



UNITED NATIONS EDUCATIONAL, SCIENTIFIC AND CULTURAL ORGANIZATION
INTERNATIONAL ATOMIC ENERGY AGENCY
INTERNATIONAL CENTRE FOR THEORETICAL PHYSICS
I.C.T.P., P.O. BOX 586, 34100 TRIESTE, ITALY, CABLE: CENTRATOM TRIESTE



SMR.961 - 3

**WORKSHOP ON:
PROTEINS, MEMBRANES and their INTERACTIONS**

22 JULY - 2 AUGUST 1996

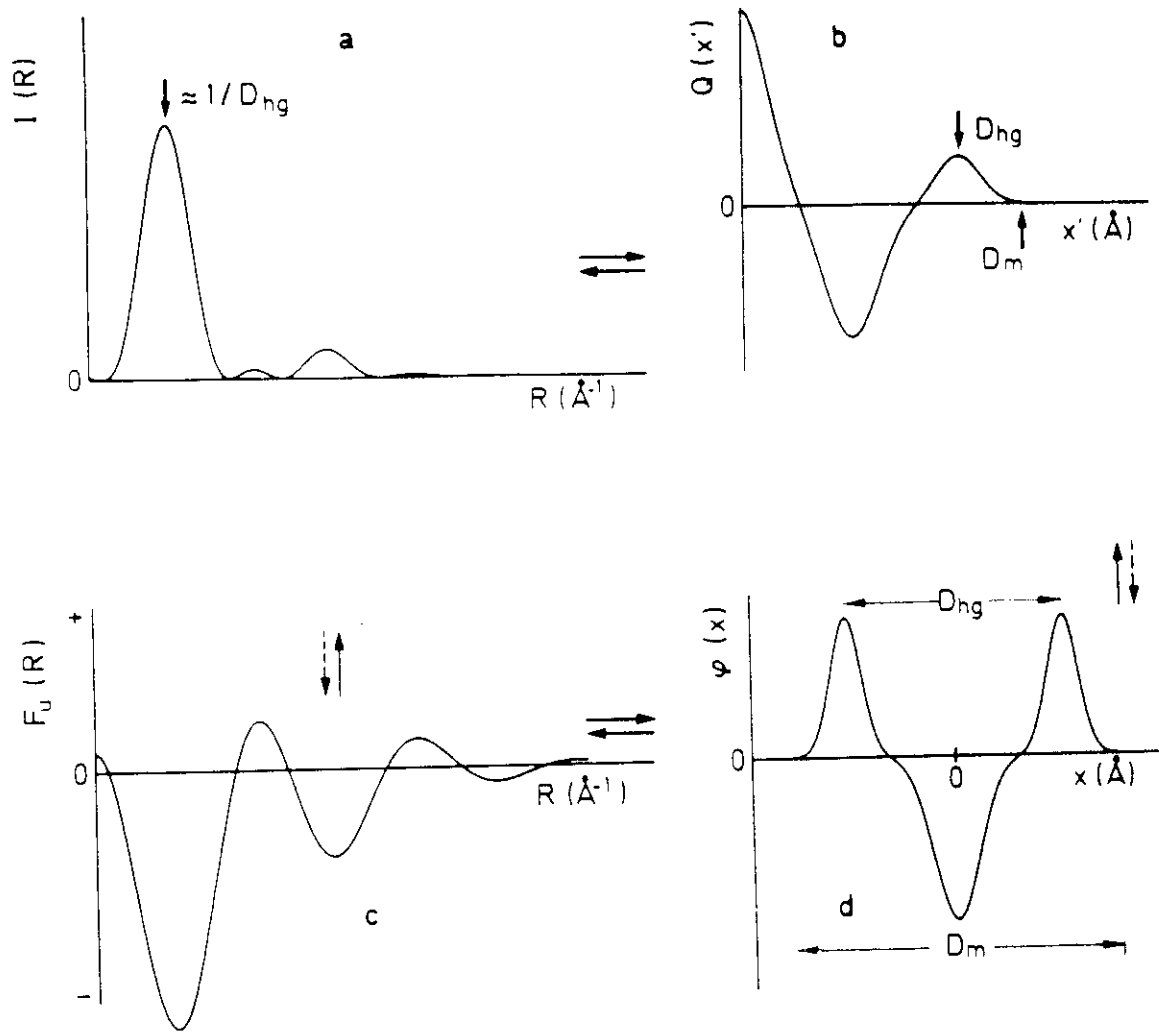
**"Lipid polymorphism and
Membrane Structure and Function"**

PART II

**John SEDDON
Imperial College of Science, Technology and Medicine
Department of Chemistry
Exhibition Road, South Kensington
SW7 2AY London
U.K.**

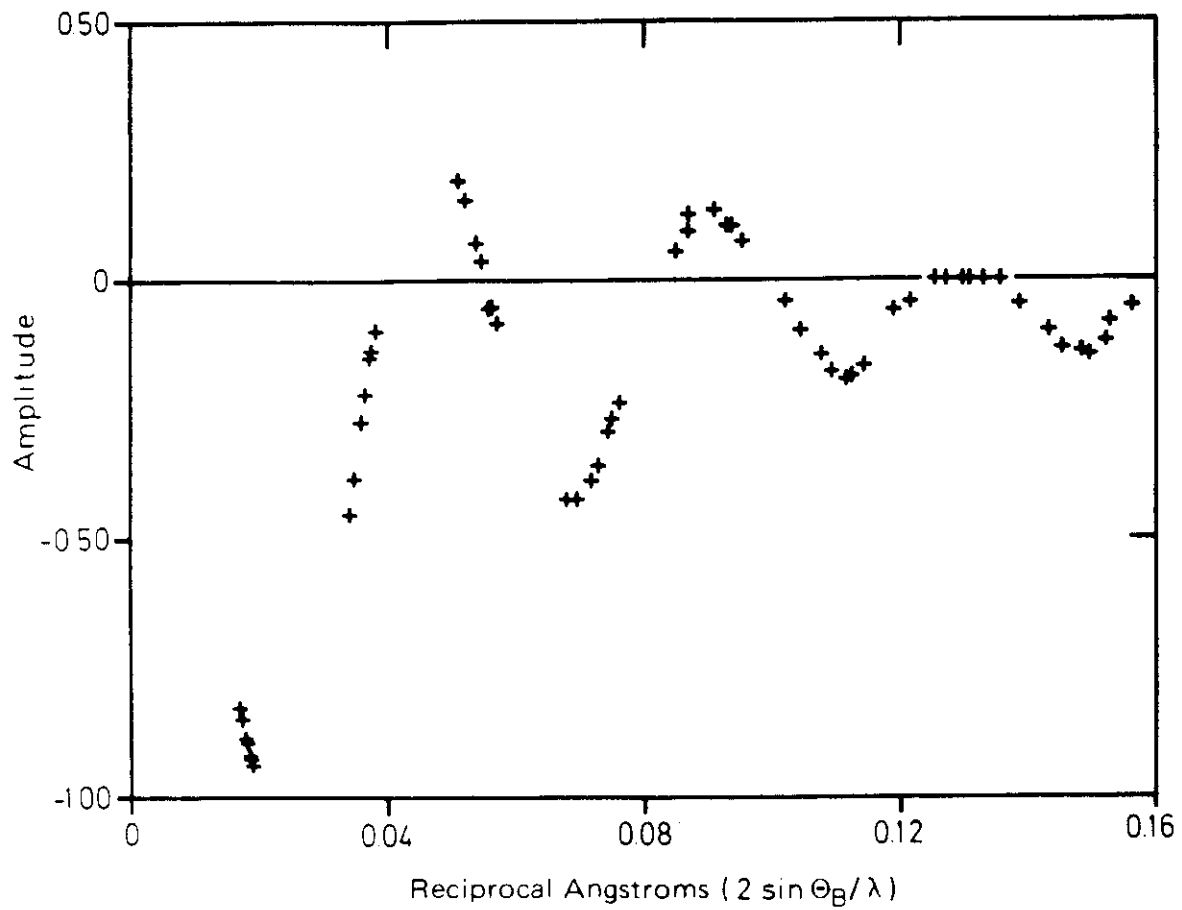
These are preliminary lecture notes, intended only for distribution to participants.

Single bilayer diffraction



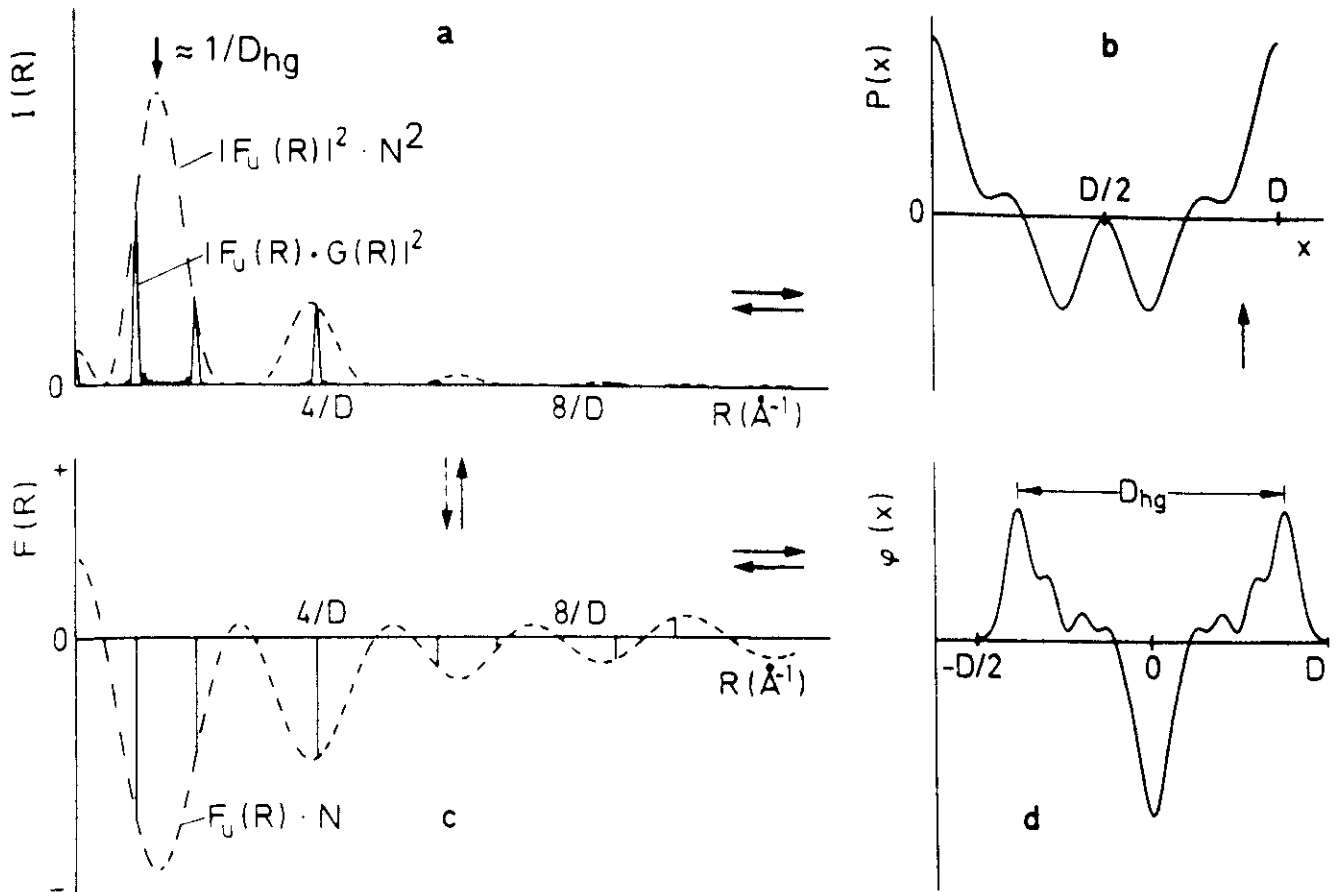
From: N. P. Franks and Y. Levine, in: *Membrane Spectroscopy* (E. Grell, Ed.), 1981, Springer Verlag, Berlin.

Lamellar swelling series



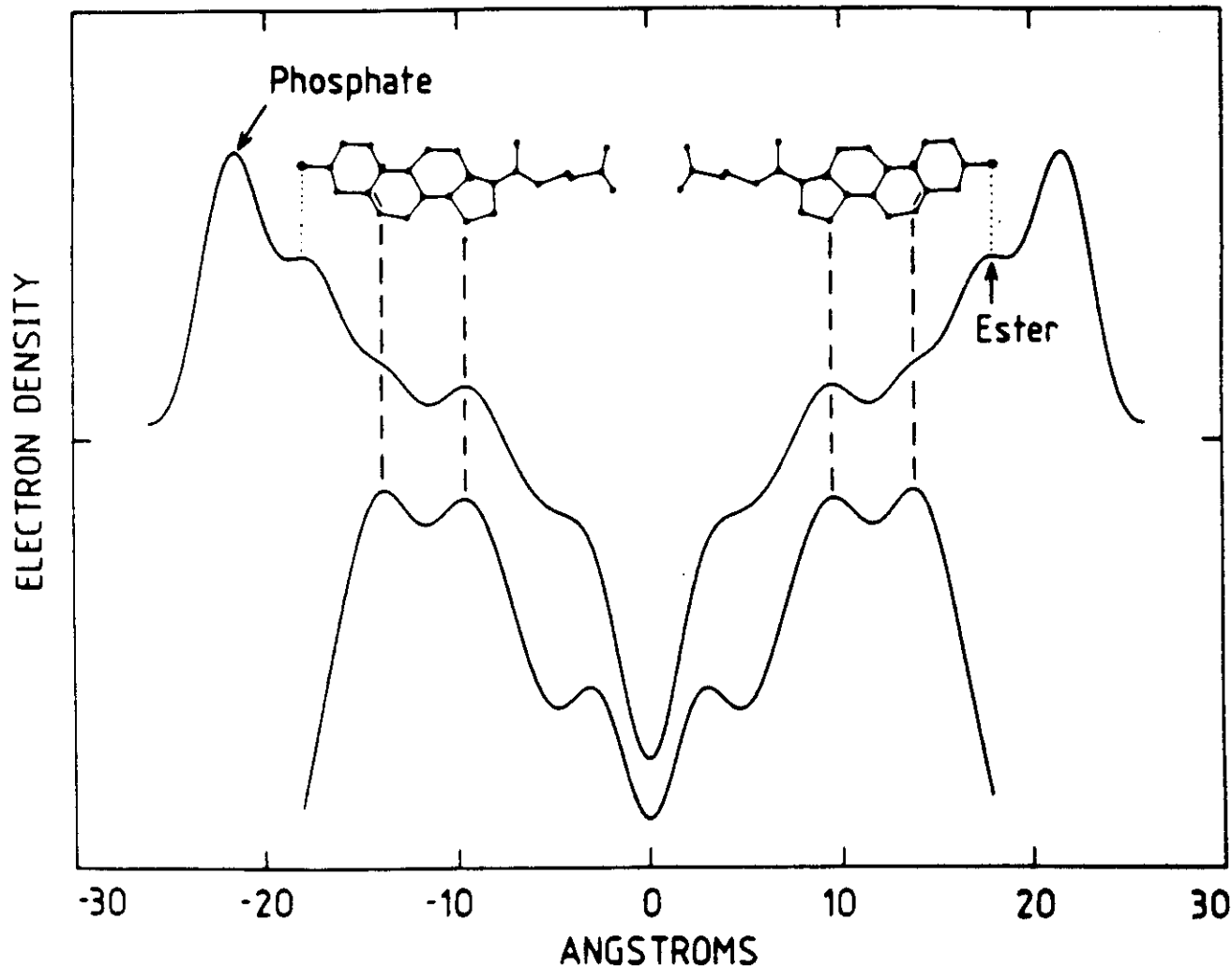
From: N. P. Franks and Y. Levine, in: *Membrane Spectroscopy*
(E. Grell, Ed.), 1981, Springer Verlag, Berlin.

Lamellar phase diffraction



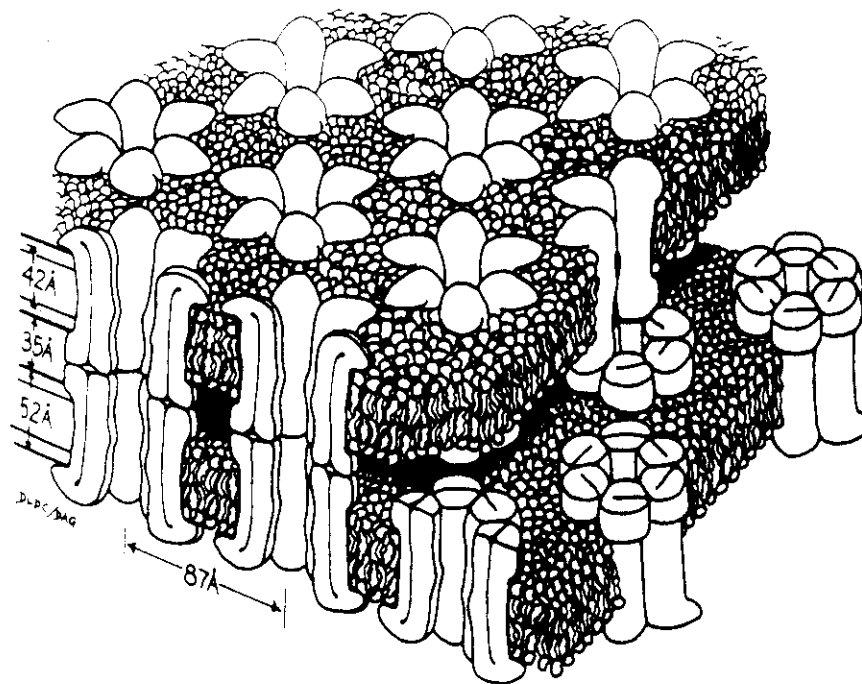
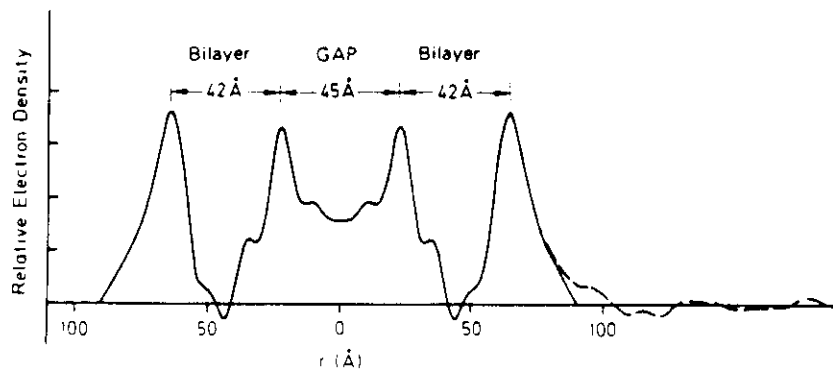
From: N. P. Franks and Y. Levine, in: **Membrane Spectroscopy**
 (E. Grell, Ed.), 1981, Springer Verlag, Berlin.

Cholesterol / DMPC bilayers: electron density profile



From: N. P. Franks and Y. Levine, in: *Membrane Spectroscopy*
(E. Grell, Ed.), 1981, Springer Verlag, Berlin.

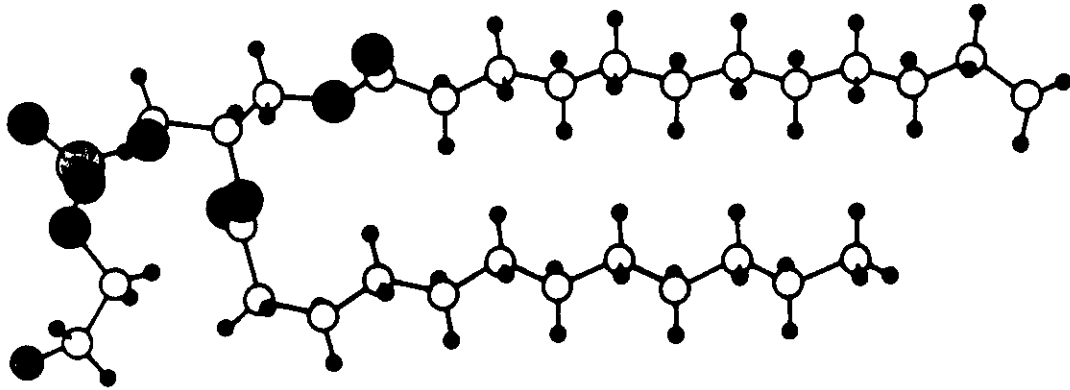
Gap junction profile



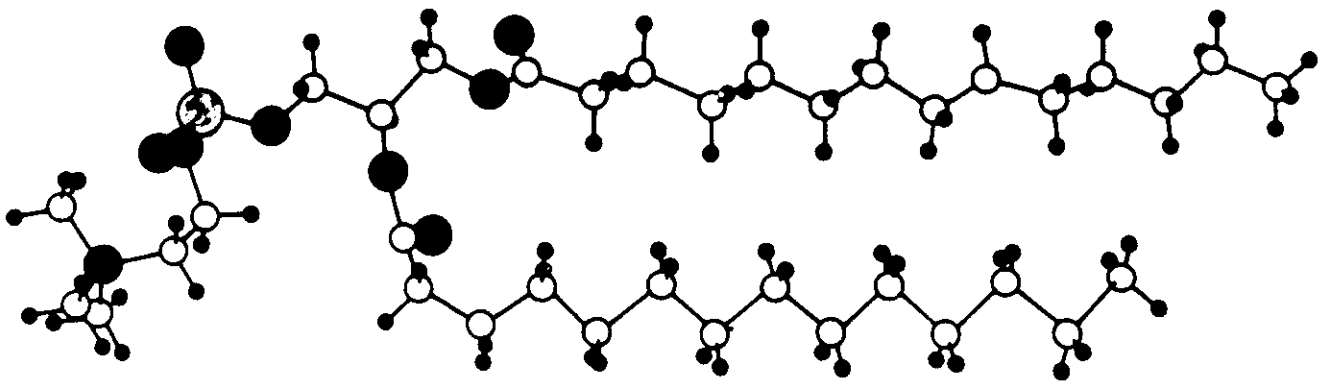
From: L. Makowski et al., 1977, *J. Cell Biol.* **74**, 629-645.

Molecular conformations of crystalline phospholipids

DLPE



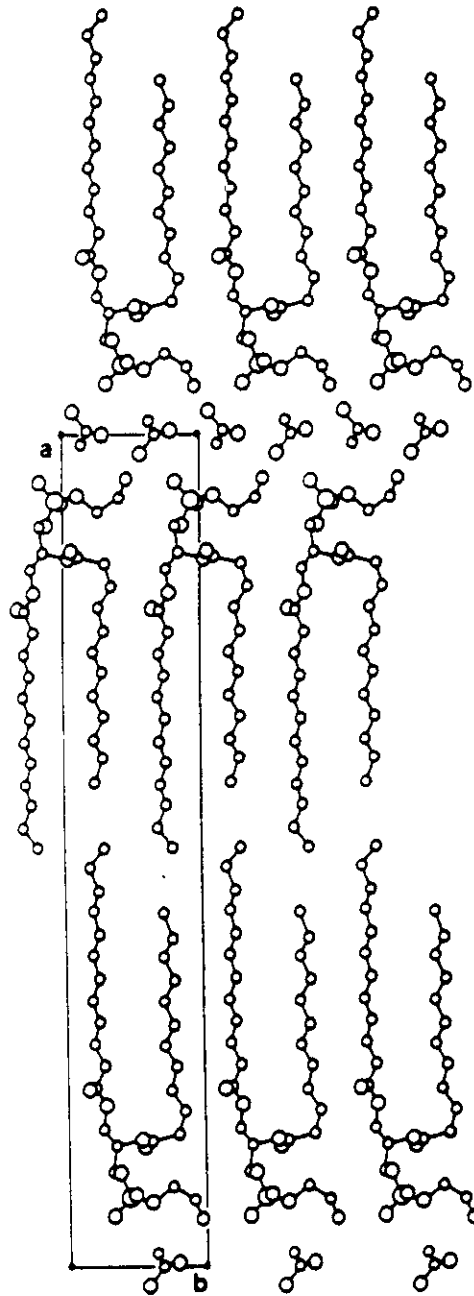
DMPC



From: I. Pascher et al., *Biochim. Biophys. Acta*, 1992, **1113**, 339-373

Packing: crystalline bilayers

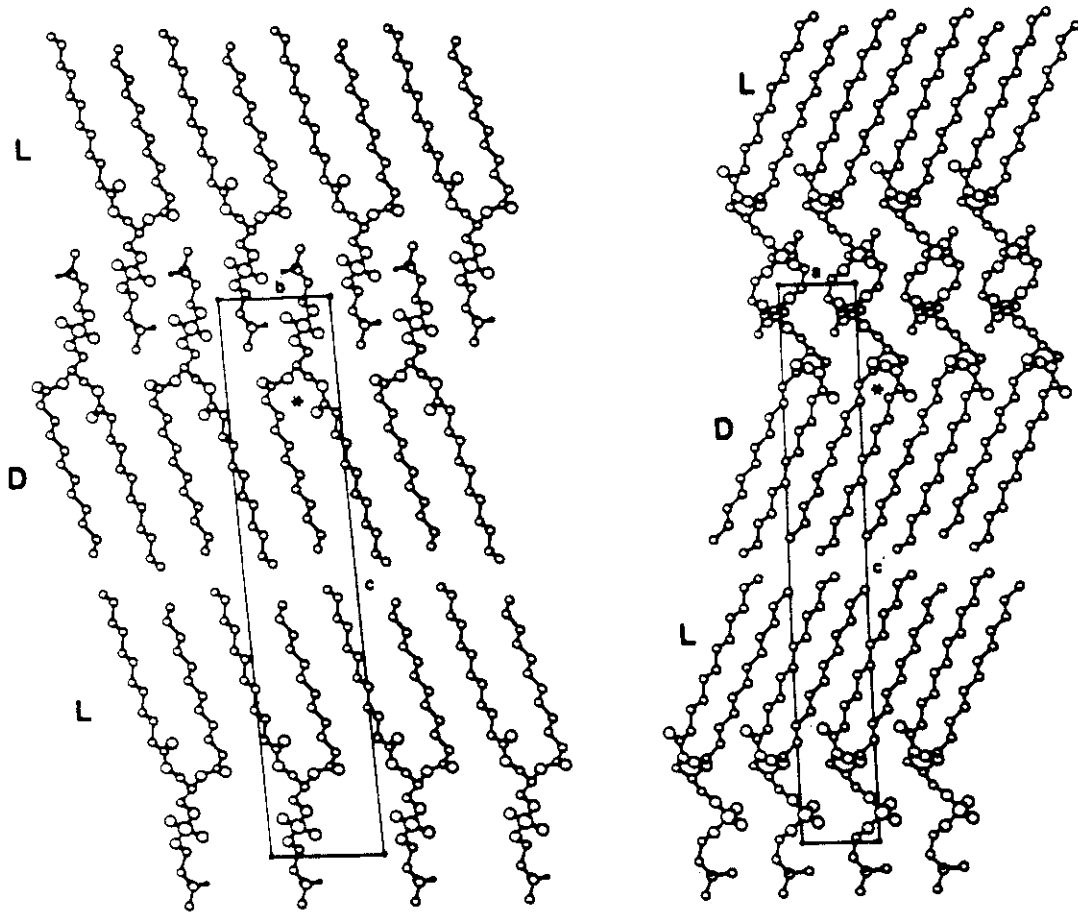
DLPE



From: I. Pascher et al., *Biochim. Biophys. Acta*, 1992, **1113**, 339-373

Packing: crystalline bilayers

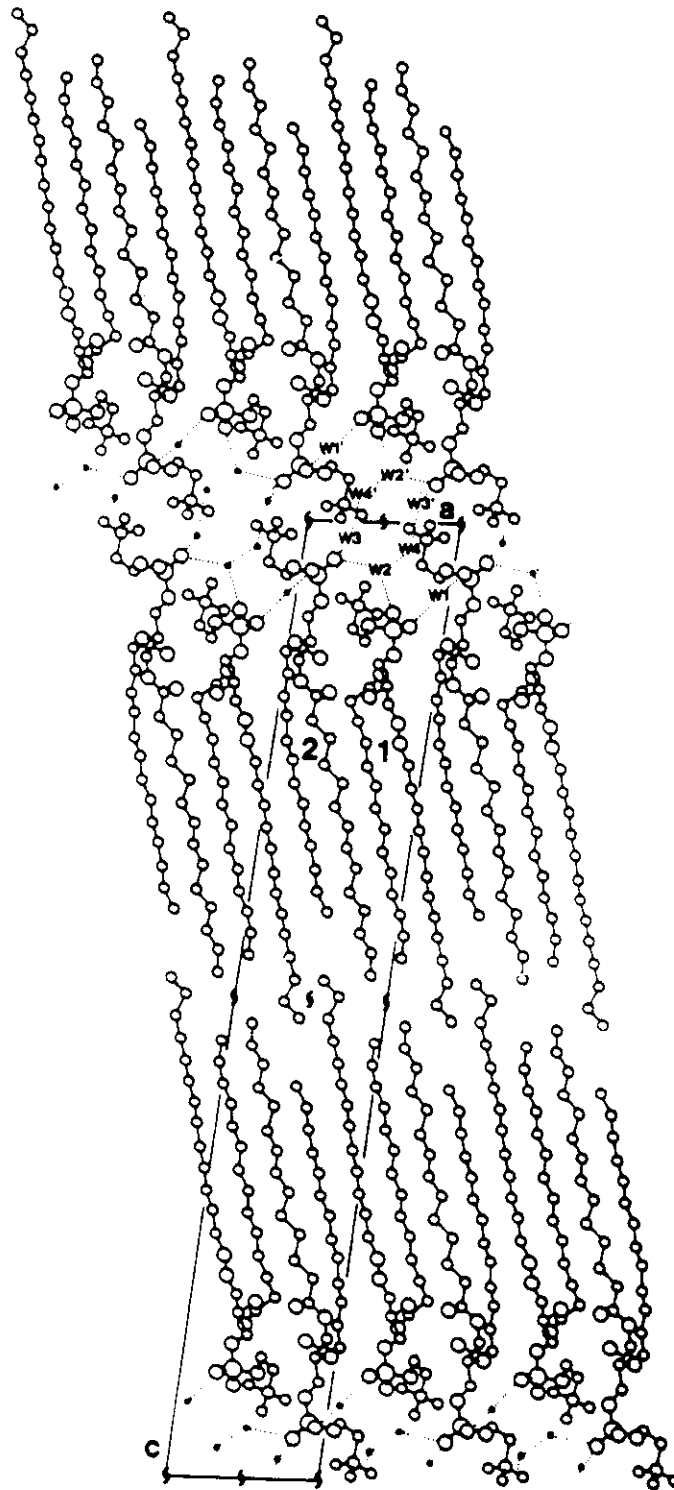
Dimethyl-DLPE



From: I. Pascher et al., *Biochim. Biophys. Acta*, 1992, **1113**, 339-373

Packing: crystalline bilayers

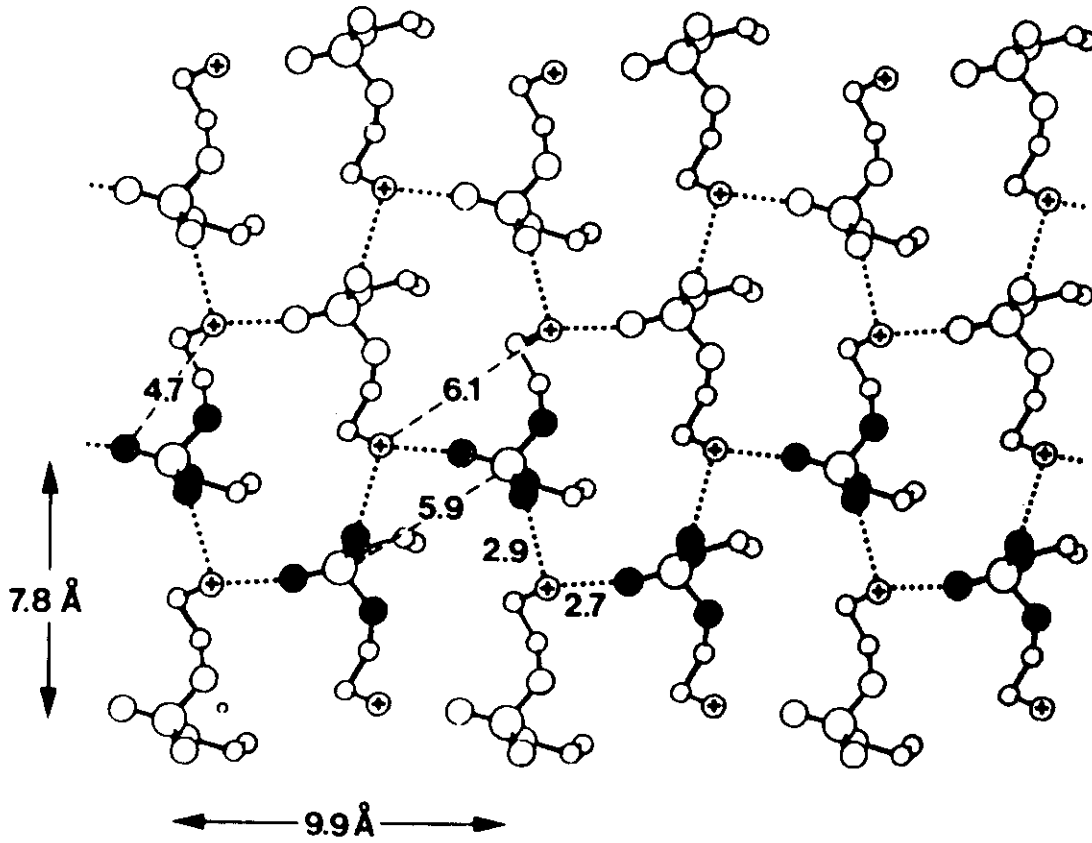
DMPC



From: I. Pascher et al., *Biochim. Biophys. Acta*, 1992, **1113**, 339-373

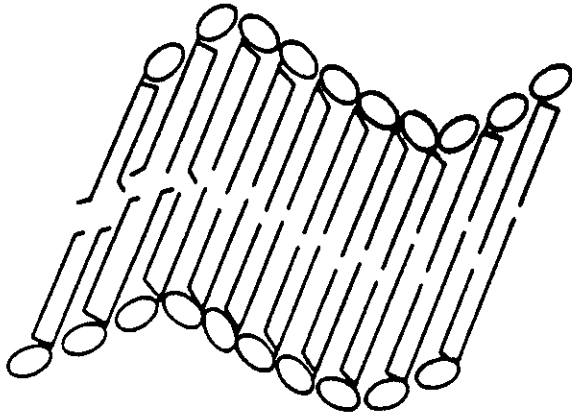
Headgroup lateral packing: crystalline bilayers

DLPE

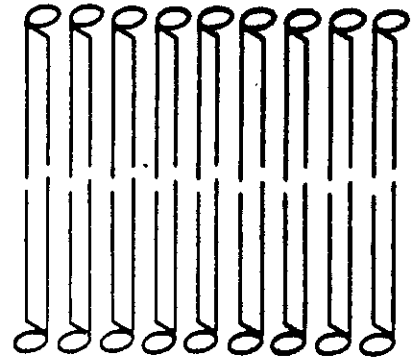


From: I. Pascher et al., *Biochim. Biophys. Acta*, 1992, **1113**, 339-373

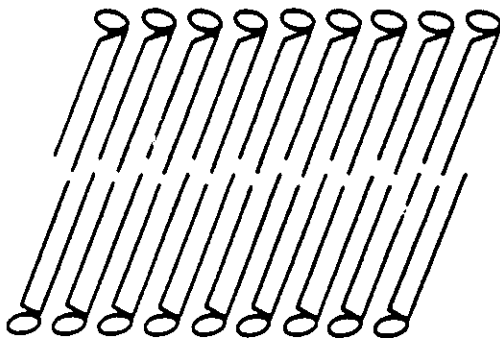
a) $P_{\beta'}$



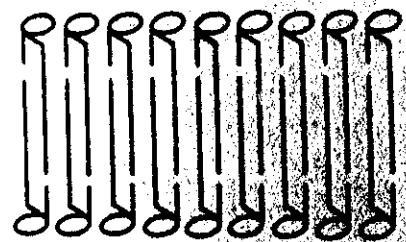
b) L_{β}



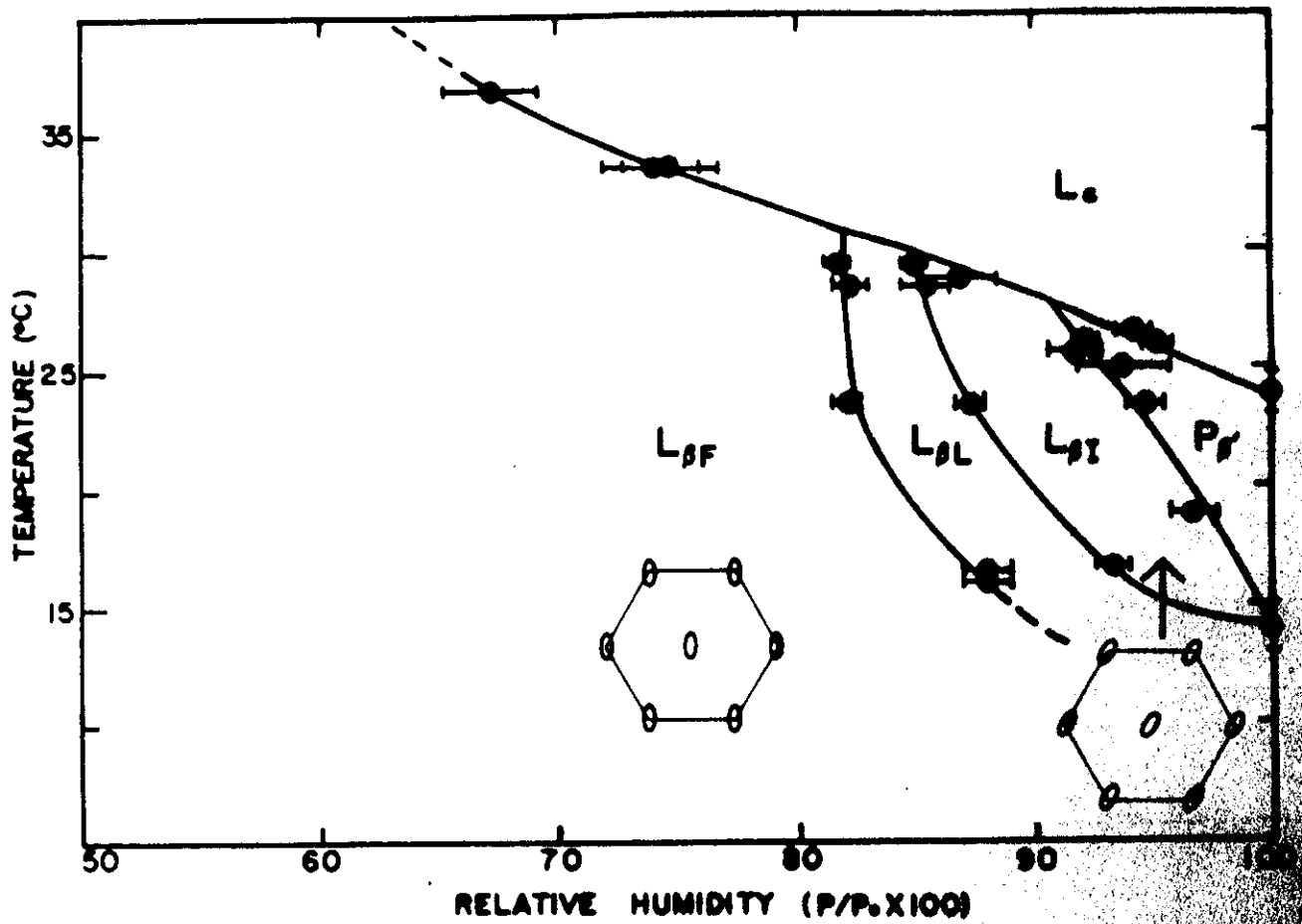
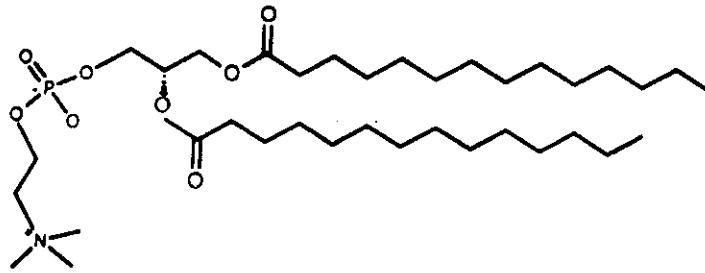
c) $L_{\beta'}$



d) $L_{\beta I}$



Dimyristoyl phosphatidylcholine / water



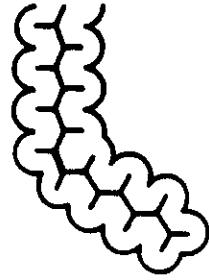
From: G. S. Smith et al. (1990) *J. Chem. Phys.* 92, 4519.

Gel-fluid transition

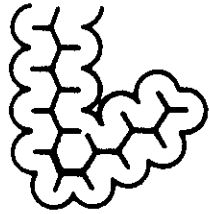
Chain rotational isomerism



all-trans



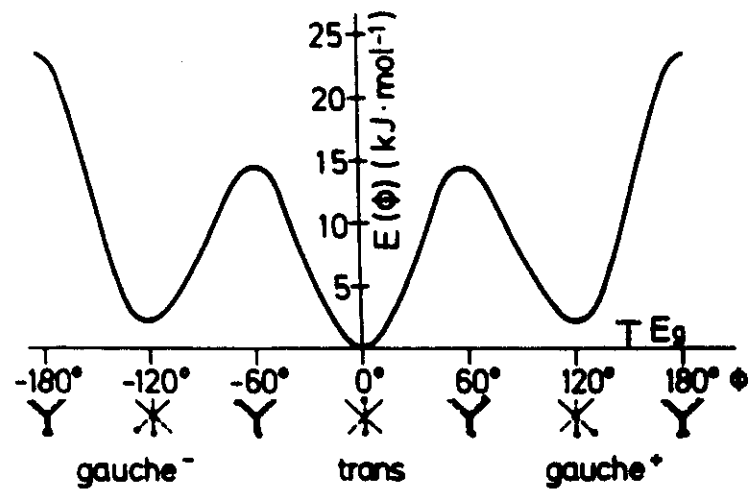
gauche (g)



$g^{\pm}g^{\mp}$

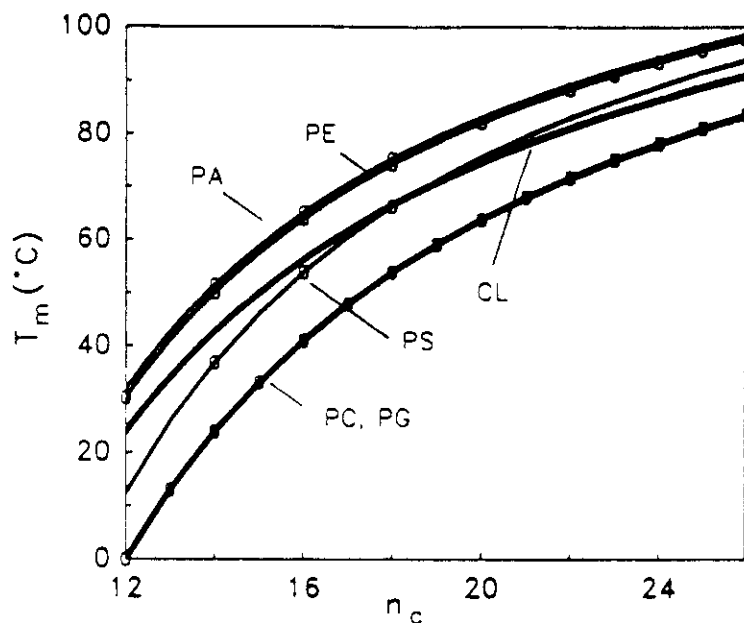


$g^{\pm}tg^{\mp}$



From: G. Ceve and D. Marsh, Phospholipid Bilayers, 1987, John Wiley, New York

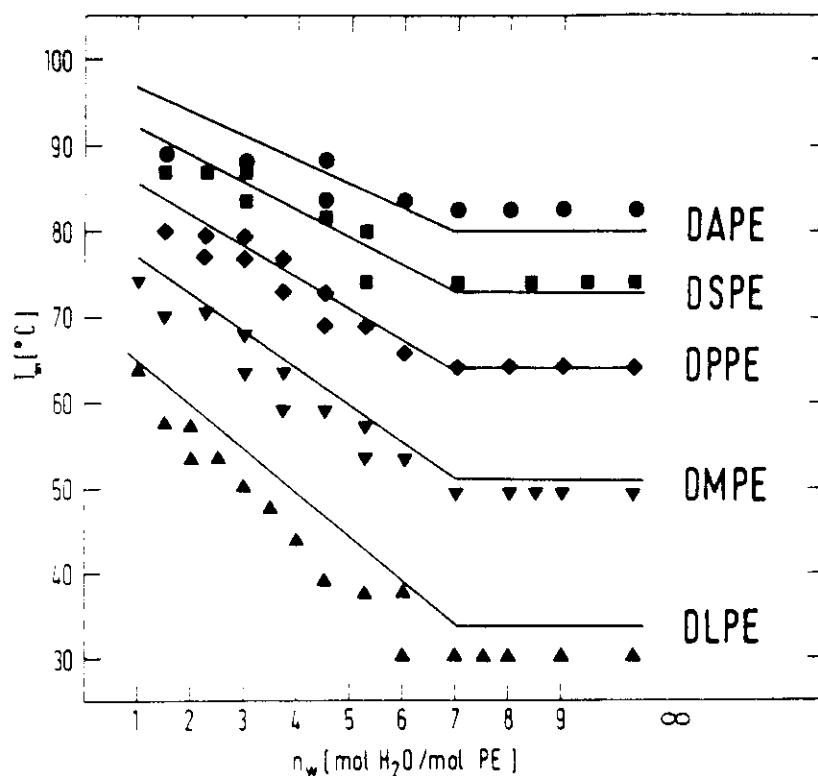
Chainlength-dependence of T_m



PA: Phosphatidic acid
PE: Phosphatidylethanolamine
PC: Phosphatidylcholine
PG: Phosphatidylglycerol
PS: Phosphatidylserine
CL: Cardiollipin

From: J. M. Seddon and G. Cevc, in *Phospholipids Handbook* (G. Cevc, Ed.), 1993, Marcel Dekker, New York.

Hydration-dependence of T_m

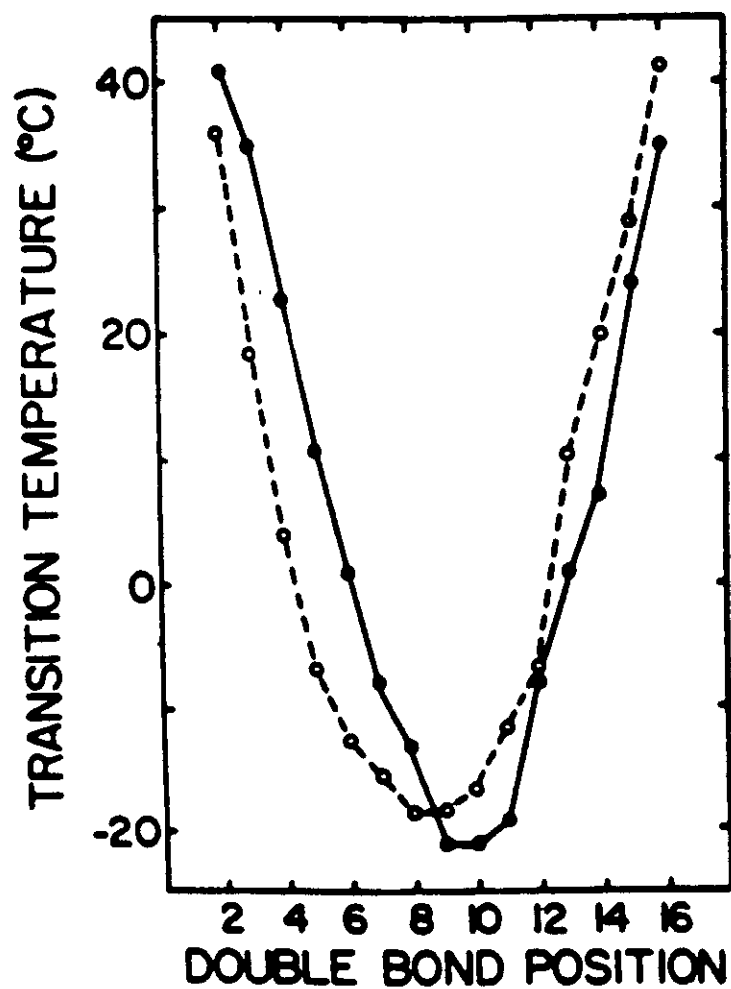


DLPE:	di-C12	Phosphatidylethanolamine
DMPE:	di-C14	"
DPPE:	di-C16	"
DSPE:	di-C18	"
DAPE:	di-C20	"

From: J. M. Seddon and G. Cevc, in *Phospholipids Handbook* (G. Cevc, Ed.), 1993, Marcel Dekker, New York.

Effect of double bond position on T_m

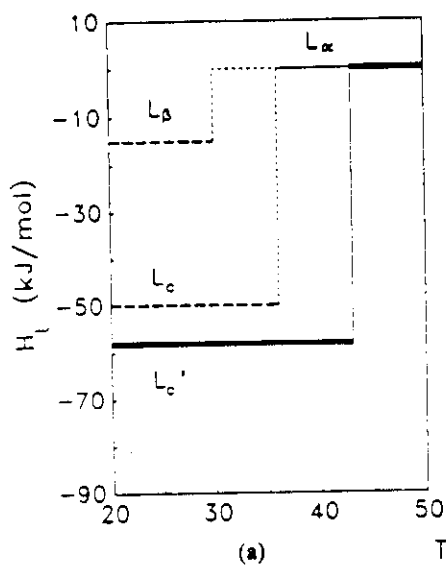
cis-unsaturated PC / H₂O



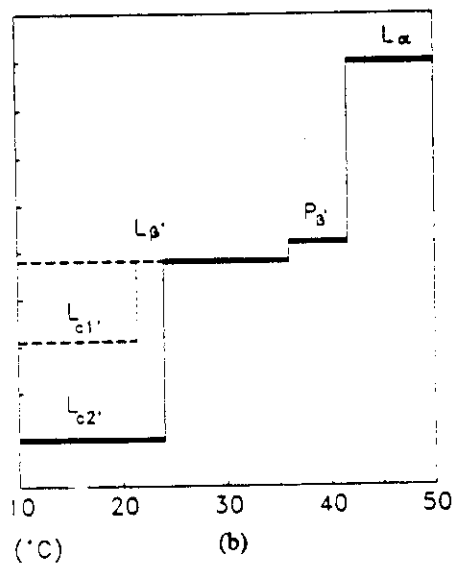
From: C. B. Berde et al., 1983, *Biochemistry*, **19**, 4279-4293.

Lipid bilayer metastability

DLPE



DPPC

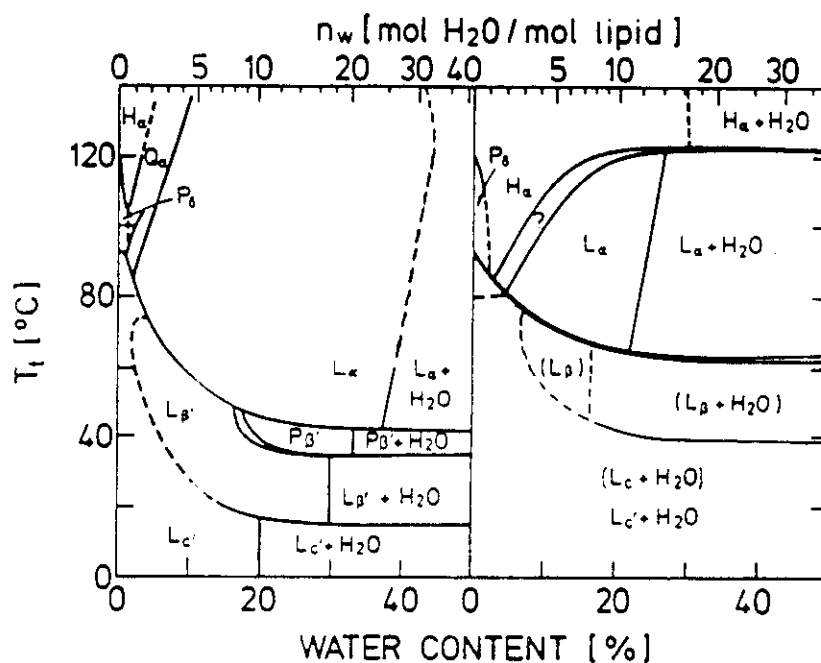


From: J. M. Seddon and G. Cevc, in *Phospholipids Handbook* (G. Cevc, Ed.), 1993, Marcel Dekker, New York.

Phospholipid / H₂O phase diagrams

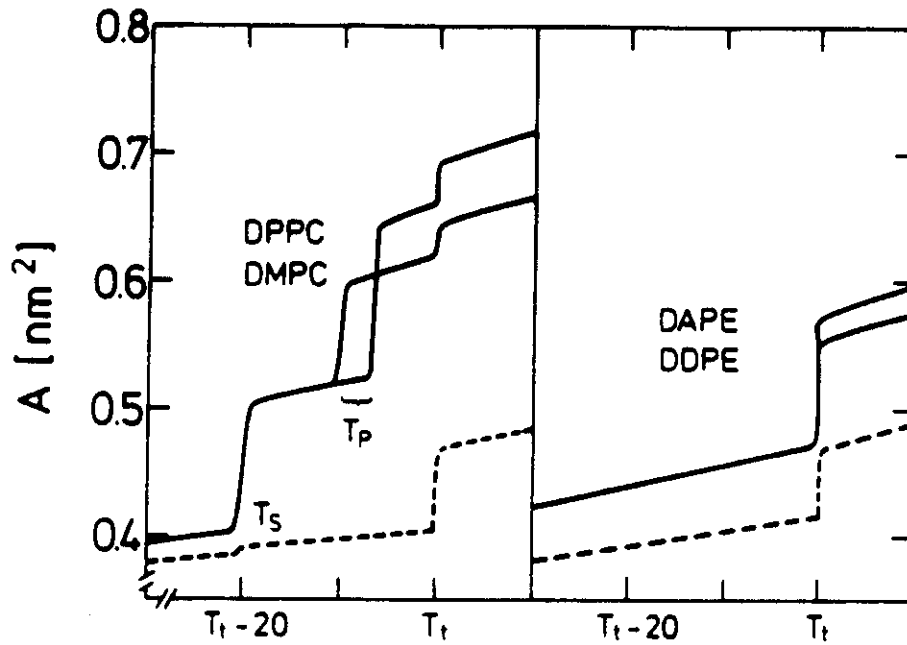
DPPE

DPPE

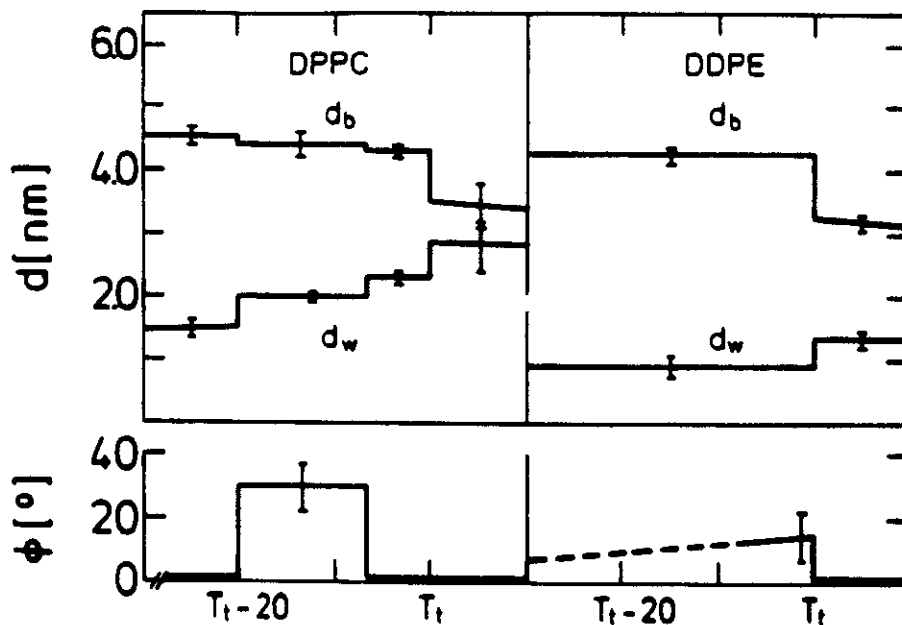


From: J. M. Seddon and G. Cevc, in Phospholipids Handbook (G. Cevc, Ed.), 1993, Marcel Dekker, New York.

Temperature-dependence of lipid molecular area for PC and PE



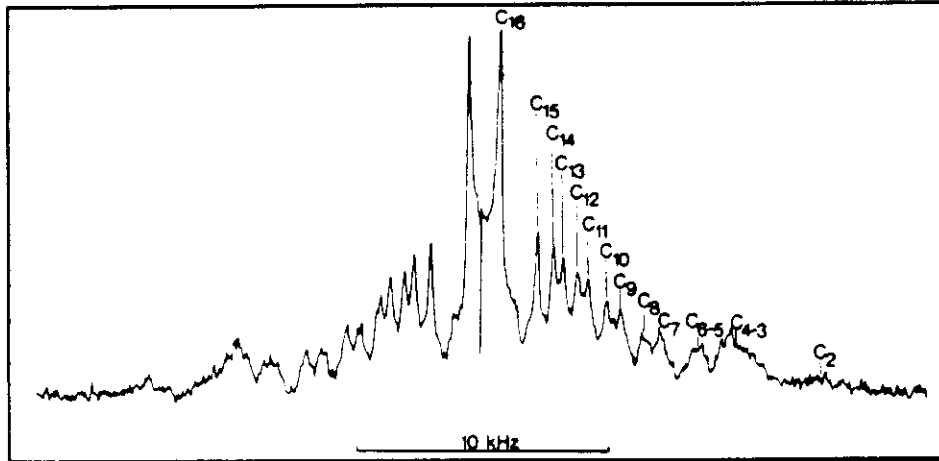
Temperature-dependence of lipid and water layer thickness for PC and PE



From: G. Cevc and D. Marsh, Phospholipid Bilayers, 1987, John Wiley, New York

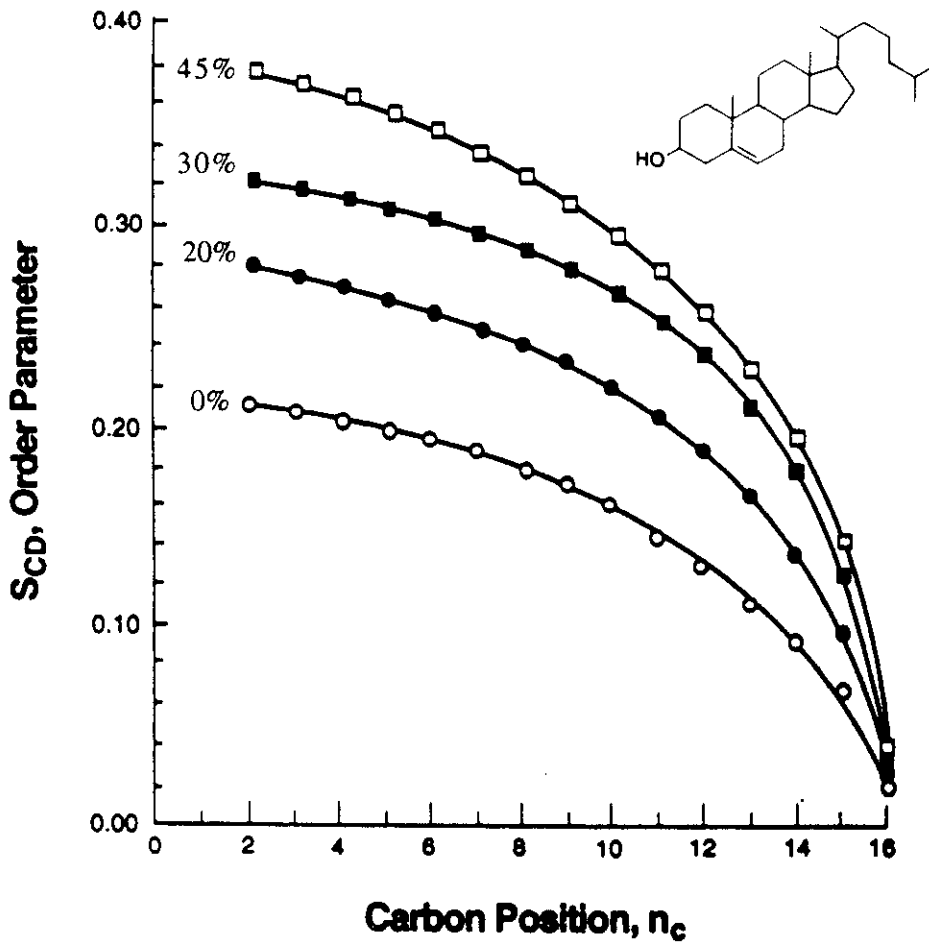
Order parameter profile of bilayer

$^2\text{H-NMR}$: POPC/cholate/water



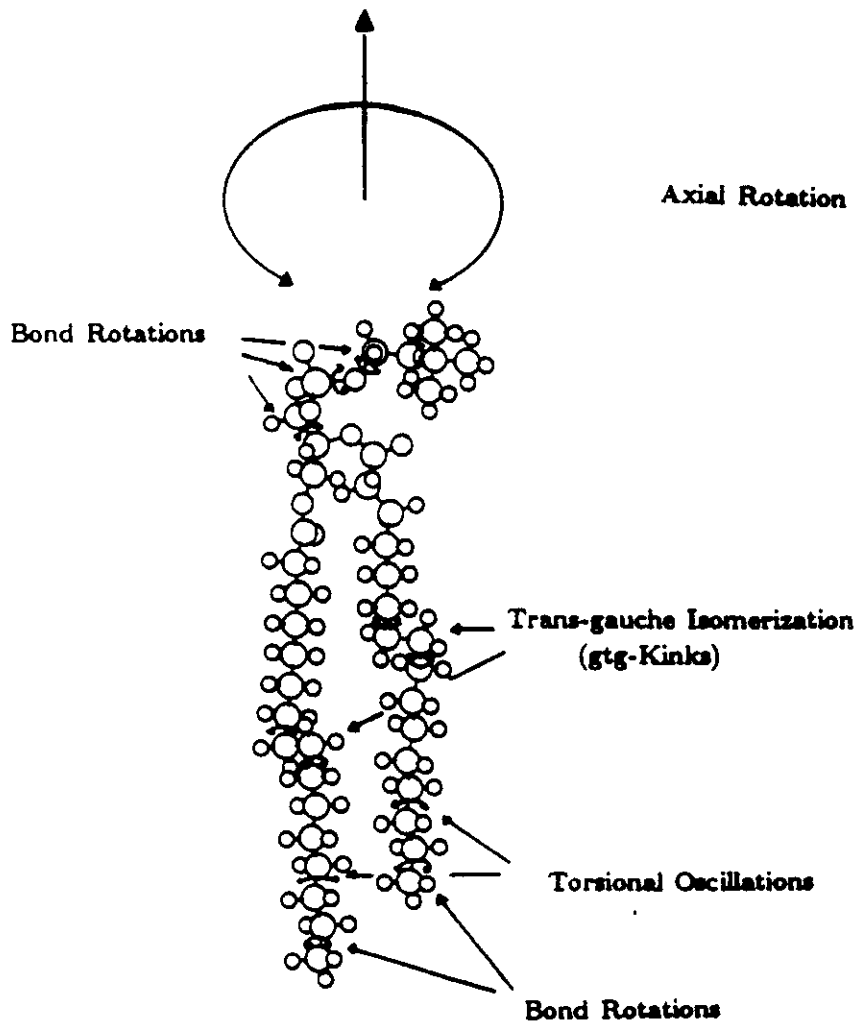
From: G. Lindblom et al., *Biochemistry*, 1982, **21**, 1553

Addition Of Cholesterol To POPC



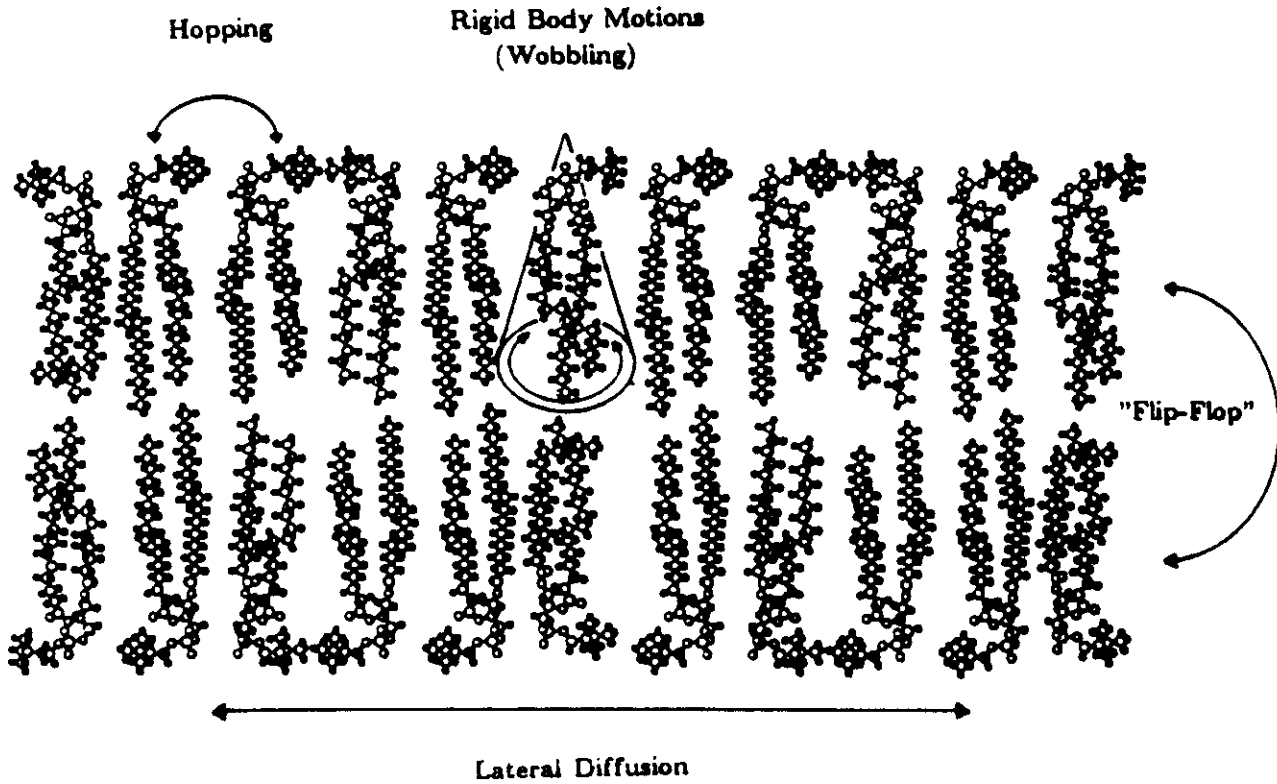
From: G. Lafleur et al., *Eur. Biophys. J.*, 1990, **19**, 55

Motions of lipids in membranes



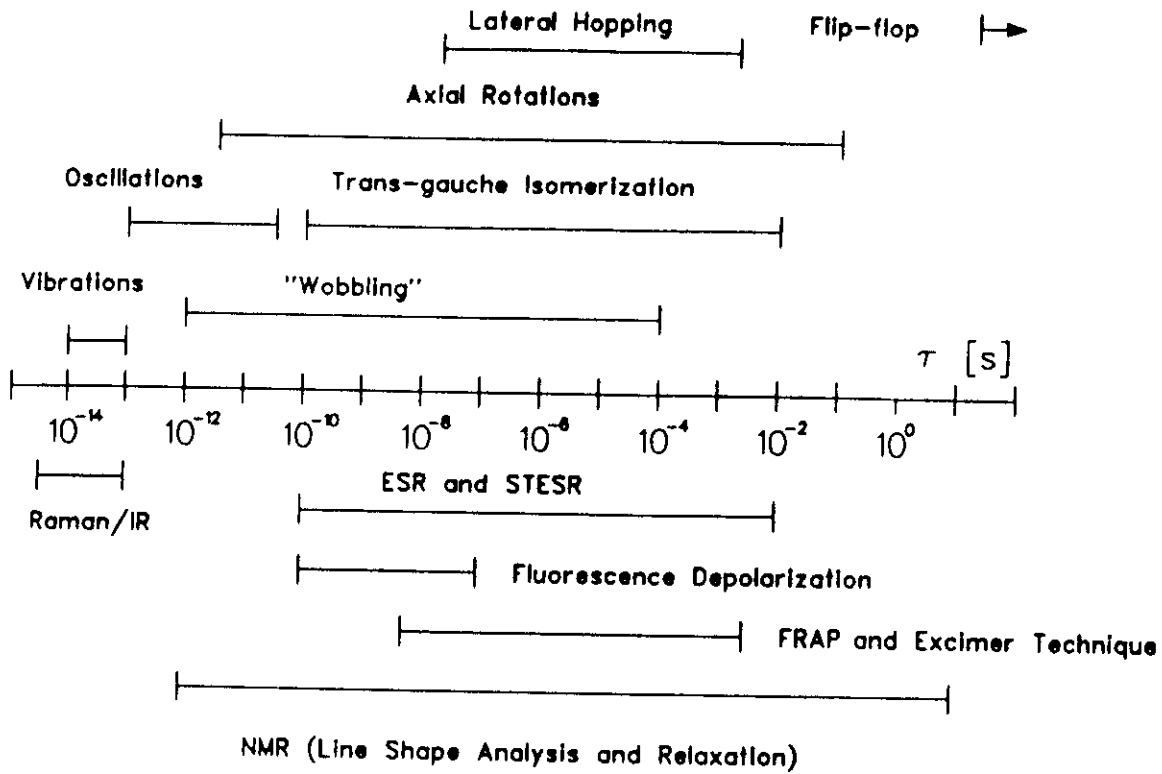
From: A. Blume, in Phospholipids Handbook (G. Cevc, Ed.), 1993, Marcel Dekker, New York.

Molecular motions in membranes



From: A. Blume, in *Phospholipids Handbook* (G. Ceve, Ed.), 1993, Marcel Dekker, New York.

Molecular motions in membranes: correlation times and time ranges



From: A. Blume, in Phospholipids Handbook (G. Cecc, Ed.), 1993, Marcel Dekker, New York.

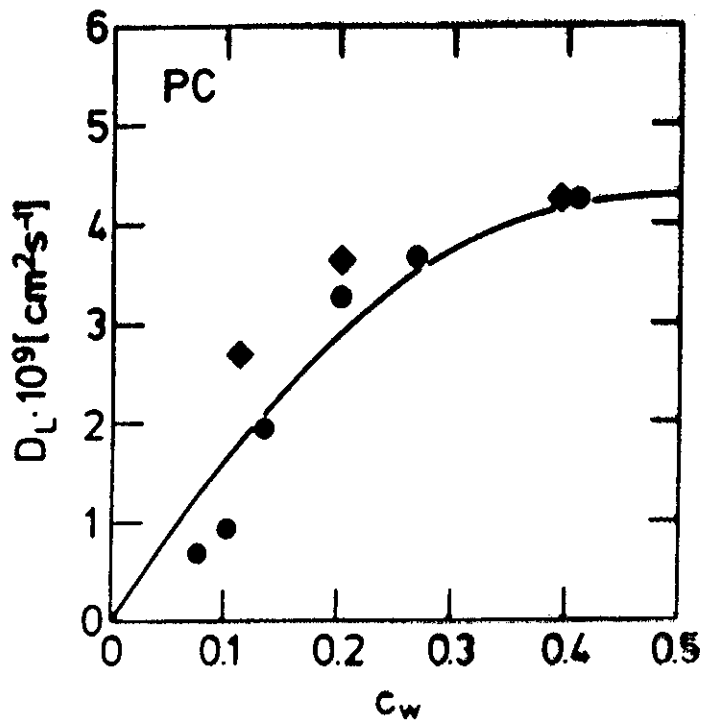
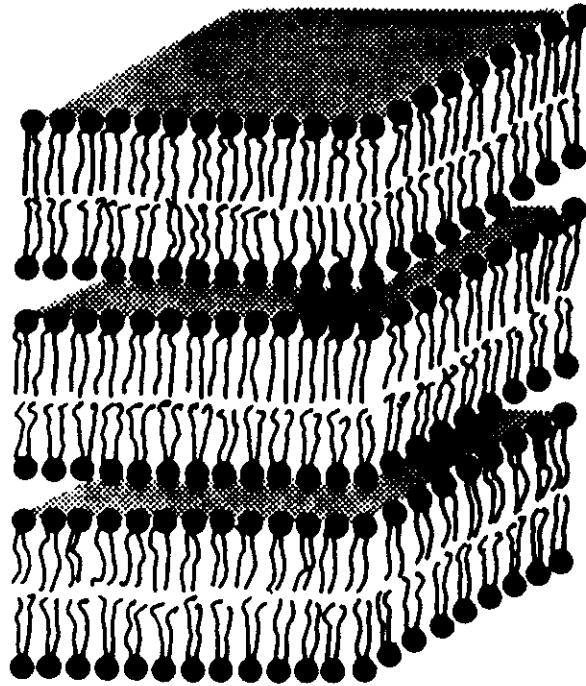


Figure 3.14. The lipid lateral diffusion constant in phosphatidylcholine bilayers as a function of the water concentration. Experimental points are deduced from the data of Seddon et al. (1978) and McCown et al. (1981), and the curves are from Eq. (3.41), with the parameters given in the text. [From Cevc and Seddon (1985), courtesy Plenum Press.]

Fluid Lamellar L α Phase



Typical values for phospholipids:

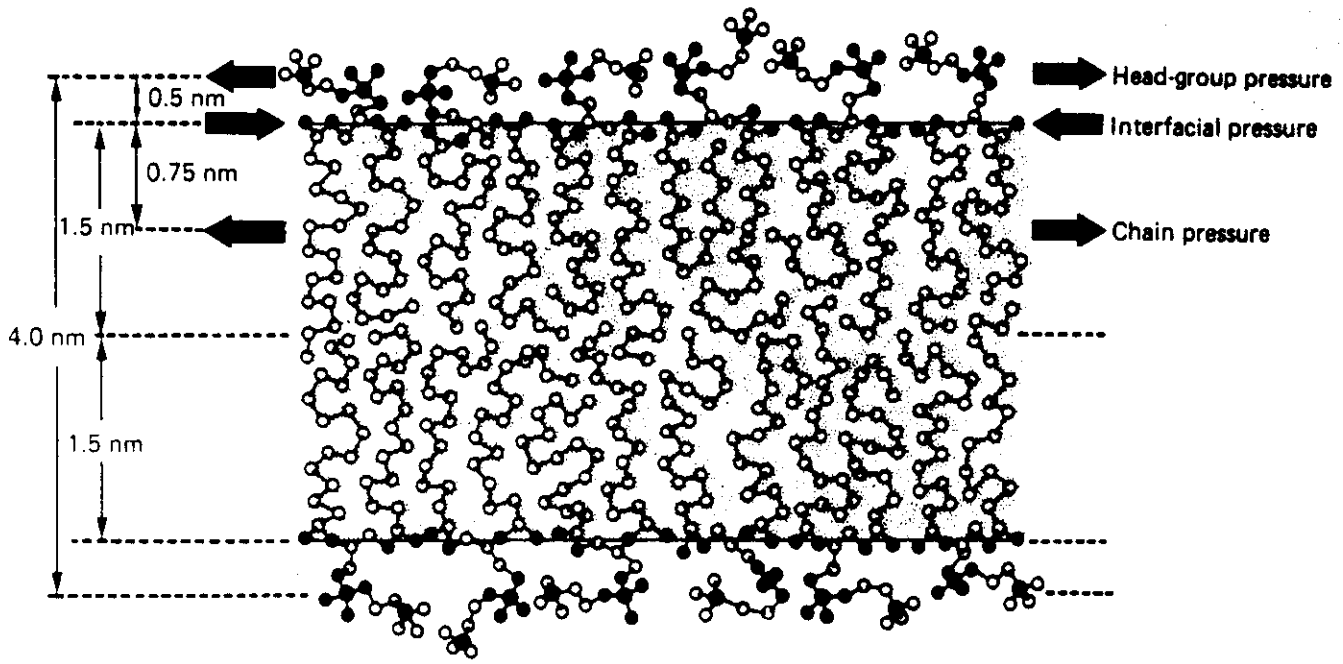
Area compressibility modulus: $0.1 - 1 \text{ Nm}^{-1}$

Bending modulus: $4 \times 10^{-20} \text{ J}$ (10 kT)

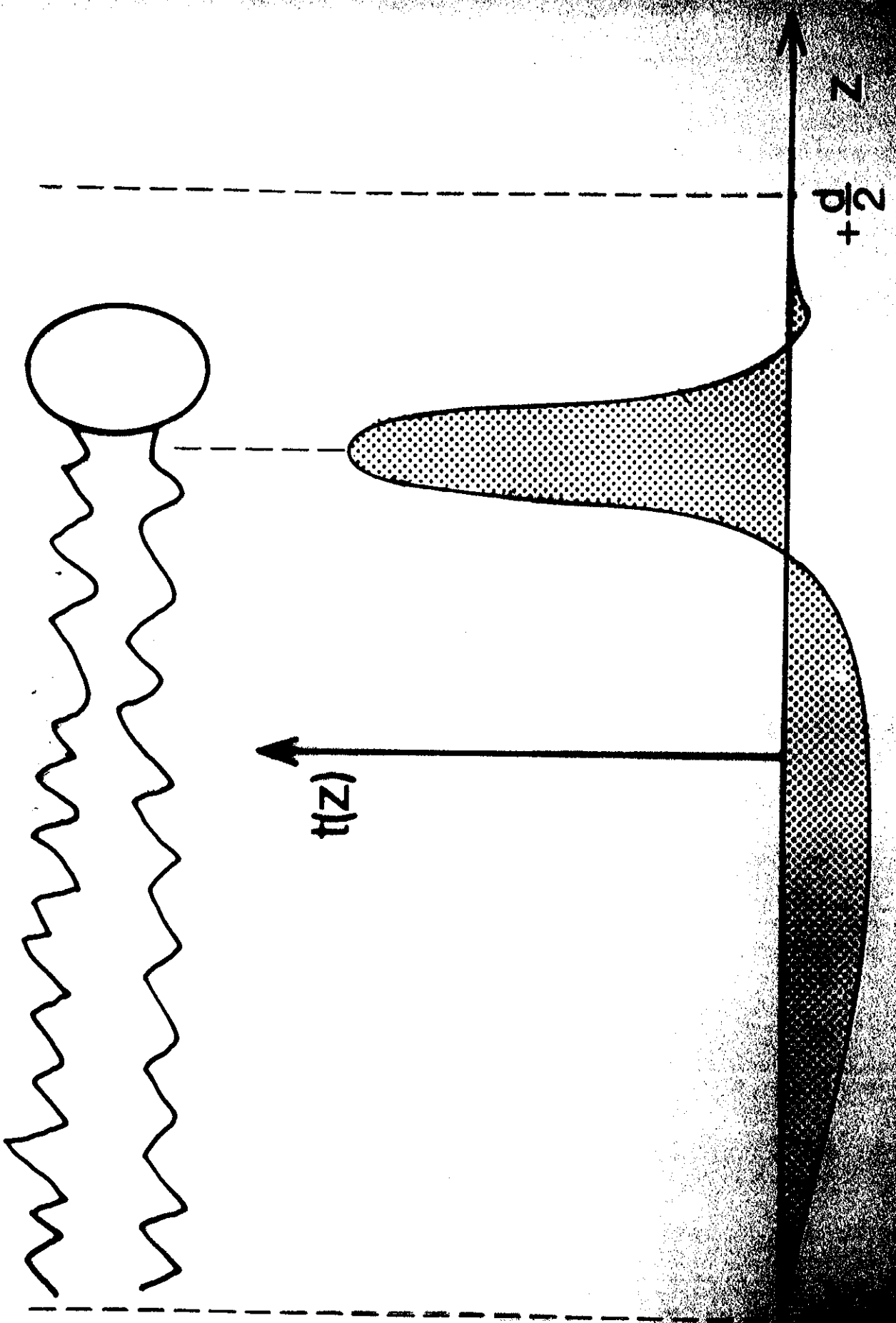
Lateral diffusion constant: $10^{-11} - 10^{-12} \text{ m}^2\text{s}^{-1}$

Rotational correlation time: $10^{-9} - 10^{-8} \text{ s}$

Balance of forces in fluid lipid bilayers

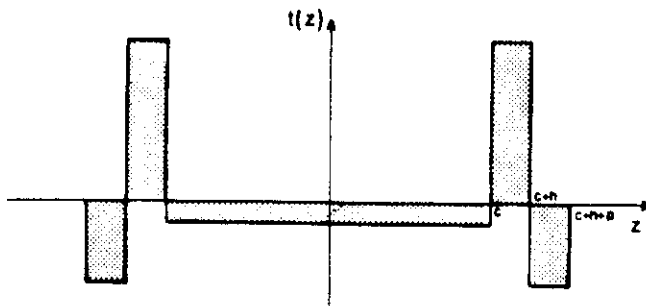


From: J. Israelachvili, 1991, *Intermolecular and Surface Forces*, Academic Press, London.

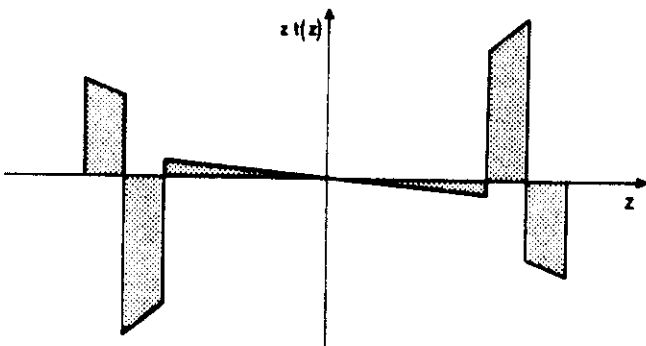


Bilayer curvature elasticity

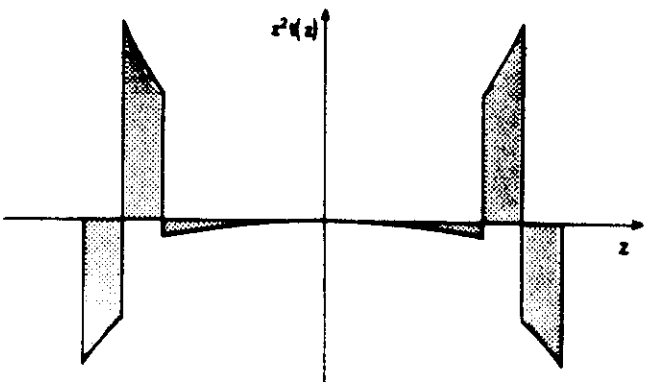
$$g = 1/2 \kappa_M (c_1 + c_2 - c_0)^2 + \kappa_G c_1 c_2$$



$$\sigma = \int t(z) dz = 0$$



$$\tau = -\kappa_M c_0 = \int t(z) z dz = 0$$



$$\kappa_G = \int t(z) z^2 dz > 0$$

INTERFACIAL CURVATURE

Mean curvature

$$H = 1/2(c_1 + c_2)$$

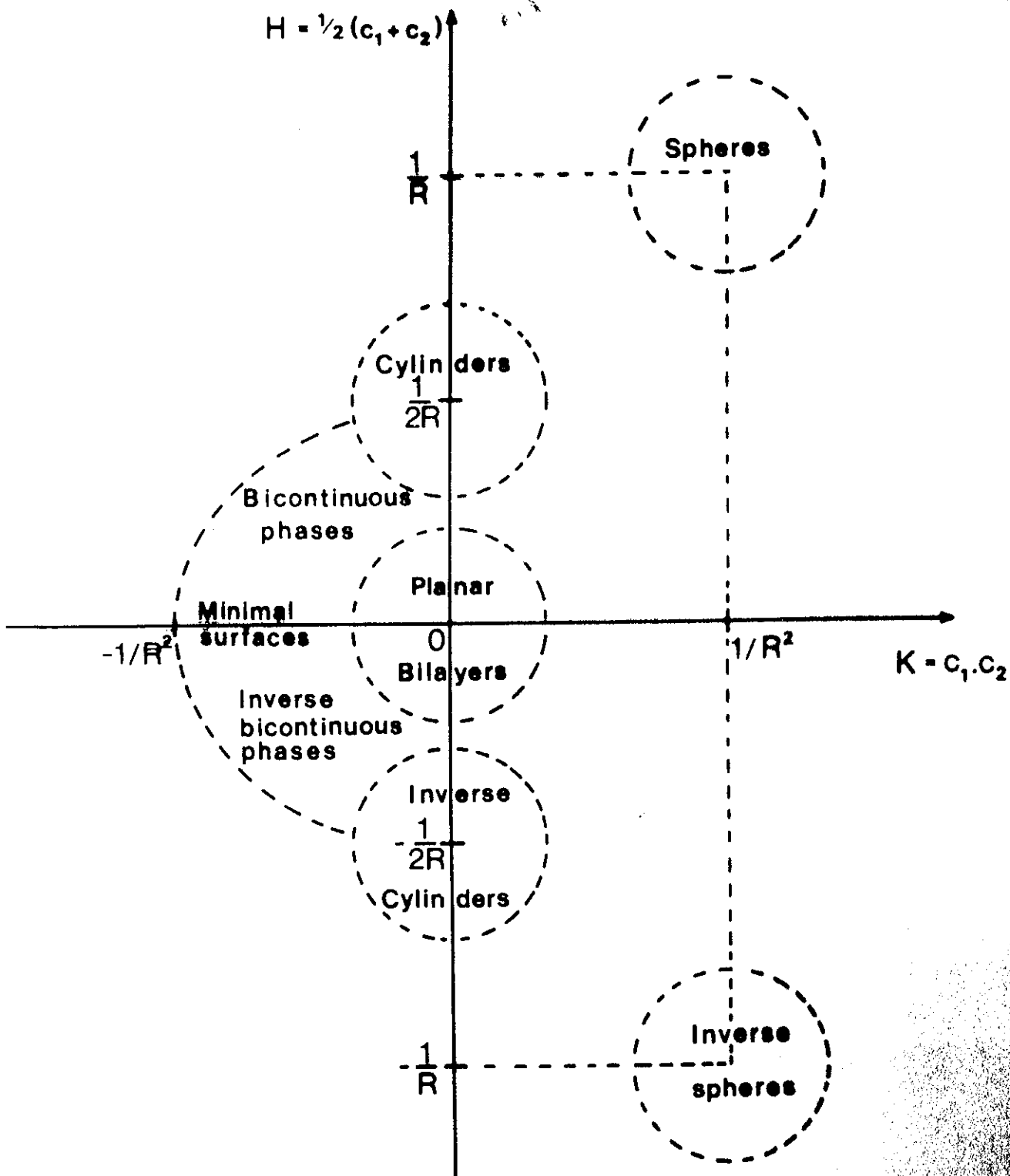
Gaussian curvature

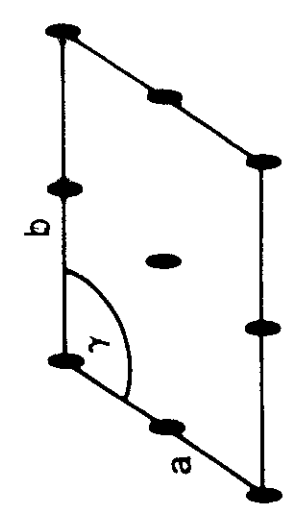
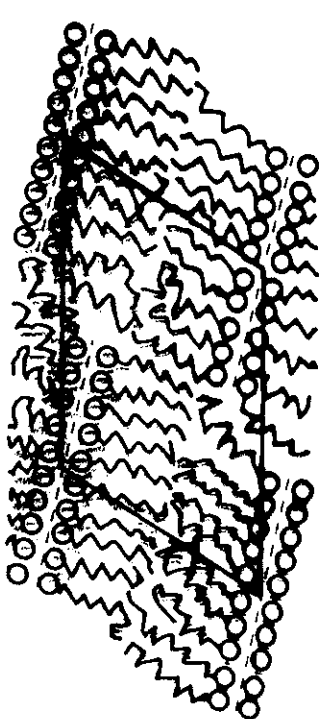
$$K = c_1 \cdot c_2$$

Principal curvatures:

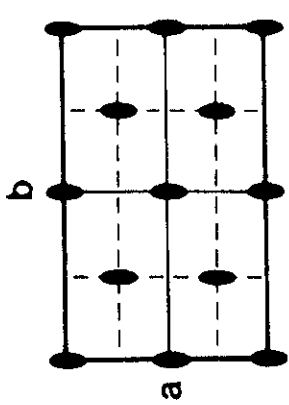
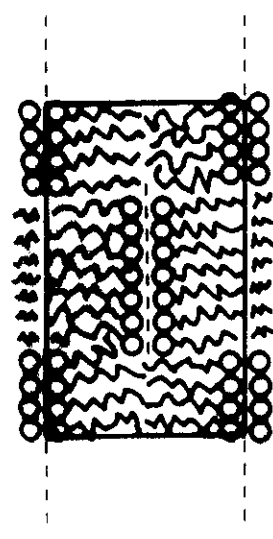
$$c_1 = 1/R_1 ; c_2 = 1/R_2$$

K	H	Form of surface	Example
>0	> or < 0	Elliptic	Sphere
0	> or < 0	Parabolic	Cylinder
0	0	Parabolic	Plane
<0	> or < or = 0	Hyperbolic	Saddle surface

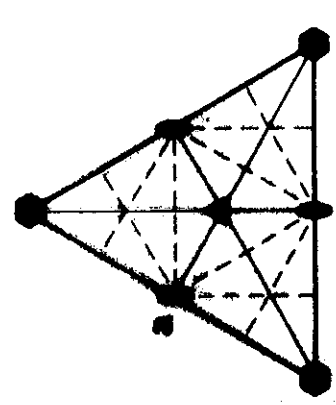
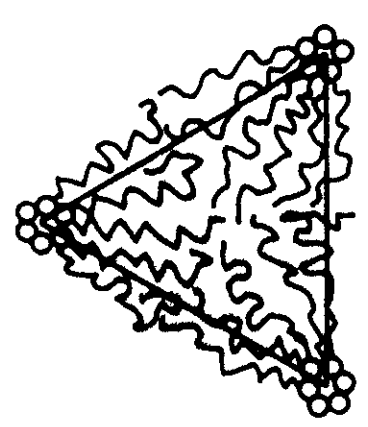




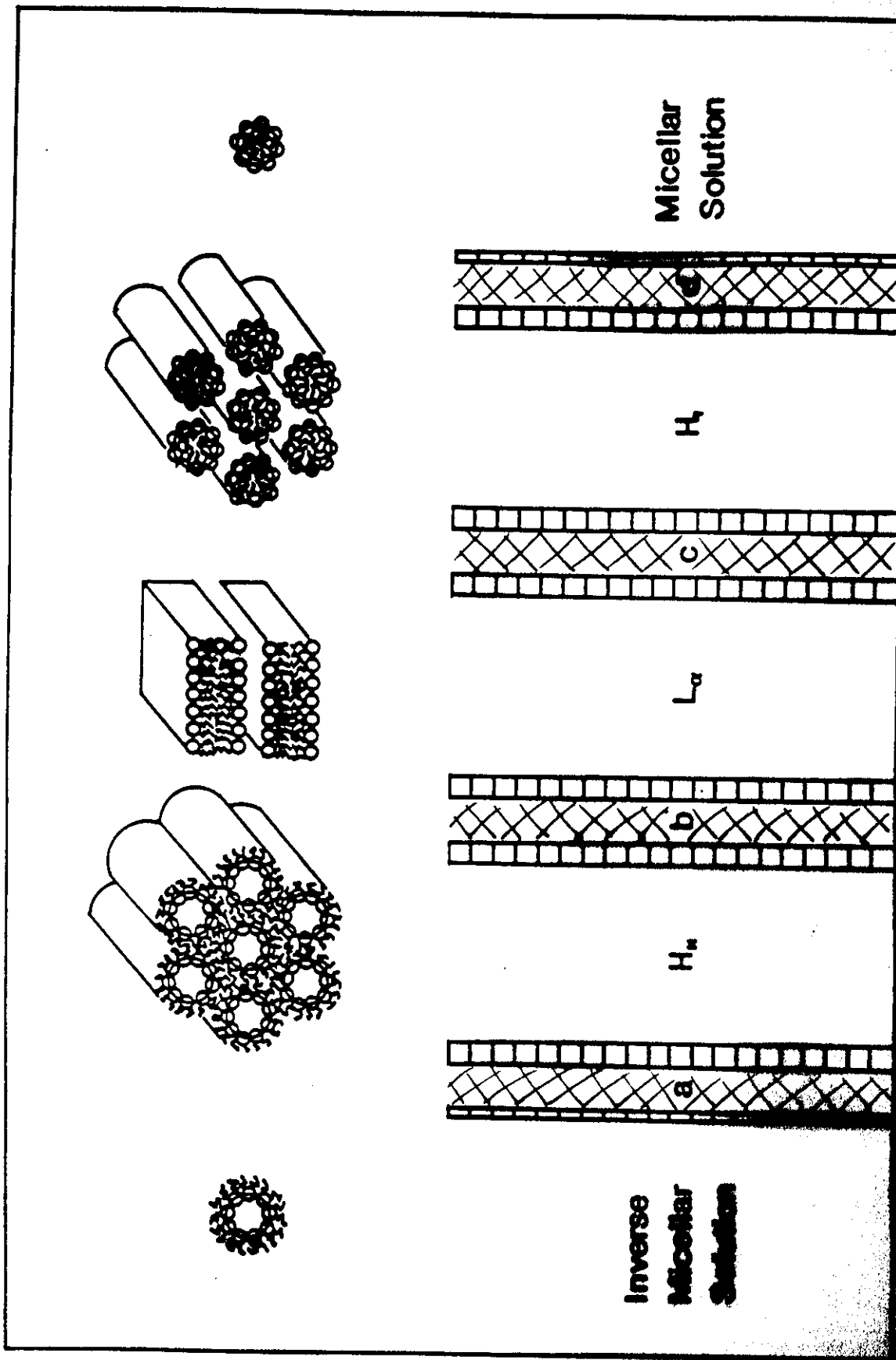
a) M_{11}
oblique
 $p2$

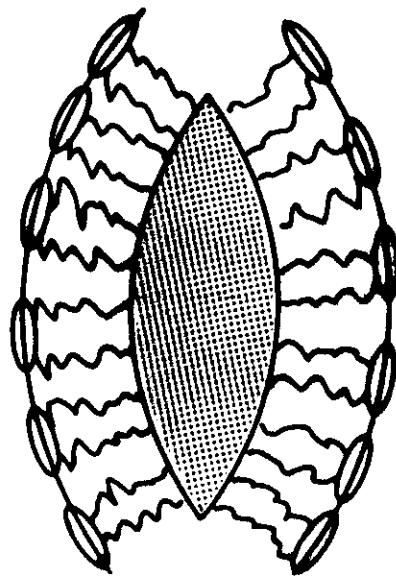


b) P_n
centred rectangular
 cm

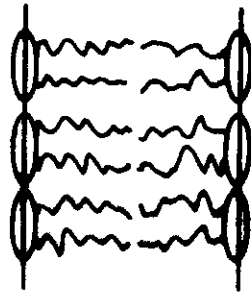


c) H_n
hexagonal
 $p6m$

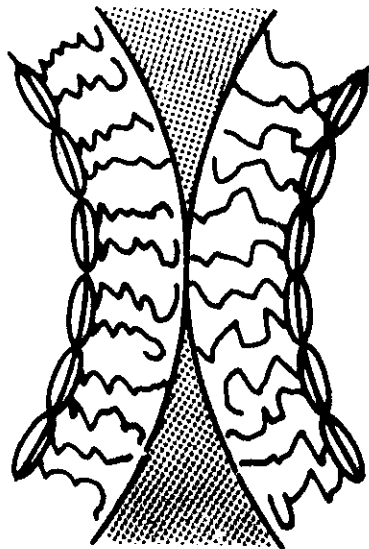


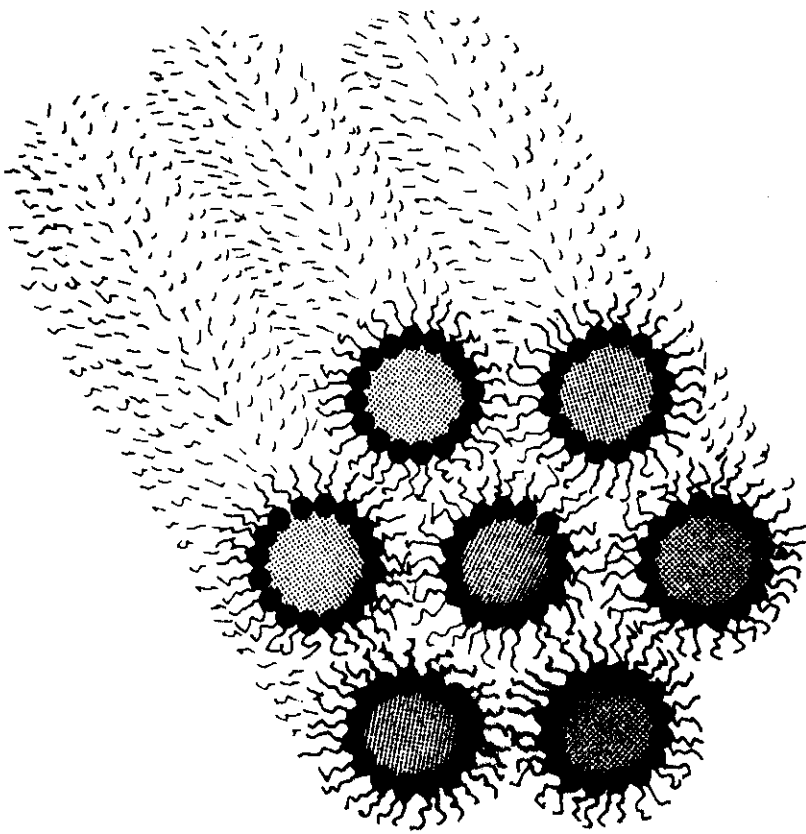


H_2O ↑

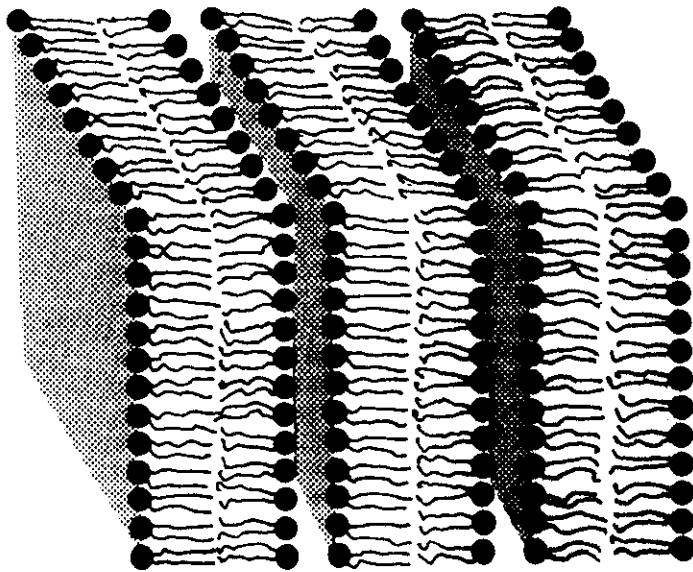


↓ T

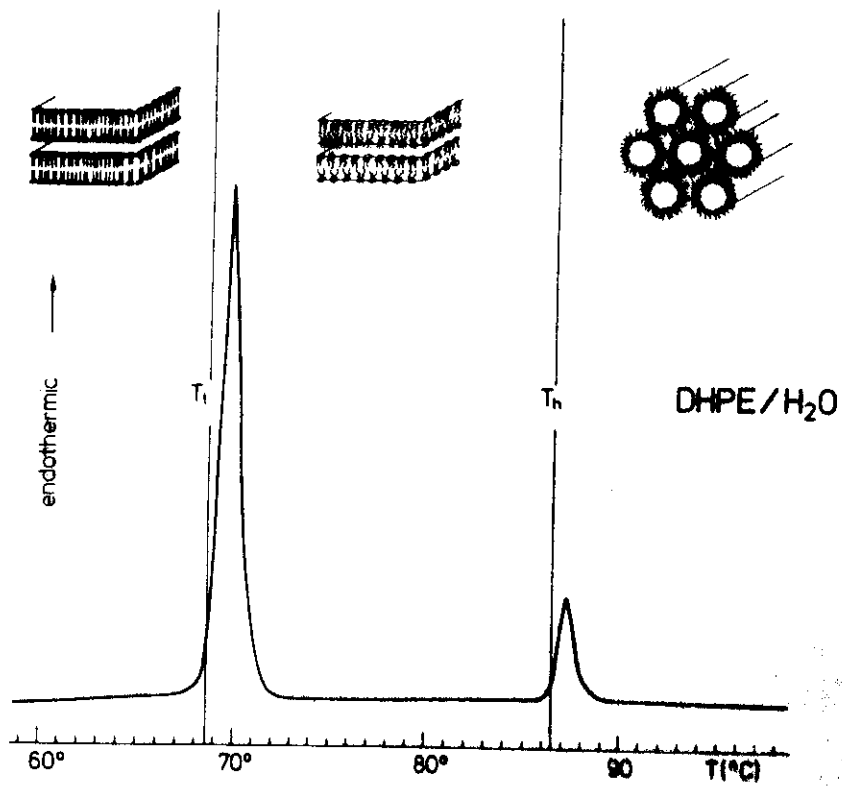


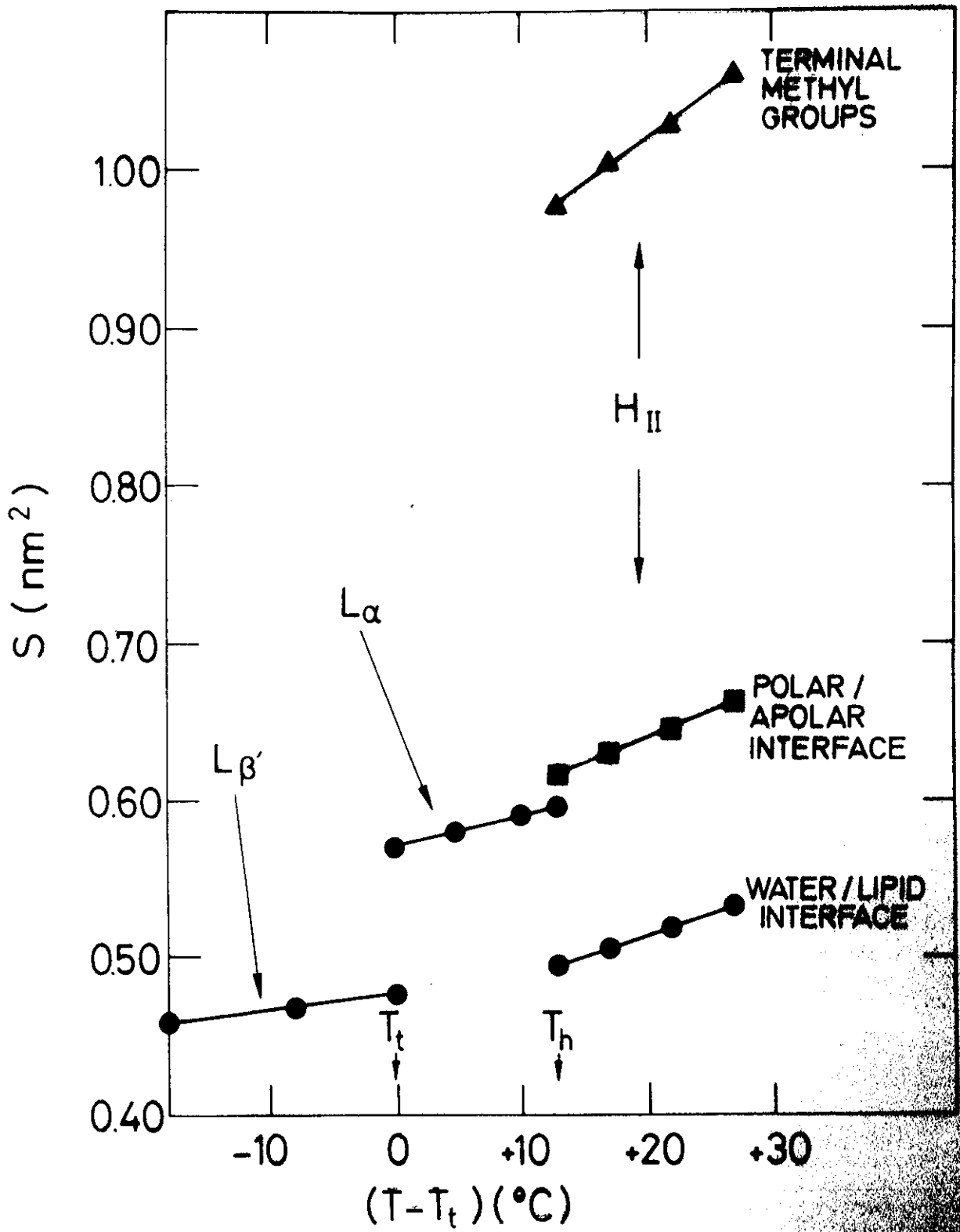


HEXAGONAL H_{II}

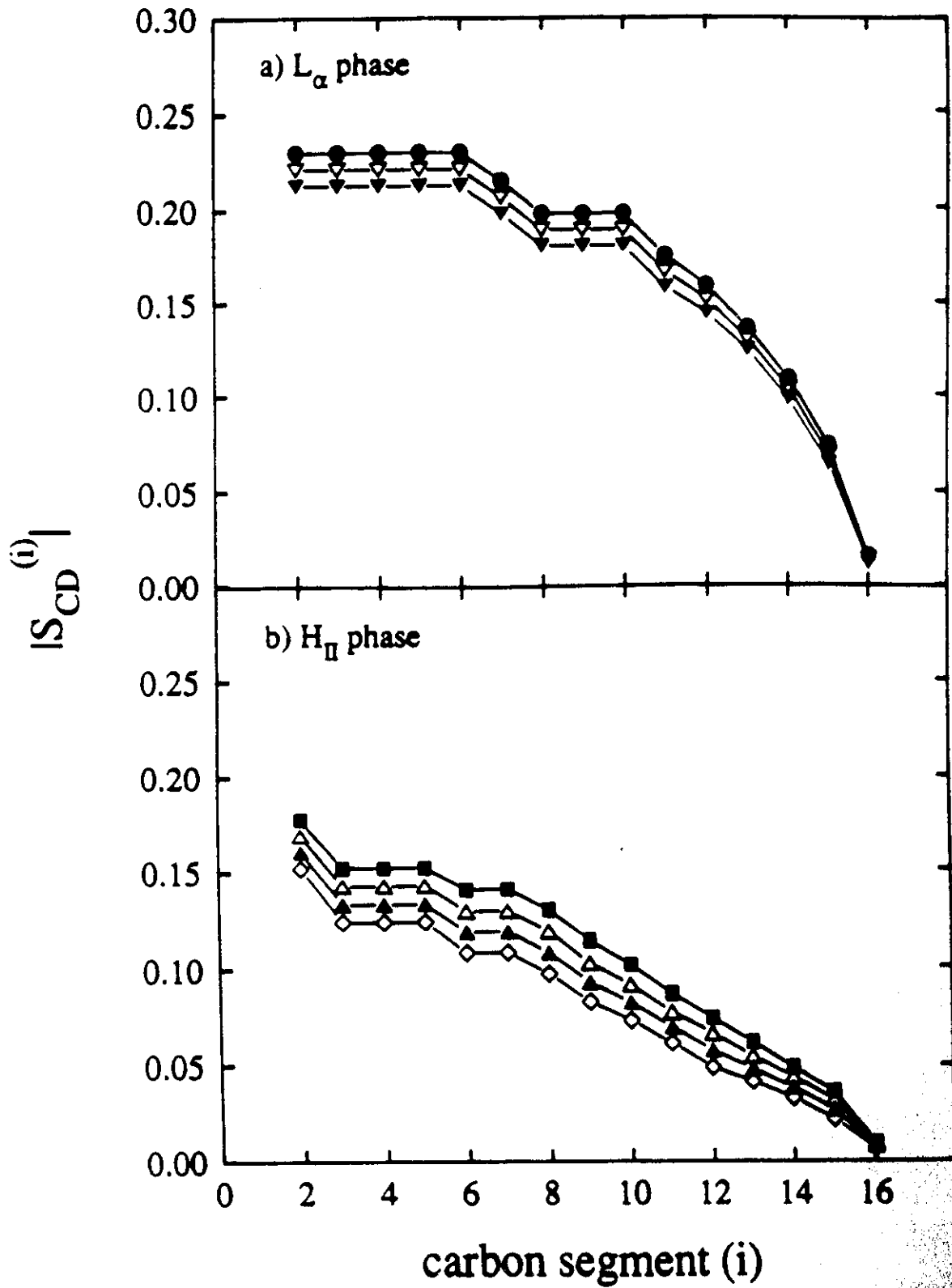


LAMELLAR L_α

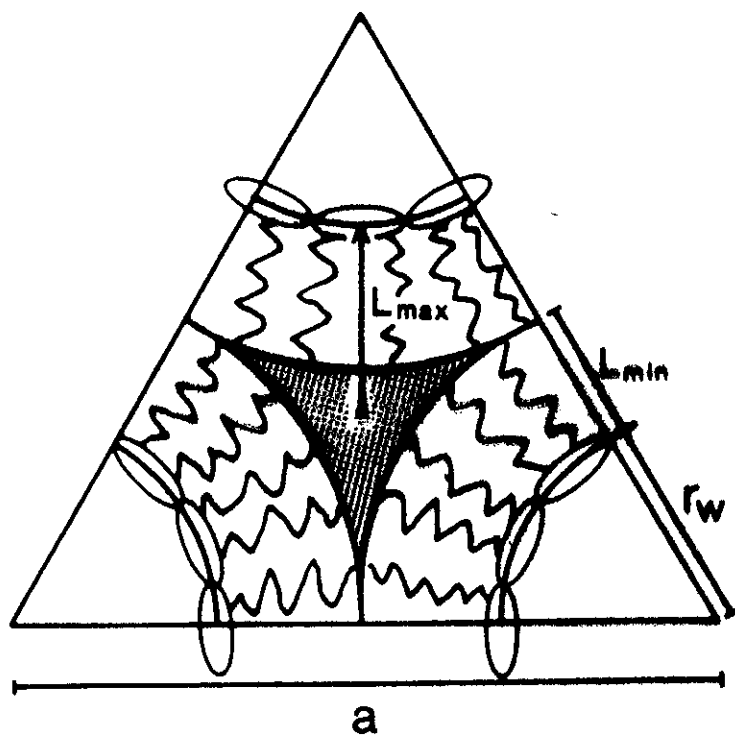


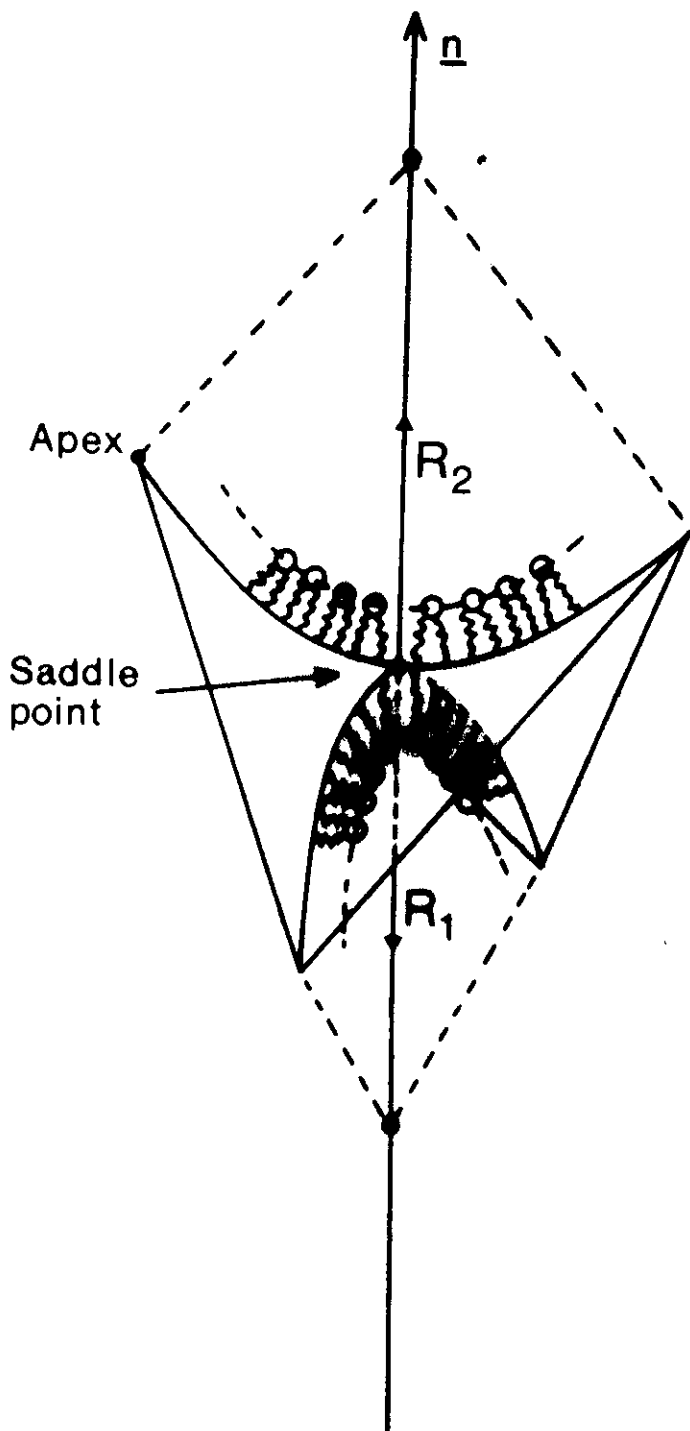


Order parameter profiles of PLPE in the L_{α} and H_{II} phases



From: R. L. Thurmond et al., *Biochemistry*, 1993, **32**, 5394-5410





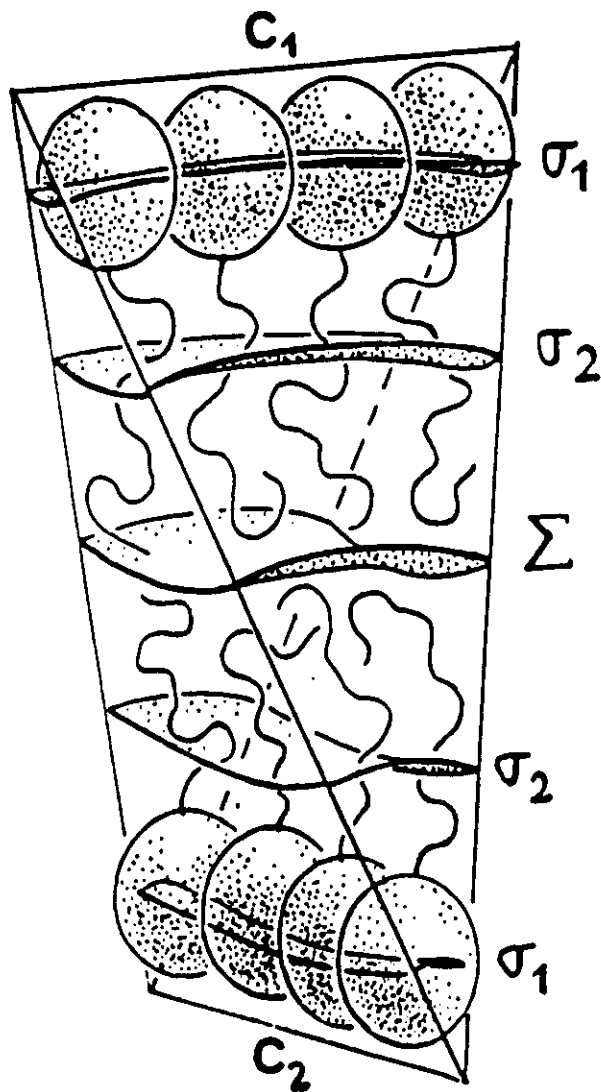
$$C_1 < 0$$

$$C_2 > 0$$

$$\underline{K < 0}$$

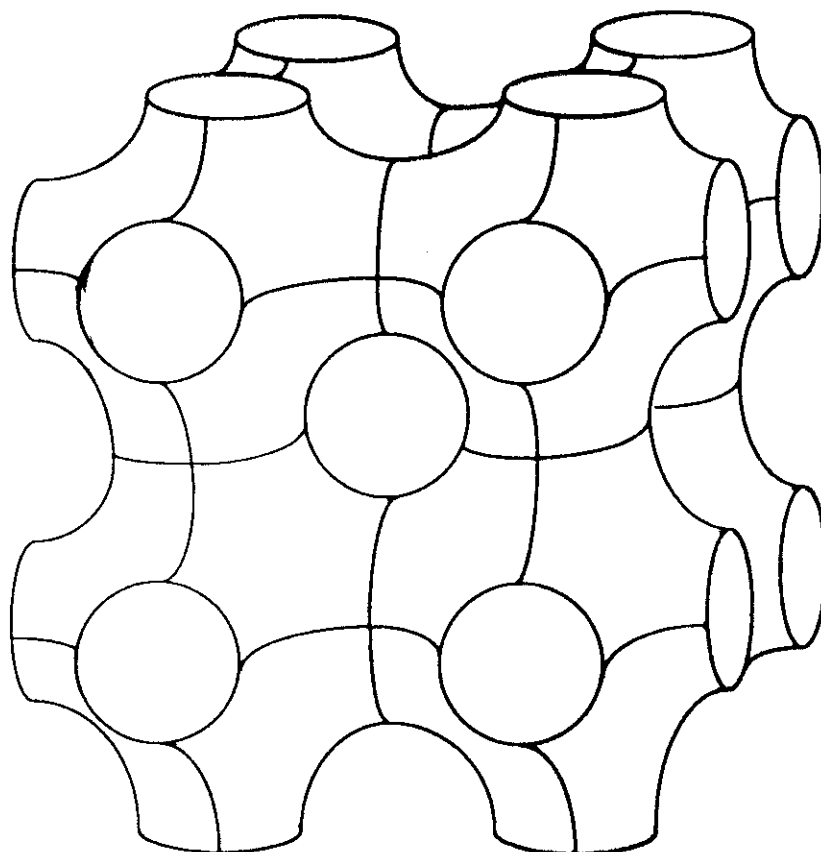
$$(H < \text{or} > \text{or} = 0)$$

Saddle deformation of a bilayer

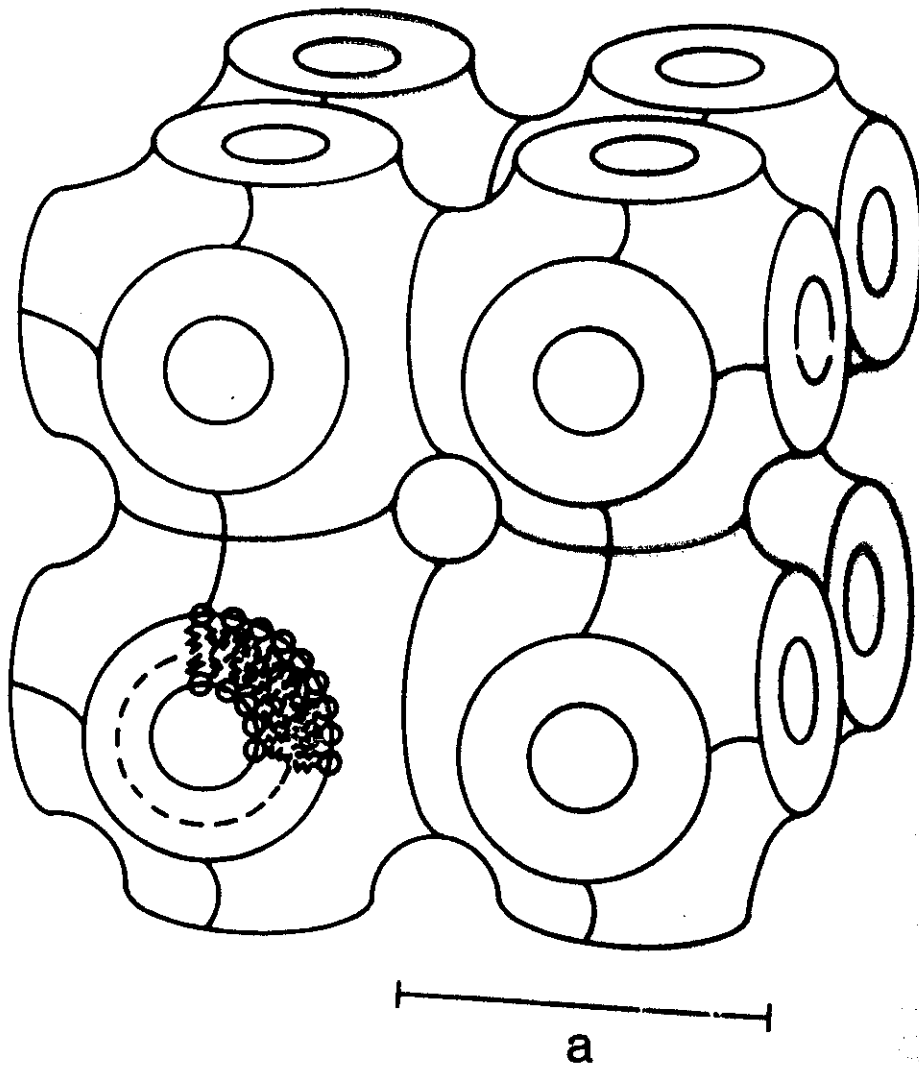


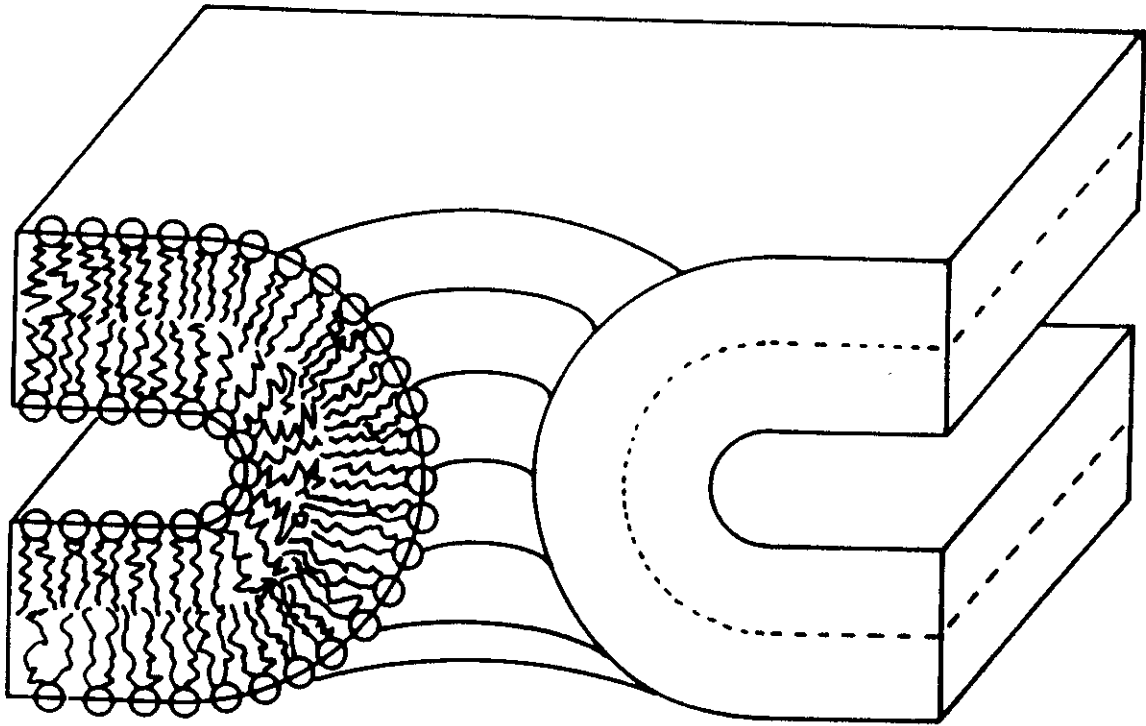
From: Y. Bouligand, J. Phys. (Paris), 1990, 51(C7), 35-52

Schwarz's P-surface ($I\bar{m}\bar{3}m$)

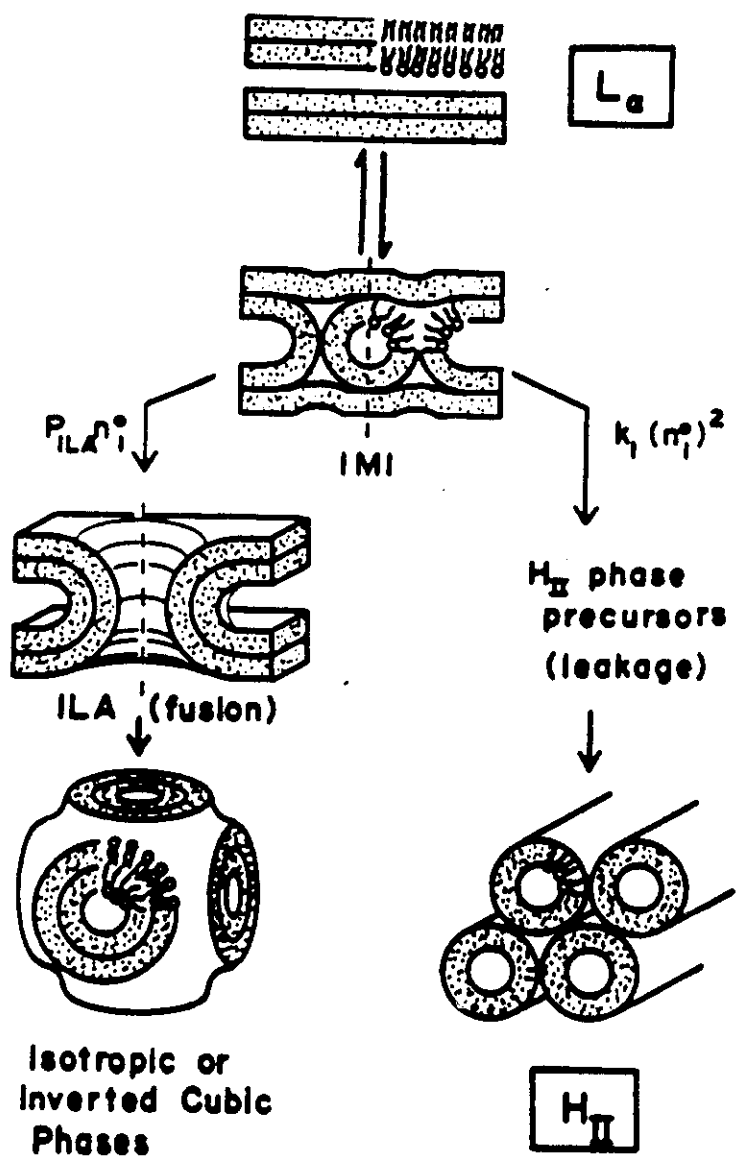


Im3m (Q²²⁹)



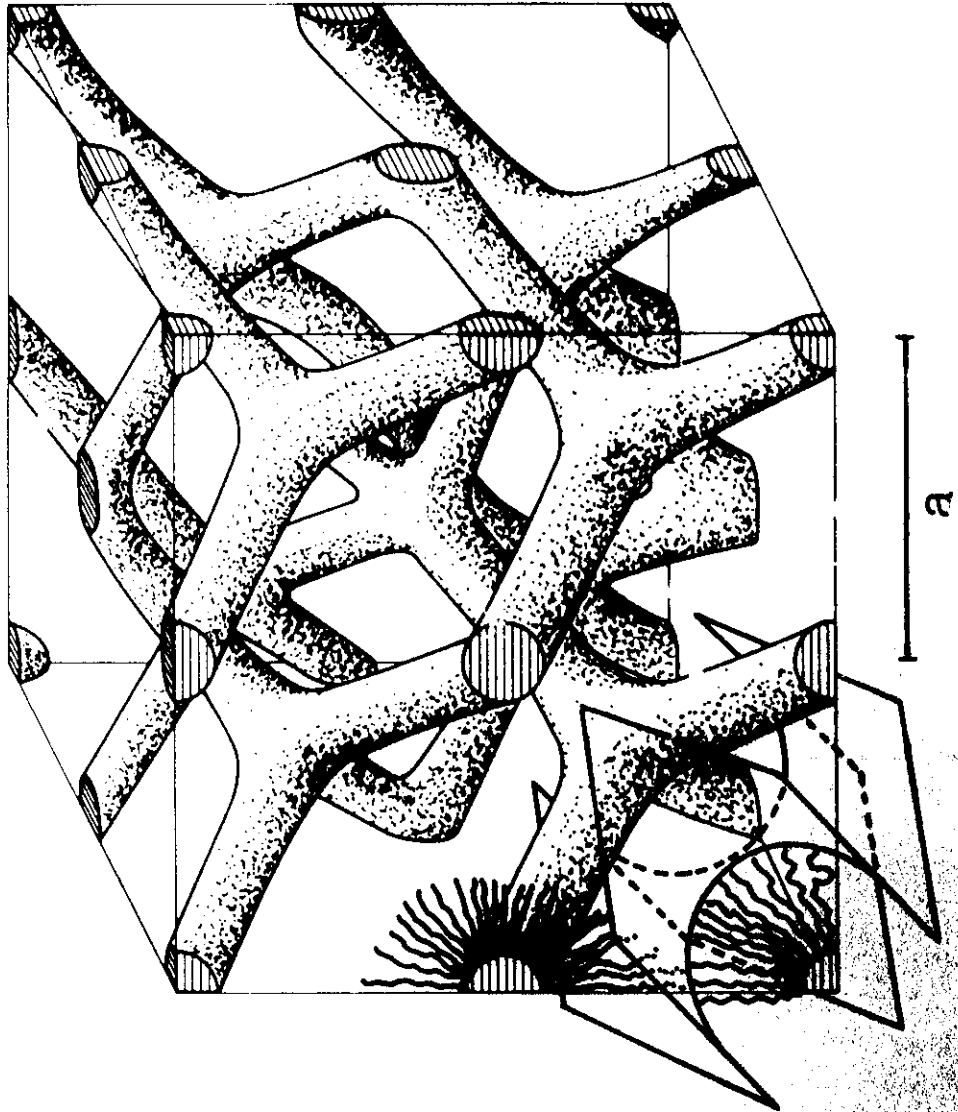


Formation paths of inverse phases



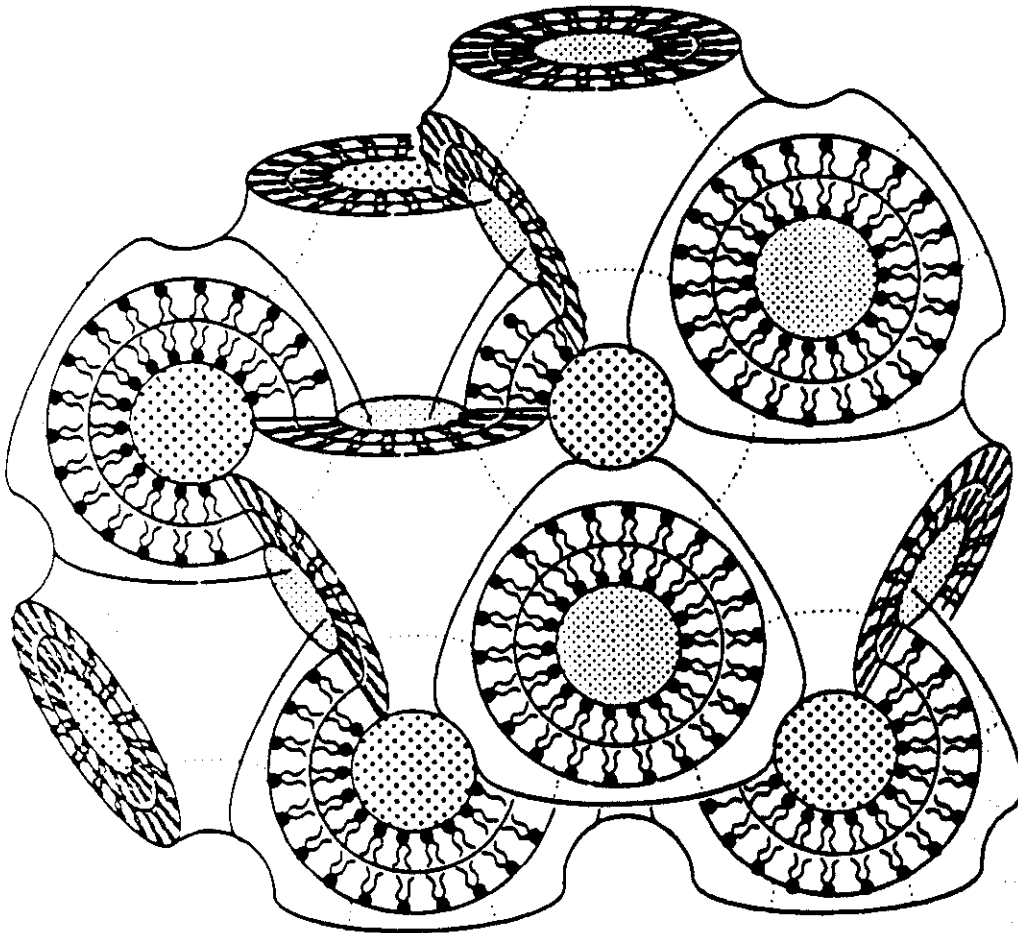
From: D. P. Siegel et al., Biophys. J., 1989, 56, 161-169

Pn 3m (Q²²⁴)



Inverse bicontinuous cubic phases

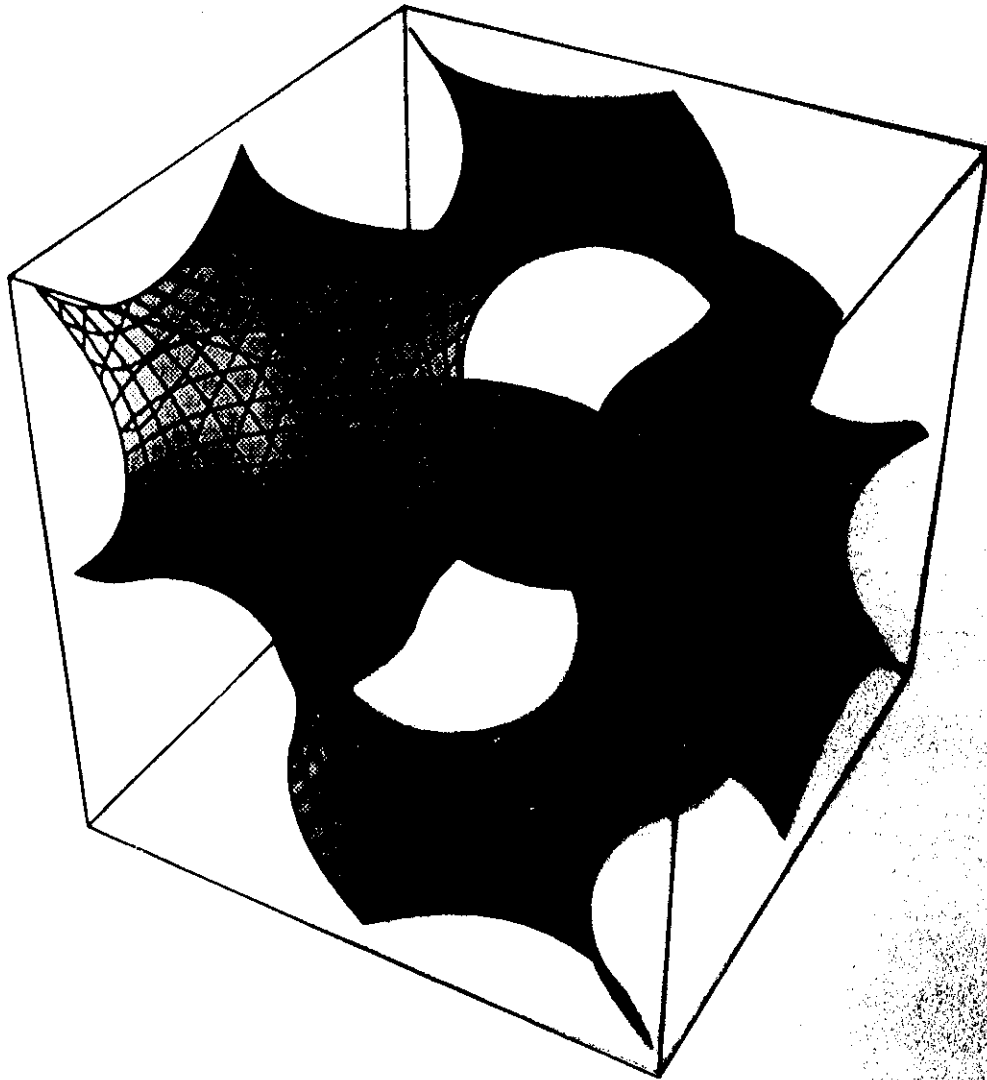
$Pn3m$ (Q224)



From: M. W. Tate et al., *Chem. Phys. Lipids*, 1991, **57**, 147-164

Gyroid Minimal Surface

(Nodal Approximation)



Ia 3d

