

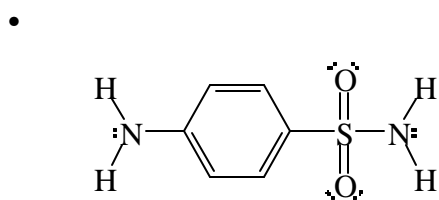
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- $4.29 \times 10^{14} \text{ Hz} - 6.67 \times 10^{14} \text{ Hz}$
 $2.84 \times 10^{-19} \text{ J} - 4.42 \times 10^{-19} \text{ J}$ ($171 \text{ kJ mol}^{-1} - 266 \text{ kJ mol}^{-1}$)
- Electron density is the probability of finding an electron at that point in space.
- F: The $2p$ electrons are not effective in screening other $2p$ electrons, so the electrons in F feel a higher effective nuclear charge in F than in O.
 O: The $2p$ electrons in O are closer to the nucleus and therefore held more tightly than the $3p$ electrons in S.
- To get strong double and triple bonds the nuclei need to be close together, since the orbital overlap in a π -type bond is not as good as in a σ -type bond. Hence only small atoms (mainly second row) can form multiple bonds.

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- $\ominus \quad \oplus$
 $:\text{C} \equiv \text{O}:$ Electronegativity values suggest that the O would be δ^- and the C would be δ^+ . The formal charges, however, counteract the electronegativity differences and a small molecular dipole results.
- 1,4-Butanediol will be more soluble in water as it can form more hydrogen bonds than 1-butanol.
 Chloroform will be more soluble as it is more polar than carbon tetrachloride. Being more polar, it will have stronger dipole-dipole interactions with the very polar solvent.

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tetrahedral

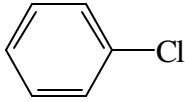
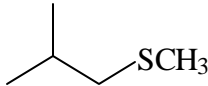
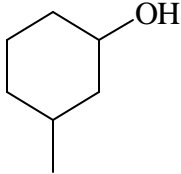
sp^2

120°

10

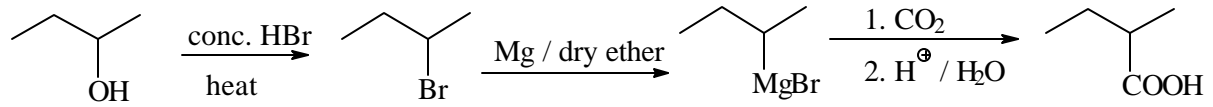
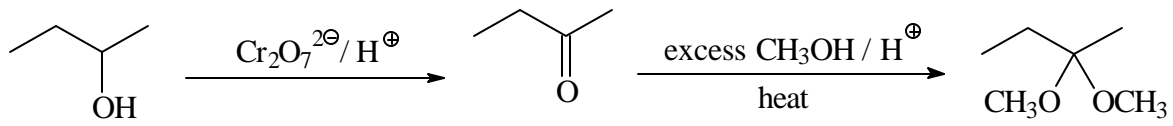
- Hydrogen bonding: O is small enough and electronegative enough to form H-bonds.
 Dipole-dipole: The H_2S molecule has a dipole - S is too big to form H-bonds, yet not big enough to have significant dispersion forces.
 Dispersion: Te is less electronegative than S, so the dipole in H_2Te is less than that in H_2S . However, the b.p. of H_2Te is greater than that of H_2S , so the dominant force must be dispersion forces - Te is much larger atom than S and therefore more polarisable.

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	1) CH_3MgI 2) $\text{H}^+ / \text{H}_2\text{O}$	2-butanol
		$\text{CH}_3\text{—CH(OH)—CH}_2\text{CH}_3$
		
3-methylcyclohexanone		
1-propyl acetate		CH_3COOH + $\text{HOCH}_2\text{CH}_2\text{CH}_3$
		$\text{CH}_3\text{—C(=O)—N—CH}_2\text{CH}_3$ H + $\text{CH}_3\text{CO}_2^- \text{NH}_3^+\text{CH}_2\text{CH}_3$
	conc. H_2SO_4 / heat	

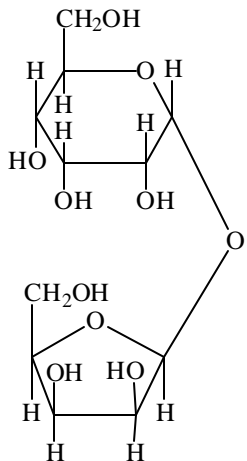
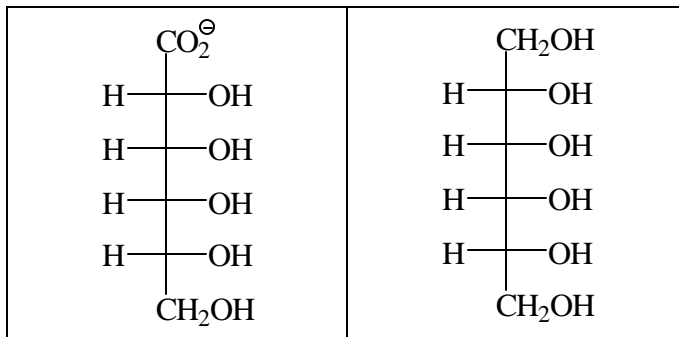
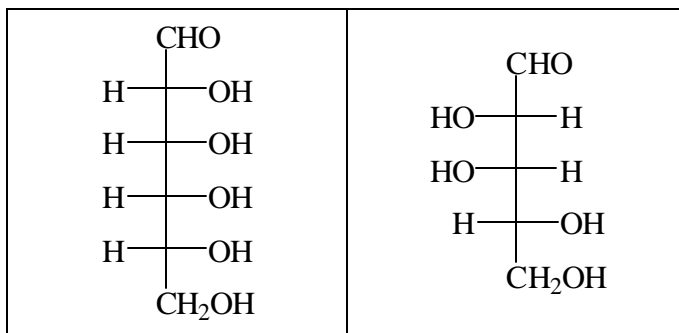
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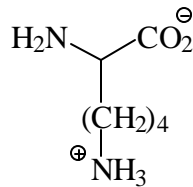
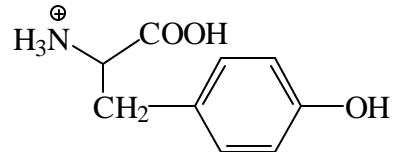
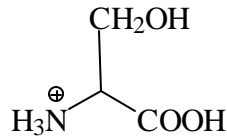
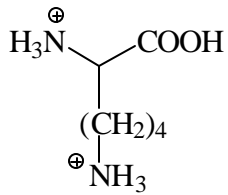
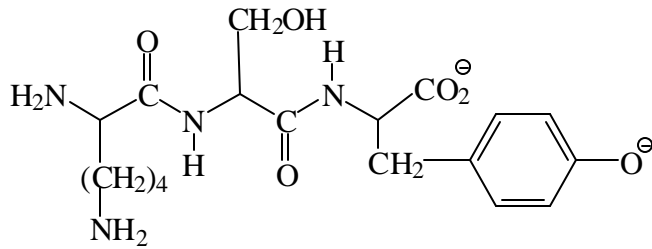
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pH = 9.74

