  'Alkaline (conditions)' or 'base' or 'high pH'</p> <p>IGNORE  Any references to temperature</p> <p><b>(Use)</b>  Dye / pigment / colouring / indicator / in foodstuff / in paint / methyl orange <b>(1)</b></p> <p>IGNORE  Any reference to medicines</p>		<b>2</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23 (d)</b>	<div data-bbox="359 336 933 430" data-label="Chemical-Block"> <math display="block">\text{C}_6\text{H}_5\text{N}_2^+ + \text{H}_2\text{O} \rightarrow \text{C}_6\text{H}_5\text{OH} + \text{N}_2 + \text{H}^+</math> </div> <p>ALLOW The + sign to be on either N atom in the benzenediazonium ion</p> <p>OR</p> $\text{C}_6\text{H}_5\text{N}_2^+ + \text{H}_2\text{O} \rightarrow \text{C}_6\text{H}_5\text{OH} + \text{N}_2 + \text{H}^+$ <p>OR</p> $\text{C}_6\text{H}_5\text{N}_2\text{Cl} + \text{H}_2\text{O} \rightarrow \text{C}_6\text{H}_5\text{OH} + \text{N}_2 + \text{HCl}$ <p>OR</p> $\text{C}_6\text{H}_5\text{N}_2^+ + 2\text{H}_2\text{O} \rightarrow \text{C}_6\text{H}_5\text{OH} + \text{N}_2 + \text{H}_3\text{O}^+$ <p>OR</p> $\text{C}_6\text{H}_5\text{N}_2^+\text{Cl}^- + \text{H}_2\text{O} \rightarrow \text{C}_6\text{H}_5\text{OH} + \text{N}_2 + \text{HCl}$ <p><b>NOTE</b> —C<sub>6</sub>H<sub>5</sub> can be written or drawn</p> <p><b>First mark</b> for <b>N<sub>2</sub></b> <span style="float: right;"><b>(1)</b></span></p> <p><b>Second mark</b> for rest of the equation correct <span style="float: right;"><b>(1)</b></span></p> <p>IGNORE State symbols, even if incorrect</p>		<b>2</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23(e)(i)</b>	<p>(Otherwise) too much (product) remains in solution OR If excess (solvent) is used, crystals might not form</p> <p>ALLOW To avoid losing (too much) product (in the filtrate when crystallization occurs) / 'to maximize the yield'/ 'will crystallize better from a concentrated solution'/ 'will recrystallize (better) when cold'</p> <p>IGNORE References to a 'saturated solution' or references to 'dilution' or references to the time taken for crystals to form</p>		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23(e)(ii)</b>	<p><b>(Insoluble impurities removed)</b> By <b>hot</b> filtration / During the <b>first</b> filtration / During the <b>second</b> step in the process <b>(1)</b></p> <p><b>(Soluble impurities removed)</b> By remaining in solution / Left in filtrate / Removed when washed (with cold solvent) <b>(1)</b></p>		<b>2</b>

Question Number	Acceptable Answers	Reject	Mark
<b>23(e)(iii)</b>	<p>Measure the melting temperature / melting point  <b>and</b>  compare with data / known value  (from a data book / literature / Internet /data base)  (BOTH points needed for the mark)</p> <p>OR</p> <p>The melting point is sharp  (Just this statement is needed for the mark)</p> <p>ALLOW  Any form of chromatography</p> <p>IGNORE  References to any types of spectroscopy</p>	<b>(0)</b> if reference to determination of the boiling point is made	<b>1</b>

**Total for Question 23 = 15 Marks**

Question Number	Acceptable Answers	Reject	Mark
<b>24(a)(i)</b>	$\text{TiCl}_4 + 4\text{Na} \rightarrow 4\text{NaCl} + \text{Ti}$  IGNORE State symbols, even if incorrect  ALLOW Multiples Reversible arrows		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24(a)(ii)</b>	<p>Ti <b>reduced</b> as oxidation number decreases from <b>+4 to 0</b>/ changes from <b>+4 to 0</b></p> <p style="text-align: right;"><b>(1)</b></p> <p>Na <b>oxidized</b> as oxidation number increases from <b>0 to +1</b> /changes from <b>0 to +1</b></p> <p style="text-align: right;"><b>(1)</b></p> <p>ALLOW Correct oxidation numbers only for one mark</p> <p><b>NOTE</b> Max (1) if no + sign included</p> <p>ALLOW '<b>4+</b>' and/or '<b>1+</b>' given instead of <b>+4</b> and <b>+1</b></p> <p><b>NOTE</b> If any of the oxidation numbers are wrong, award max (1) for the idea that during oxidation the oxidation number increases <b>AND</b> during reduction the oxidation number decreases</p> <p>IGNORE References to loss and /or gain of electrons</p>		<b>2</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24(b)</b>	<p>(Ti [Ar]) <math>3d^2 4s^2 / 4s^2 3d^2</math> <b>(1)</b></p> <p>(Ti<sup>3+</sup> [Ar]) <math>3d^1 / 3d^1 4s^0</math>            (Ti<sup>4+</sup> [Ar]) 'nil' / <math>3d^0 4s^0 / 3d^0</math>            space left blank by candidate</p> <p>BOTH Ti<sup>3+</sup> and Ti<sup>4+</sup> correct for second mark <b>(1)</b></p> <p>Mark CQ on Ti electron configuration for the second mark</p> <p>ALLOW            Upper case (e.g. 'D' for 'd' in electronic configurations)            Subscripts for numbers of electrons</p> <p>Full correct electronic configurations  <math>1s^2, 2s^2, \dots</math></p>		<b>2</b>



Question Number	Acceptable Answers	Reject	Mark
<b>24(c)(i)</b>	<p><b>(d-block element)</b></p> <p><b>EITHER</b>            Ti has (two) electrons in the 3d subshell /            Ti has a partially filled d-subshell /            Ti has a partially filled d-orbital /            Ti has electrons in d-orbital(s) /            Ti has electrons in d-subshell            (During the build up of its atoms) last added / valence electron is in a d-subshell / d-orbital</p> <p><b>OR</b>            (During the build up of its atoms) last added / valence electron is in a d-subshell / d-orbital</p>	<p><b>Outer / highest energy</b> electrons are in a d-orbital /  <b>Outer / highest energy</b> electrons are in a d-subshell</p> <p>Electrons in the 'd-block' /            'electrons in the d-shell'</p>	<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24(c)(ii)</b>	<p>(transition element)</p> <p>Forms one (or more stable) <b>ions</b> /            forms <b>Ti<sup>3+</sup></b> (ions) which have</p> <p>incomplete d-orbital(s) /            an incomplete d-subshell /            a partially filled d-subshell /            an unpaired d electron</p> <p>IGNORE            References to variable oxidation states</p>		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24(d)(i)</b>	<p><b>First mark:</b> d-subshell splits /d-orbitals split (in energy by ligands) /d energy level(s) split(s) <b>(1)</b></p> <p><b>Second mark:</b> <b>absorbs light</b> (in visible region) <b>(1)</b></p> <p><b>Third mark:</b> Electron transitions from lower to higher energy / electron(s) jump from lower to higher energy OR Electron(s) promoted (within d) <b>(1)</b> Mark independently</p> <p><b>NOTE</b> Maximum of (1) mark (i.e. the first mark only) if refers to electrons falling back down again</p>	<p>d-<b>orbital</b> / d-<b>shell</b> splits</p> <p>absorbs <b>purple</b> light</p>	<b>3</b>

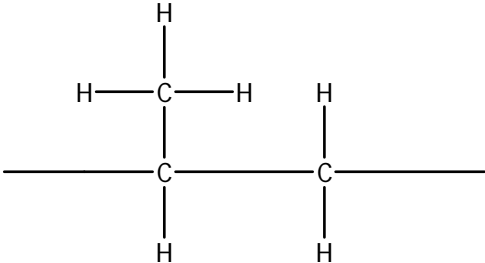
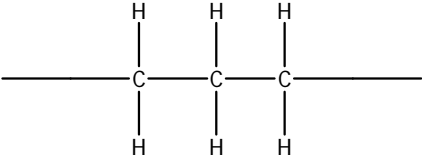
Question Number	Acceptable Answers	Reject	Mark
<b>24(d)(ii)</b>	No d-electrons / empty d-subshell		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24(e)(i)</b>	<p><b>TiO<sub>2</sub></b> <b>'Structure' mark</b></p> <p><b>EITHER</b></p> <p>Giant (structure) <b>OR</b> Lattice (structure)</p> <p>IGNORE Whether stated as ionic or covalent for this mark <b>(1)</b></p> <p><b>TiO<sub>2</sub></b> <b>'Bonding' mark</b></p> <p>EITHER</p> <p><b>Strong</b> (electrostatic) attraction between ions</p> <p>ALLOW <b>Strong</b> ionic bonds / ionic bonds require a lot of energy to break</p> <p><b>OR</b></p> <p><b>Strong</b> covalent bonds/covalent bonds require a lot of energy to break <b>(1)</b></p> <p><b>TiCl<sub>4</sub></b> <b>'Structure' mark</b></p> <p>(Simple) molecules / (small) molecules /molecular <b>(1)</b></p> <p><b>TiCl<sub>4</sub></b> <b>'Bonding' mark</b></p> <p><b>Weak</b> London / dispersion / van der Waals' forces (between molecules) / London /dispersion / van der Waals' forces (between molecules) require little energy to break</p>	<p>TiO<sub>2</sub> (small) molecules / simple molecular</p> <p>For TiO<sub>2</sub> mention of any type of intermolecular forces between molecules of TiO<sub>2</sub></p> <p>TiCl<sub>4</sub> giant structure</p> <p>Covalent bonds broken (on melting) in TiCl<sub>4</sub></p> <p>Ionic bonding in TiCl<sub>4</sub></p> <p>Hydrogen bonding <b>(0)</b> for this mark</p>	<p><b>4</b></p>

	<p><b>NOTE</b></p> <p>If candidates assumes <math>\text{TiO}_2</math> and <math>\text{TiCl}_4</math> are both simple molecular, can score last mark for saying that the <b>named</b> intermolecular forces in <math>\text{TiO}_2</math> are stronger than those in <math>\text{TiCl}_4</math></p> <p>IGNORE (Permanent) dipole-dipole forces <b>(1)</b></p> <p>Mark the four scoring points independently</p>		
--	---	--	--

Question Number	Acceptable Answers	Reject	Mark
<b>24(e)(ii)</b>	Amphoteric ALLOW Recognisable spellings		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24(e)(iii)</b>	$\text{TiO}_2 + 2\text{H}_2\text{O} + 2\text{KOH} \rightarrow \text{K}_2\text{Ti}(\text{OH})_6$ OR $\text{TiO}_2 + 2\text{H}_2\text{O} + 2\text{OH}^- \rightarrow \text{Ti}(\text{OH})_6^{2-}$  IGNORE state symbols even if incorrect		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24(e)(iv)</b>	 <p><b>MUST</b> have continuation bonds at each end ALLOW CH<sub>3</sub></p> <p>IGNORE n and any brackets</p>	 <p><b>Two</b> (or more) repeat units shown</p>	<b>1</b>
<b>24(f)(i)</b>	$(\text{H}_2\text{O}_2 + 2\text{H}^+ +) \mathbf{2e^{(-)}} \rightarrow \mathbf{2\text{H}_2\text{O}}$ <p><b>BOTH</b> <math>2e^{(-)}</math> <b>and</b> <math>2\text{H}_2\text{O}</math> needed for the mark</p>		<b>1</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24(f)(ii)</b>	<p>(Moles <math>\text{H}_2\text{O}_2 = \frac{0.0200 \times 22.50}{1000}</math>  <math>\Rightarrow 4.5 \times 10^{-4} \text{ mol H}_2\text{O}_2</math> <b>(1)</b></p> <p>(Moles <math>\text{Ti}^{3+}</math> reacting in <math>25.0 \text{ cm}^3</math>) <math>= 9.0 \times 10^{-4} \text{ mol Ti}^{3+}</math></p> <p>(Moles <math>\text{Ti}^{3+}</math> in <math>250 \text{ cm}^3</math>) <math>= 9.0 \times 10^{-3} \text{ mol Ti}^{3+}</math> <b>(1)</b></p> <p>(Original concentration of <math>\text{Ti}^{3+}</math>  <math>= \frac{9.0 \times 10^{-3}}{0.00500}</math>  <math>\Rightarrow 1.8 \text{ (mol dm}^{-3}\text{)}</math> <b>(1)</b></p> <p><math>1.8 \text{ (mol dm}^{-3}\text{)}</math> with or without working scores <b>(3)</b></p> <p><b>NOTES:</b>          If mole ratio <math>\text{H}_2\text{O}_2 : \text{Ti}^{3+}</math> is 1:1          final answer for concentration of <math>\text{Ti}^{3+}</math> is <math>0.9 \text{ (mol dm}^{-3}\text{)}</math> scores <b>(2)</b> overall</p> <p>If mole ratio <math>\text{H}_2\text{O}_2 : \text{Ti}^{3+}</math> is 2:1          final answer for concentration of <math>\text{Ti}^{3+}</math> is <math>0.45 \text{ (mol dm}^{-3}\text{)}</math> scores <b>(2)</b> overall</p> <p>If candidate forgets to multiply no. of moles of <math>\text{Ti}^{3+}</math> by 10 then answer is <math>0.18 \text{ (mol dm}^{-3}\text{)}</math> this scores <b>(2)</b></p> <p>If volume of <math>\text{H}_2\text{O}_2</math> used is 25.0 no first mark, but can score <b>(2)</b> if final answer CQ is <math>2(.0) \text{ (mol dm}^{-3}\text{)}</math></p>		<b>3</b>

Question Number	Acceptable Answers	Reject	Mark
<b>24</b> <b>(f)(iii)</b>	(It/titanium(III)/Ti <sup>3+</sup> ) <b>oxidized</b> (by oxygen in the air)  ALLOW 'It is a <b>strong</b> reducing agent'	Hydrolysis	<b>1</b>

Question Number	Correct Answer	Mark
25	D	1

Question Number	Correct Answer	Mark
26	B	1

Question Number	Correct Answer	Mark
27	A	1

Question Number	Correct Answer	Mark
28 (a)	C	1

Question Number	Correct Answer	Mark
28 (b)	B	1

Question Number	Correct Answer	Mark
28 (c)	D	1

Question Number	Correct Answer	Mark
29	B	1



Question Number	Correct Answer	Mark
30	D	1

Question Number	Correct Answer	Mark
31	D	1

Question Number	Correct Answer	Mark
32	B	1

Question Number	Correct Answer	Mark
33	A	1

Question Number	Acceptable Answers	Reject	Mark
34 (a)(i)	(Acid) hydrolysis	substitution	1

Question Number	Acceptable Answers	Reject	Mark
(a)(ii)	$K_2Cr_2O_7$ / $Na_2Cr_2O_7$ / $Cr_2O_7^{2-}$ Potassium dichromate(VI) / sodium dichromate(VI) / dichromate(VI) ions  <i>ALLOW</i> manganate(VII) ions, etc	Just "dichromate"  chromates  Correct formula with wrong name and vice versa  Incorrect oxidation number	1

Question Number	Acceptable Answers	Reject	Mark
(a)(iii)	Lithium tetrahydridoaluminate/ lithium aluminium hydride/ $LiAlH_4$ (in dry ether)	Just $[H^-]$	1

Question Number	Acceptable Answers	Reject	Mark
(a)(iv)	Methyl butanoate (1)  $CH_3CH_2CH_2COOH + CH_3OH \rightarrow CH_3CH_2CH_2COOCH_3 + H_2O$ (1)  <i>ALLOW</i> $\rightleftharpoons$ <i>IGNORE</i> state symbols even if wrong	Methyl butoate	2

Question Number	Acceptable Answers	Reject	Mark
34 (a)(v)	$\text{CH}_3-\text{CH}_2-\text{CH}_2-\text{C} \begin{array}{l} \text{O} \\ \parallel \\ \text{Cl} \end{array}$ <p>Don't penalise undisplayed methyl groups as here. COCl must be displayed as above.</p>	C <sub>3</sub> H <sub>7</sub> for CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub>	1

Question Number	Acceptable Answers	Reject	Mark
(b)(i)	<p>Nitrogen inert / unreactive / less reactive (than oxygen)</p> <p>OR</p> <p>Oxygen might react with chemicals going through column / sample might oxidise</p>		1

Question Number	Acceptable Answers	Reject	Mark
(b)(ii)	<p>Solubility (in liquid / stationary phase)</p> <p>OR</p> <p>Interaction with liquid / stationary phase</p> <p>OR</p> <p>Interaction between mobile and stationary phase</p> <p>OR</p> <p>Attraction for liquid / stationary phase</p> <p>OR</p> <p>Strength of (named) intermolecular forces</p> <p>OR</p> <p>Adsorption on liquid / stationary phase</p> <p>OR</p> <p>Absorption on liquid / stationary phase</p>	<p>Size of molecule / molar mass</p> <p>Polarity, unless with explanation</p> <p>Boiling point / volatility</p> <p>Viscosity</p> <p>Attraction for carrier gas</p> <p>Just a named intermolecular force</p> <p>Just 'retention time'</p> <p>Density</p>	1

Question Number	Acceptable Answers	Reject	Mark
34 (c)(i)	$\left[ \text{O}-\underset{\text{CH}_3}{\overset{\text{H}}{\text{C}}}-\text{CH}_2-\overset{\text{O}}{\underset{\text{  }}{\text{C}}}-\text{O}-\underset{\text{CH}_3}{\overset{\text{H}}{\text{C}}}-\text{CH}_2-\overset{\text{O}}{\underset{\text{  }}{\text{C}}} \right]$ <p>OR</p> $\left[ \underset{\text{CH}_3}{\overset{\text{H}}{\text{C}}}-\text{CH}_2-\overset{\text{O}}{\underset{\text{  }}{\text{C}}}-\text{O}-\underset{\text{CH}_3}{\overset{\text{H}}{\text{C}}}-\text{CH}_2-\overset{\text{O}}{\underset{\text{  }}{\text{C}}}-\text{O} \right]$ <p>Ester link including C=O (1)  Rest of polymer with oxygens at end correct (1)</p> <p>All H atoms must be shown.</p> <p><i>PENALISE</i> lack of displayed C=O once only  <i>ACCEPT</i>  Without brackets around formula but bonds at end should be shown  More than two correct units  <i>IGNORE</i> n after brackets</p>		2

Question Number	Acceptable Answers	Reject	Mark
(c)(ii)	Hydrolysis OR Splits / breaks ester link OR polymer breaks down to monomers OR equation showing hydrolysis	Just 'breaks polymer down'	1