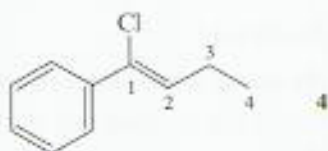


Figure 2 Two layers of close-packed spheres.

Question 4

Ignoring the ring carbon atoms, identify the longest chain in Compound 4 as C₄ and therefore derived from butane. Numbering the carbon chain from the end attached to the benzene ring gives



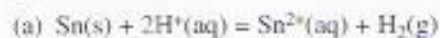
Name the chloro and phenyl groups, and put them in the correct alphabetical order.

Assign both of these substituents to C-1, and identify the double bond as 'ene'. (Note The double bond need not be labelled 1 (as in but-1-ene).)

The priorities on C-1 are Cl > phenyl; and on C-2 are ethyl > hydrogen. The two higher-priority groups, Cl and CH₂CH₃, are on the *same* side of the double bond, so Compound 4 is the *Z*-isomer.

The systematic name of Compound 4 is therefore *Z*-1-chloro-1-phenylbutene.

Question 5



(b) $\Delta S_m^\ominus = (-17.0 + 130.7 - 51.6) \text{ J K}^{-1} \text{ mol}^{-1} = 62.1 \text{ J K}^{-1} \text{ mol}^{-1}$

$$\Delta H_m^\ominus = -8.8 \text{ kJ mol}^{-1}$$

So $\Delta G_m^\ominus = \Delta G_f^\ominus = \Delta H_m^\ominus - T\Delta S_m^\ominus = -8.8 \text{ kJ mol}^{-1} - (298.15 \times 62.1) \text{ J mol}^{-1}$

$$= -8.8 \text{ kJ mol}^{-1} - 18\,515 \text{ J mol}^{-1}$$

$$= -8.8 \text{ kJ mol}^{-1} - 18.5 \text{ kJ mol}^{-1} = -27.3 \text{ kJ mol}^{-1}$$

Question 6

Molecularity provides a means of classifying elementary reactions in terms of the number of reactant species that take part in the reaction.

If Reaction 1 were elementary, the theoretical rate equation would be

$$J = k_{\text{theory}}[\text{F}_2][\text{CCl}_3\text{F}]$$

However, this is not the case, since the partial order with respect to F₂ in the given experimental rate equation is fractional (1/2).

The reaction must therefore be composite.