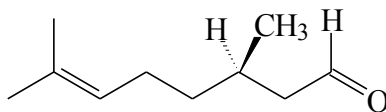


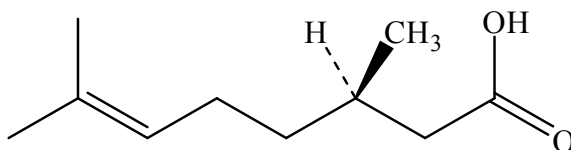
1.



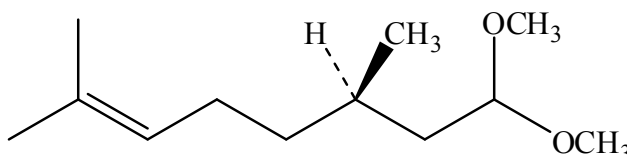
(a) **C₁₀H₁₈O**

(b) **Aldehyde (-CHO, carbonyl) and alkene (C=C)**

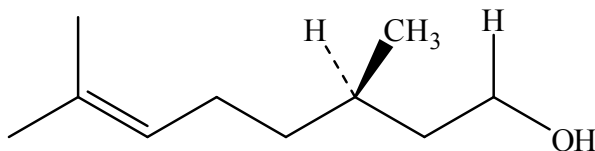
(c) (i) **Cr₂O₇²⁻ / H⁺: aldehyde → carboxylic acid:**



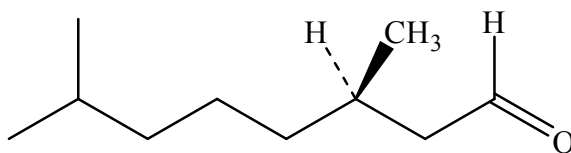
(ii) **excess CH₃OH / catalytic amount H₂SO₄: aldehyde → acetal:**



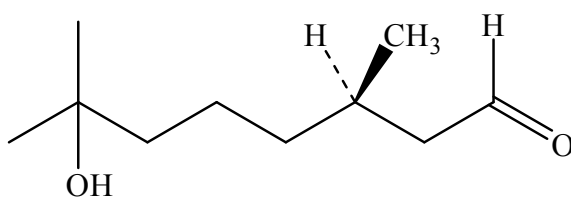
(iii) **NaBH₄ in CH₃OH followed by H⁺/H₂O: aldehyde → primary alcohol:**



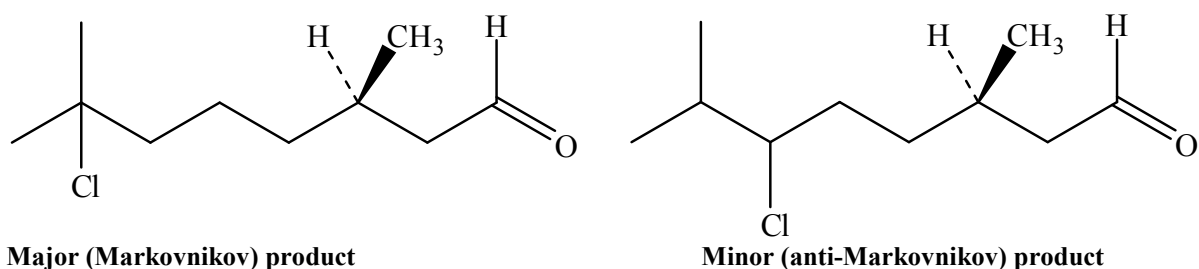
(iv) **H₂/Pd in ethanol: alkene → alkane:**



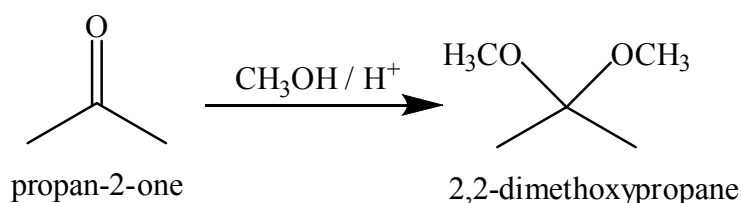
(v) **3 M H₂SO₄: alkene → alcohol:**



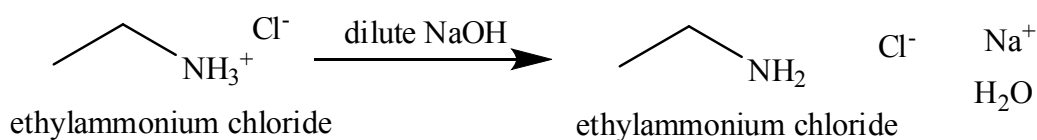
(vi) HCl in CCl₄ solvent: **addition of HCl across C=C:**



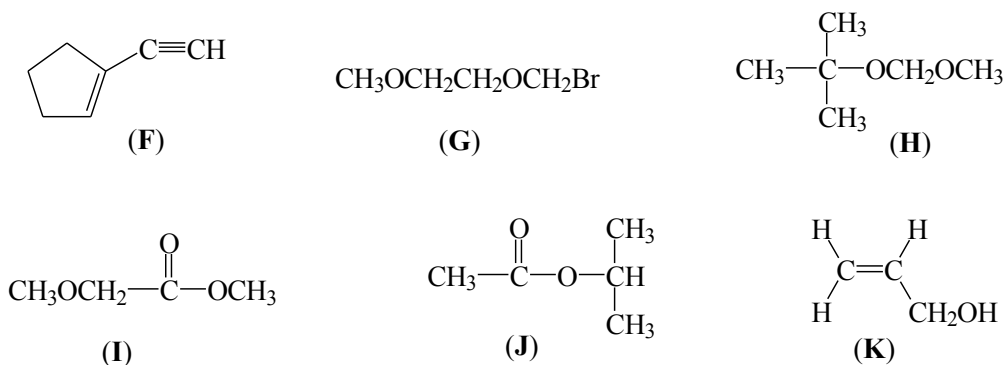
2. (a) **Acid catalysed addition of CH₃OH to ketone → acetal:**



(b) **Deprotonation:**



3.



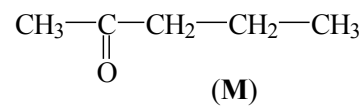
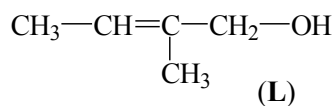
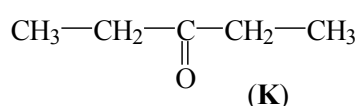
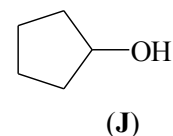
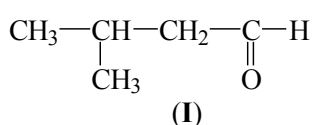
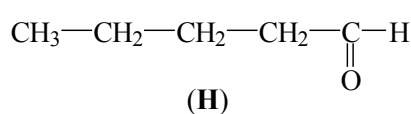
- (a) **IR absorption in the range 1650-1800 cm⁻¹ is associated with the presence of the C=O group: I and J**
- (b) **Strong IR absorption in the range 3200-3700 cm⁻¹ is associated with the presence of the O-H and N-H groups: K only**
- (c) **Strong UV absorption is associated with the presence of conjugated systems (i.e. those with alternating single and multiple bonds): F only**
- (d) **The presence of only three signals in the ¹H NMR means that three (and only three) distinct types of hydrogen – H, I and J.**

As these are singlets, there are *not* chemically different hydrogen atoms on immediately adjacent carbon atoms. This is not the case for J where the signal on hydrogen of the CH group will be split into a septet by the 2CH₃ groups (and the signal due to hydrogen atoms on these CH₃ will be split into a doublet by the CH group).

Overall, only H and I fit.

- (e) Bromine has two main isotopes ⁷⁹Br (51%) and ⁸¹Br (49%). The mass spectrum of G therefore shows peaks with approximately equal intensity at separated by 2 mass units.

4.



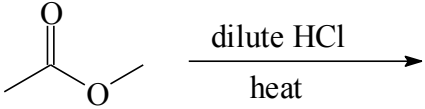
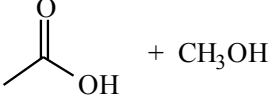
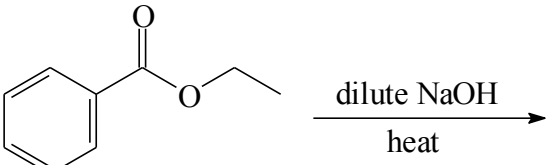
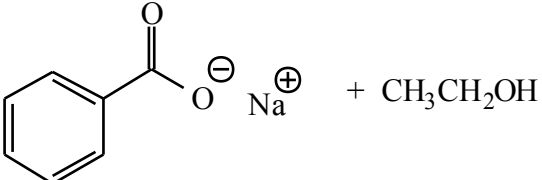
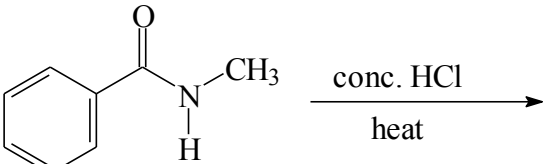
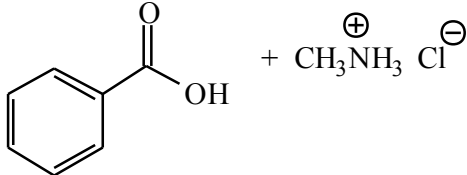
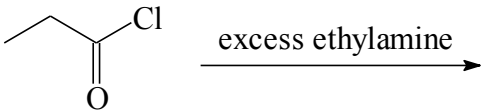
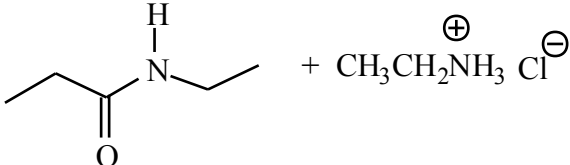
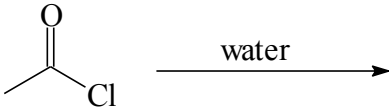

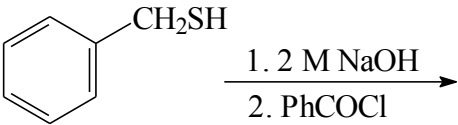
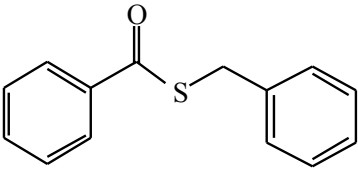
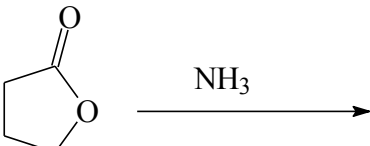
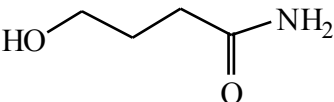
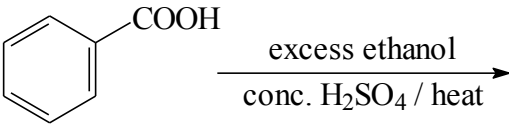
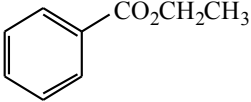
- (a) All of them have molar mass = 86.
- (b) The number shows two signals in the ¹H NMR spectrum – a quartet ('a') and a singlet ('b'). The molecule therefore has only two types of hydrogen atom. This is only true for K.
- (c) The signal due to the hydrogen atoms on the CH₂ group is split into a quartet by the CH₃ group. Hence signal 'a' is due to the two hydrogen atoms on the CH₂ group. The relative intensity is therefore 2.
- The signal due to the hydrogen atoms on the CH₃ group is split into a triplet by the CH₂ group. Hence signal 'b' is due to the three hydrogen atoms on the CH₃ group. The relative intensity is therefore 3.
- (d) IR absorption in the ranges 1650-1750, 2850-2950 and 3200-3600 cm⁻¹ are associated with vibrations of C=O, C-H and O-H groups respectively.

1650-1750 cm⁻¹: H, I, K and M

2850-2950 cm⁻¹: all of them

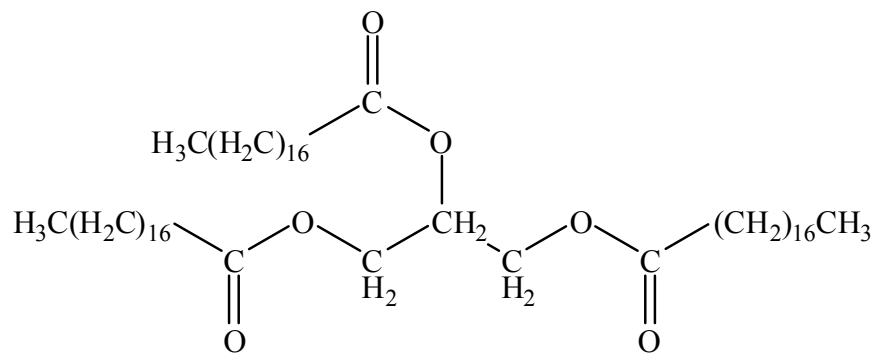
3200-3600 cm⁻¹: J and L.

5.

reactants and reaction conditions	major organic product(s)
 <chem>CC(=O)OC >> CC(=O)O + CO</chem>	 <chem>CC(=O)O + CO</chem>
 <chem>CCOC(=O)c1ccccc1 >> CCOC(=O)[O-] + [Na+]</chem>	 <chem>CCOC(=O)[O-] + [Na+] + CCO</chem>
 <chem>CNC(=O)c1ccccc1 >> OC(=O)c1ccccc1 + CN</chem>	 <chem>OC(=O)c1ccccc1 + CN + [Cl-]</chem>
 <chem>CCC(=O)Cl >> CCC(=O)NCC + CCN</chem>	 <chem>CCC(=O)NCC + CCN + [Cl-]</chem>
 <chem>CC(=O)Cl >> CC(=O)O</chem>	 <chem>CC(=O)O</chem>
 <chem>c1ccccc1CS >> c1ccccc1C(=O)SCc2ccccc2</chem>	 <chem>c1ccccc1C(=O)SCc2ccccc2</chem>
 <chem>O=C1OCCC1 >> NCC(=O)CCO</chem>	 <chem>NCC(=O)CCO</chem>
 <chem>c1ccccc1C(=O)O >> c1ccccc1C(=O)OCC</chem>	 <chem>c1ccccc1C(=O)OCC</chem>

6. **Fats and oils are triesters of glycerol:**

glycerol + 3 fatty acid → triglyceride



The triglyceride can be converted back to stearic acid and glycerol using either strong aqueous acid and heat or strong aqueous base and heat.