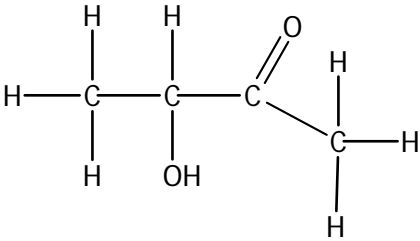




Question Number	Acceptable Answers	Reject	Mark
19(c)(i)	<p><b>First mark:</b> (X is neutral) so not a (carboxylic) acid (1)</p> <p><b>IGNORE</b> "X doesn't have a charge as it is neutral" / "X is not an alkali" / "X is not a base"</p> <p><b>Second mark:</b> (X does not react with Tollens') so is not an aldehyde / is a ketone (1)</p> <p><b>Third mark:</b> (X reacts with <math>H^+</math> / <math>Cr_2O_7^{2-}</math> so) is an alcohol / contains an OH (group) / contains R—OH / contains hydroxyl (group) (1)</p> <p><b>IGNORE</b> 'not an acid' if this is deduced <b>solely</b> from the <math>H^+</math> / <math>Cr_2O_7^{2-}</math> information</p> <p><b>Fourth mark:</b> a primary or a secondary (alcohol) <b>both needed</b> <b>OR</b> (X is) not tertiary (alcohol) (1)</p> <p><b>Mark each point separately</b></p> <p><b>NOTE:</b> 'X is a primary or a secondary alcohol' scores both the third and fourth marks</p> <p><b>ALLOW</b> Correct formulae for the functional groups, instead of their names</p>	<p>X is an aldehyde scores (0) for this scoring point / X is not a ketone scores (0) for this scoring point</p>	4

Question Number	Acceptable Answers	Reject	Mark
<b>19(c)(ii)</b>	(primary or secondary) alcohol <b>and</b> ketone  <b>NOTE</b> <b>BOTH names</b> are required here	Just 'hydroxyl for 'alcohol' and/or 'C=O /carbonyl' for ketone/	<b>1</b>

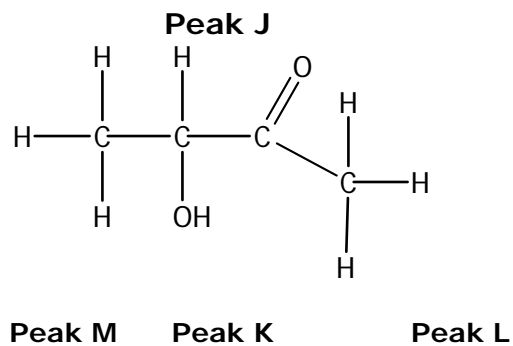
Question Number	Acceptable Answers	Reject	Mark
19(d)	<p><b>MARKING ADVICE</b> Check answer for the suggested structure of X. If the correct structure is shown</p>  <p>Mark answer according to the following. However if no structure for <b>X</b> is shown or an incorrect structure for <b>X</b> is proposed, mark answer according to "COMMENTS" scheme below</p> <p><b>MARKS CAN BE AWARDED FROM SUITABLY ANNOTATED FORMULAE FOR X.</b></p> <p><b>First mark:</b></p> <p>Four different H / hydrogen / proton environments (1)</p> <p><b>Any five from following seven points:</b></p> <p><b>Either</b> Application of the (n+1) rule to peak <b>J</b> (which is a quartet / splits into <b>four</b>) <b>or</b> application of the (n+1) rule peak <b>M</b> (which is a doublet / splits into <b>two</b>) (1)</p> <p>Any mention to explain <b>no</b> splitting for peak <b>L</b> as there is no H is attached to the adjacent carbon (1)</p> <p>Peak <b>L</b> (<b>CH</b><sub>3</sub>) next to C=O (1)</p> <p>Peak <b>M</b> (<b>CH</b><sub>3</sub>) next to CH (1)</p> <p>Peak <b>K</b> <b>OH</b> (1)</p> <p>Peak <b>J</b> (<b>CH</b>) next to CH<sub>3</sub> (1)</p> <p>Any <b>one</b> correct <math>\delta</math> value quoted within <math>\pm 0.2</math> of the following chemical shifts: 1.4(<b>M</b>) or 2.2 (<b>L</b>) or 3.7(<b>K</b>) or 4.2 (<b>J</b>) (ppm) (1)</p>	<p><b>Just 'four different chemical environments'</b></p> <p>If any incorrect chemical shift <b>OR A RANGE</b> of chemical shifts is quoted, this scoring point is not available</p>	7

**Final mark**

(Compound **X** is)  $\text{CH}_3\text{CH}(\text{OH})\text{COCH}_3$   
**NO** other compound allowed.

**ACCEPT**

any unambiguous formula, e.g. displayed formula



**ACCEPT**

3-hydroxybutan-2-one

**(1)**