

for each two carbon atom $\rightarrow l_{C-C} = 0.154 \times 10^{-9} \times \sin\left(\frac{109.5}{2}\right) = 0.126 \text{ nm}$

for C-H bonds $\rightarrow l_{C-H} = 0.107 \times 10^{-9} \times \sin\left(\frac{109.5}{2}\right) = 0.087 \text{ nm}$

$$L(n_c) = 0.174 + (n_c - 1) \times 0.126 \text{ nm}$$

b) for $n_c = 14 \rightarrow L = 1.812 \text{ nm}$

for $n_c = 20 \rightarrow L = 2.568 \text{ nm}$

7.2) entropy per molecule for a gas (ideal) is:

$$S_{\text{gas}} = K_B \left\{ \frac{5}{2} - \ln \left(\rho \cdot \left[\frac{h}{(2\pi m K_B T)^{1/2}} \right]^3 \right) \right\}$$

where for an ideal hydrogen gas: $PV = \frac{NRT}{N_A} \rightarrow PN_A = PRT \rightarrow \rho = \frac{PN_A}{RT}$

So $\rightarrow S_{\text{gas}} = K_B \left\{ \frac{5}{2} - \ln \left(\frac{PN_A}{RT} \cdot \left(\frac{h}{(2\pi m K_B T)^{1/2}} \right)^3 \right) \right\}$

$P = 10^5 \text{ Pa}$, $N_A = 6.022 \times 10^{23}$, $T = 273 \text{ K}$, $R = 8.31 \frac{\text{J}}{\text{mol} \cdot \text{K}}$, $m = 3.32 \times 10^{-27} \text{ kg}$ hydrogen molecule

$K_B = 1.38 \times 10^{-23} \frac{\text{m}^2 \text{kg}}{\text{s}^2 \text{K}}$, $h = 6.62 \times 10^{-34} \frac{\text{J}}{\text{Hz}}$

$\rightarrow \frac{S_{\text{gas}}}{K_B} = \frac{5}{2} - \ln \left(\right) = 13.91 \gg \frac{PV}{K_B T} = 1 \text{ (per molecule } N=1)$

$$\ln \frac{CMC}{690} = -1.15 n_c$$

(7.3) according to Figure 7.9, for single chain lipids the CMC is found from function:

$$CMC \text{ (molar)} = 690 \times e^{-1.15 n_c} \text{ molar} \rightarrow n_c = -\frac{1}{1.15} \ln \frac{CMC}{690}$$

$$\text{for lauroyl pc} \rightarrow n_c = -\frac{1}{1.15} \times \ln \frac{7 \times 10^{-4}}{690} = 12$$

$$\text{for myristoyl pc} \rightarrow n_c = 14$$

$$\text{for palmitoyl pc} \rightarrow n_c = 16$$

$$\text{for } T = 300K, m = 500 \text{ Da} \rightarrow \gamma = ?$$

$$\text{eq 7.6 } \rho_{agg} = \left(\frac{(2\pi m k_B T)^{3/2}}{h} \right)^3 e^{5/2} \exp \left(\frac{-2\pi n_c R_{hc} l_{cc} \gamma}{k_B T} \right)$$

$$7.4) \quad V_{hc=0.2} \text{ [nm]} \quad \text{for a single hydrocarbon chain}$$

the a_0 expected for lipid bilayer formation with 1,2,3 chains is found from

$$\frac{1}{2} \leq \frac{V_{hc}}{a_0 l_{hc}} \leq 1$$

$$\text{for single chain: } \frac{1}{2} \leq \frac{0.2}{a_0} \leq 1 \rightarrow 0.2 \text{ nm}^2 \leq a_0 \leq 0.4 \text{ nm}^2$$

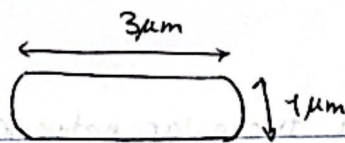
$$\text{double chain: } \frac{1}{2} \leq \frac{0.4}{a_0} \leq 1 \rightarrow 0.4 \text{ nm}^2 \leq a_0 \leq 0.8 \text{ nm}^2$$

$$\text{triple chain: } \frac{1}{2} \leq \frac{0.6}{a_0} \leq 1 \rightarrow 0.6 \text{ nm}^2 \leq a_0 \leq 1.2 \text{ nm}^2$$

the results are slightly smaller than those of Koenig but still close enough to be reasonable for a triple chain.

$$\begin{matrix} 0.5 a_0 < 0.12 \\ 0.12 < a_0 \end{matrix}$$

7.5) $k_b = k_G = 10 k_B T$



The spherocylinder is constructed from two half spheres and a cylinder with $L = 2 \mu m$ & $R = 0.5 \mu m$

$$E_{\text{bending}} = E_{\text{spheres}} + E_{\text{cylinder}} = 4\pi(2k_b + k_G) + \frac{\pi L}{R} k_b$$

$$= (120\pi + 40\pi) k_B T = 160\pi k_B T$$

Bending energy for daughter cells:

$$E_{\text{bending}} = 2 \times \left((2k_b + k_G) \times 4\pi + \frac{\pi L}{R} k_b \right) \rightarrow L = ? \quad R = ?$$

and diameter

Since the total volume remains constant:

$$V_i = (2 \mu m \times \pi R^2) + \left(\frac{4}{3} \pi R^3 \right) = 2 \times \left[(L' \times \pi R^2) + \left(\frac{4}{3} \pi R^3 \right) \right]$$

$$2\pi R^2 + \frac{4}{3}\pi R^3 = 2L'\pi R^2 + \frac{8}{3}\pi R^3 \xrightarrow{R=0.5} L' = \frac{2\pi R^2 - \frac{8}{3}\pi R^3}{2\pi R^2} = \frac{4}{3} \mu m$$

$$E_{\text{bend}} = 2 \times \left(120\pi + \frac{80\pi}{3} \right) k_B T = \frac{880\pi}{3} k_B T$$

7.6) in a triangular network (6-fold symmetry):

$$K_A = \left(\frac{\sqrt{3}}{2} K_{sp} \right) \cdot \text{zero temperature \& zero stress} \quad \mu = \frac{\sqrt{3}}{4} K_{sp}$$

$$\mu = 10^{-5} \frac{\mu}{m^2} \rightarrow K_{sp} = \frac{4}{\sqrt{3}} \mu \rightarrow K_A = 2 K_{sp} \mu = 2 \times 10^{-5} \frac{\mu}{m^2}$$

$$K_b \rightarrow = \frac{2 \times 10^{-5} \times (25 \times 10^{-9})^2}{12} = 1.04 \times 10^{-21} = 0.25 K_B T$$

$$K_b \text{ dampc} \sim 5.33 K_B T$$

the structure helps with
the flexibility of the RBC