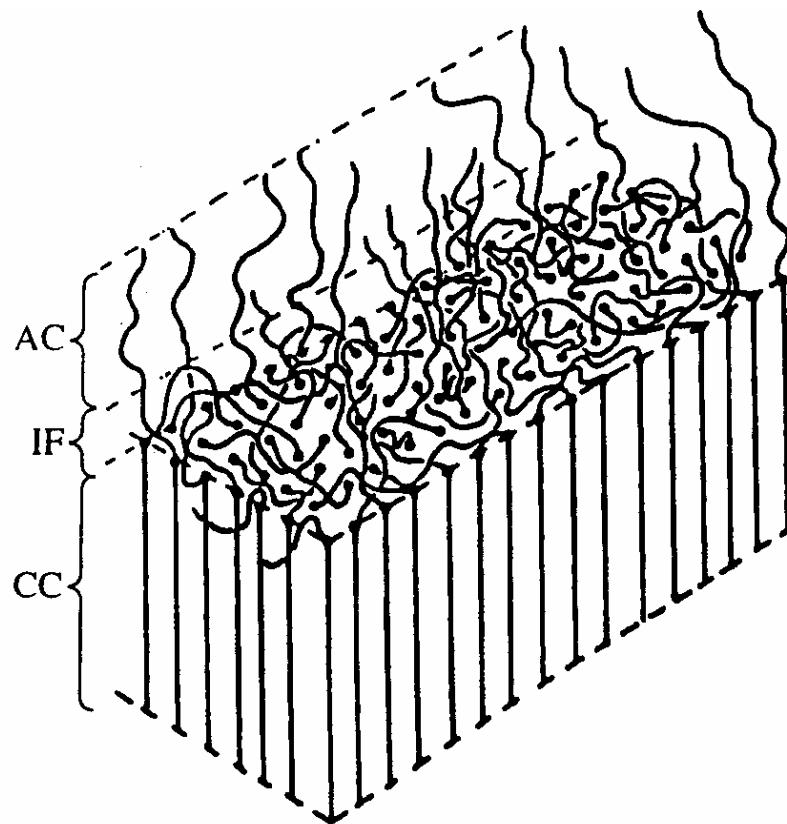


polymer structure and micromechanics

structure

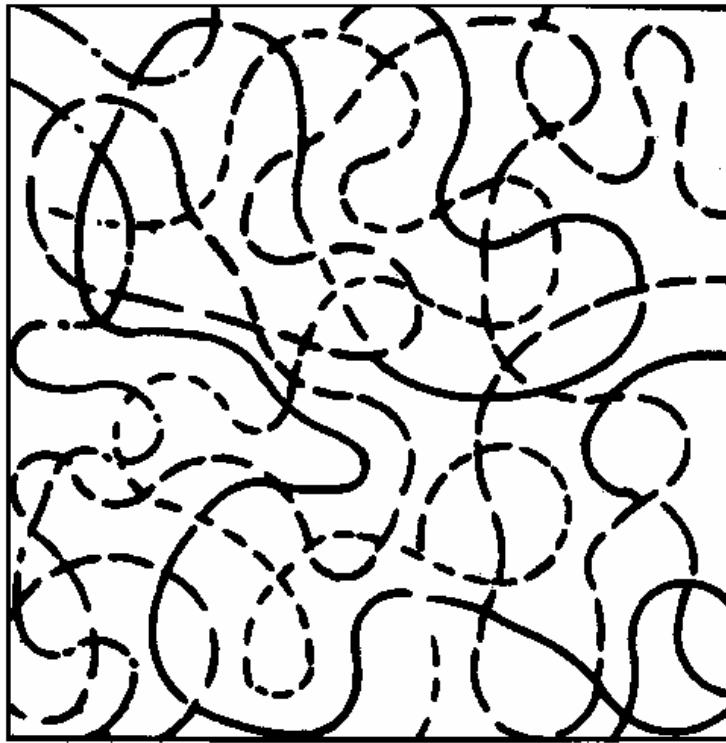
Physical structure



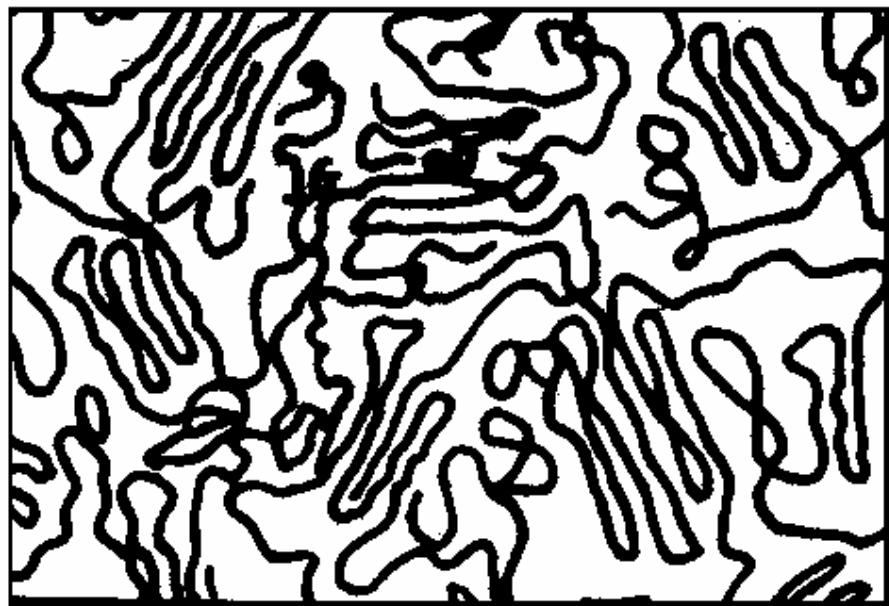
Amorphous components (AC)

Interface (IF)

Crystalline components (CC)

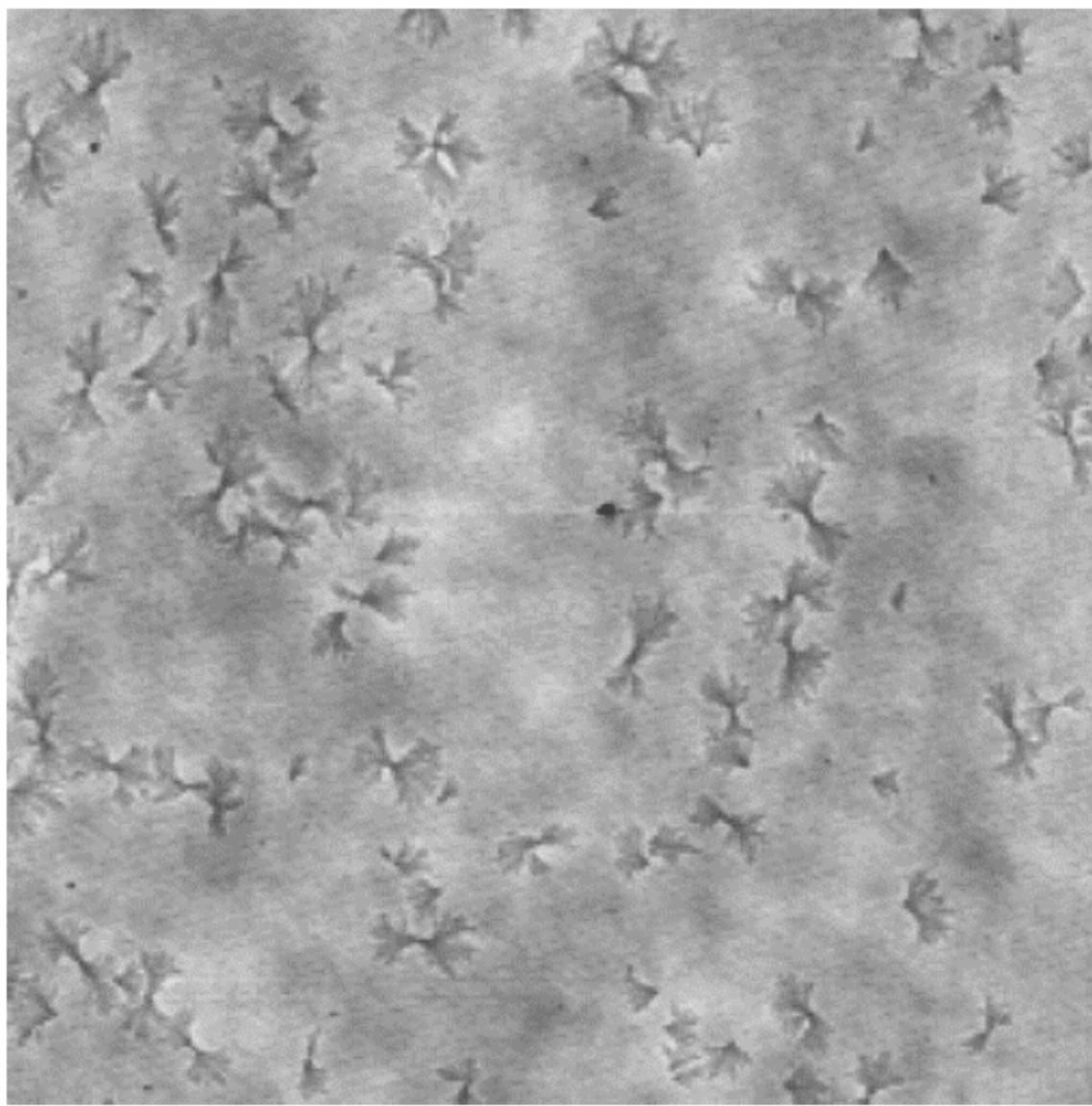


P.J. Flory, 1949



G.S.Y. Yeh, 1972

Amorphous components (AC)



15 nm
0 nm

PET / AFM: M Durell, J E Macdonald, D Trolley, Europhysics Letters (2002)

Crystal “grains”: crystallites of varying orientation



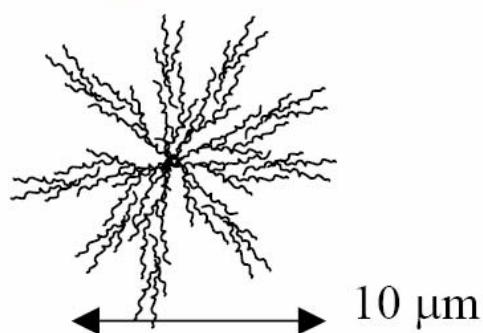
Ex: Stainless steels Fe-Ni-Cr

Depletes at grain boundaries causing corrosion

used for fracture fixation plates, etc., & angioplasty stents

Spherulites: radially oriented crystallites interspersed w/ amorphous phase

*semicrystalline polymers,
glass-ceramics*



Structure

Crystalline Amorphous components (CC)

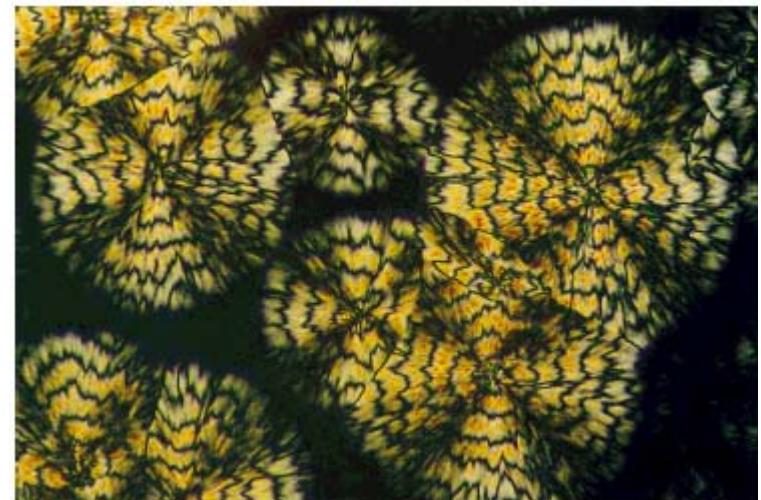
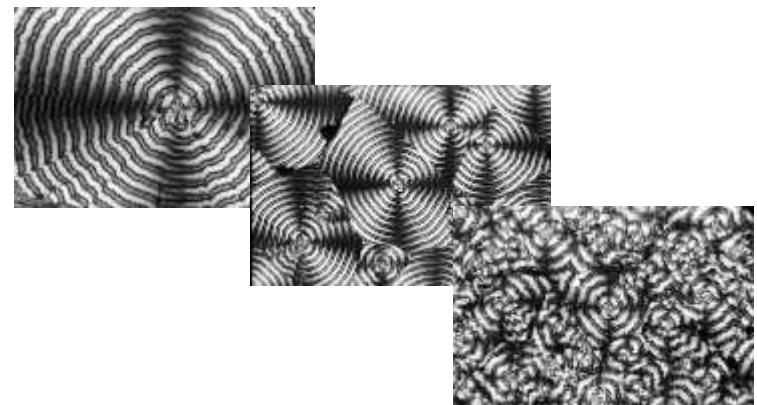
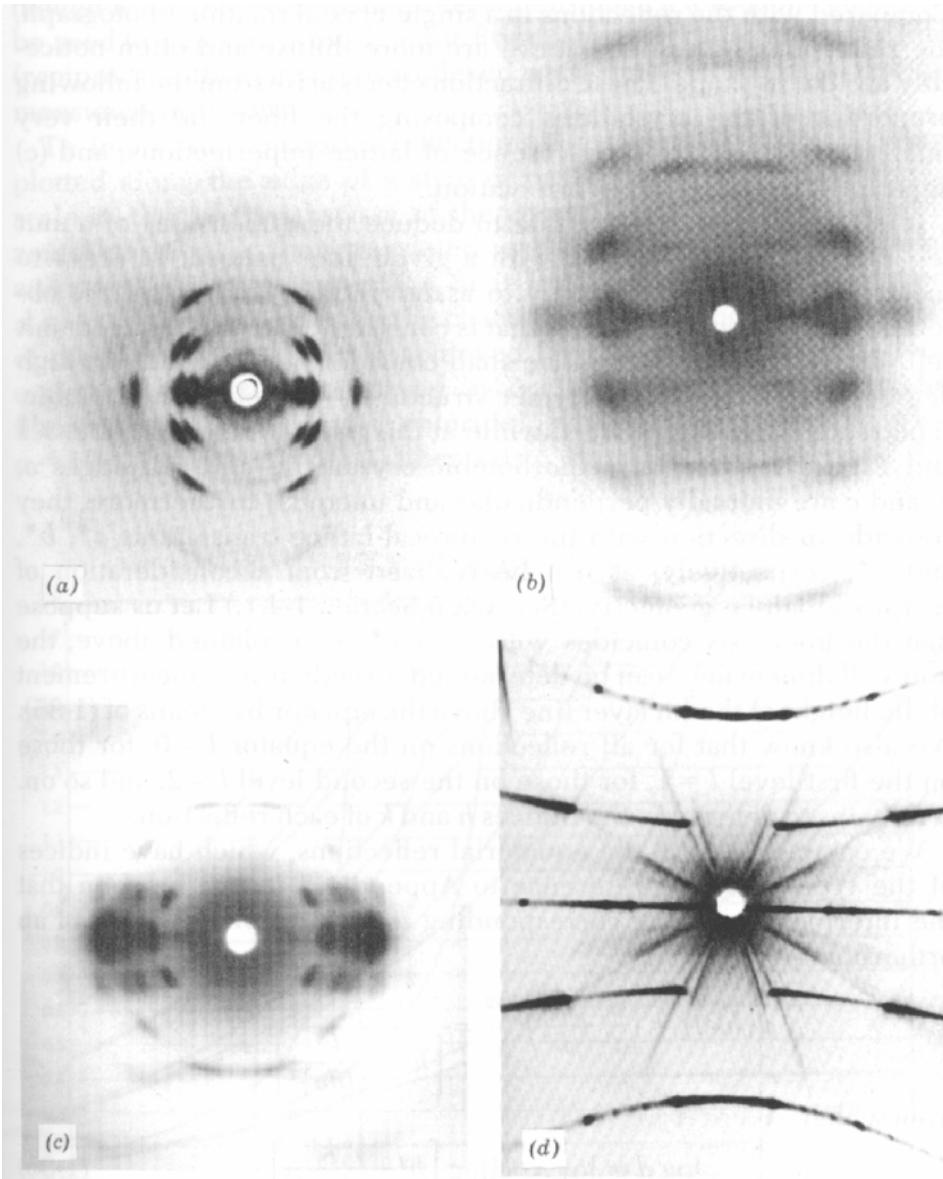
➤ Single

crystalline



A.Keller, 1953

Structure, scales, partially crystalline polymers

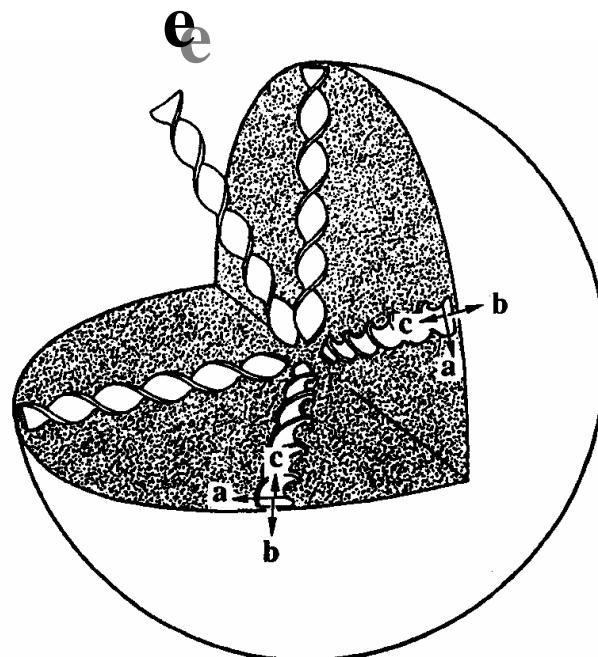


1 millimeter square, OM

Structure

Crystalline Amorphous components (CC)

➤ Pherulit



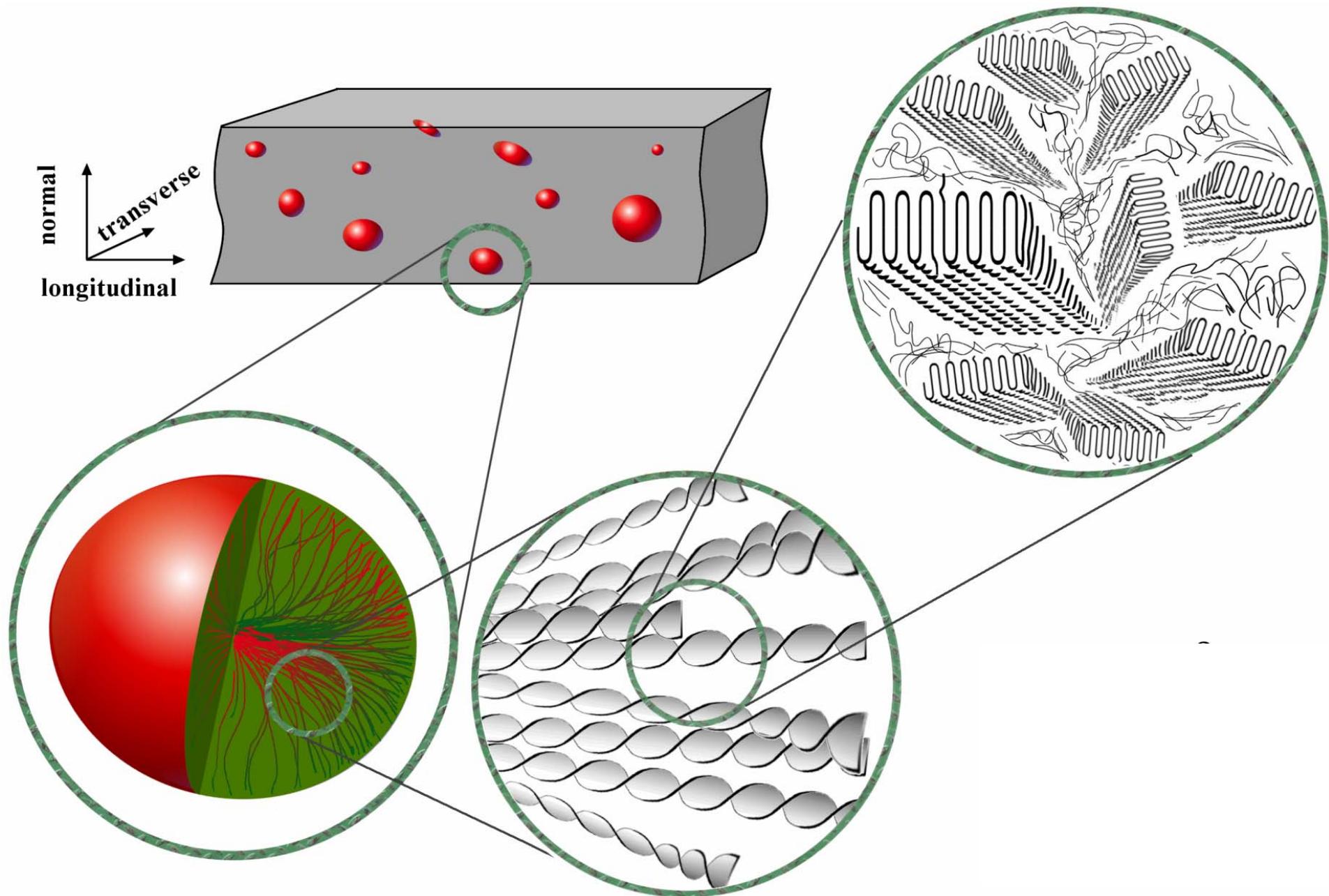
Schultz, J.M., 1984



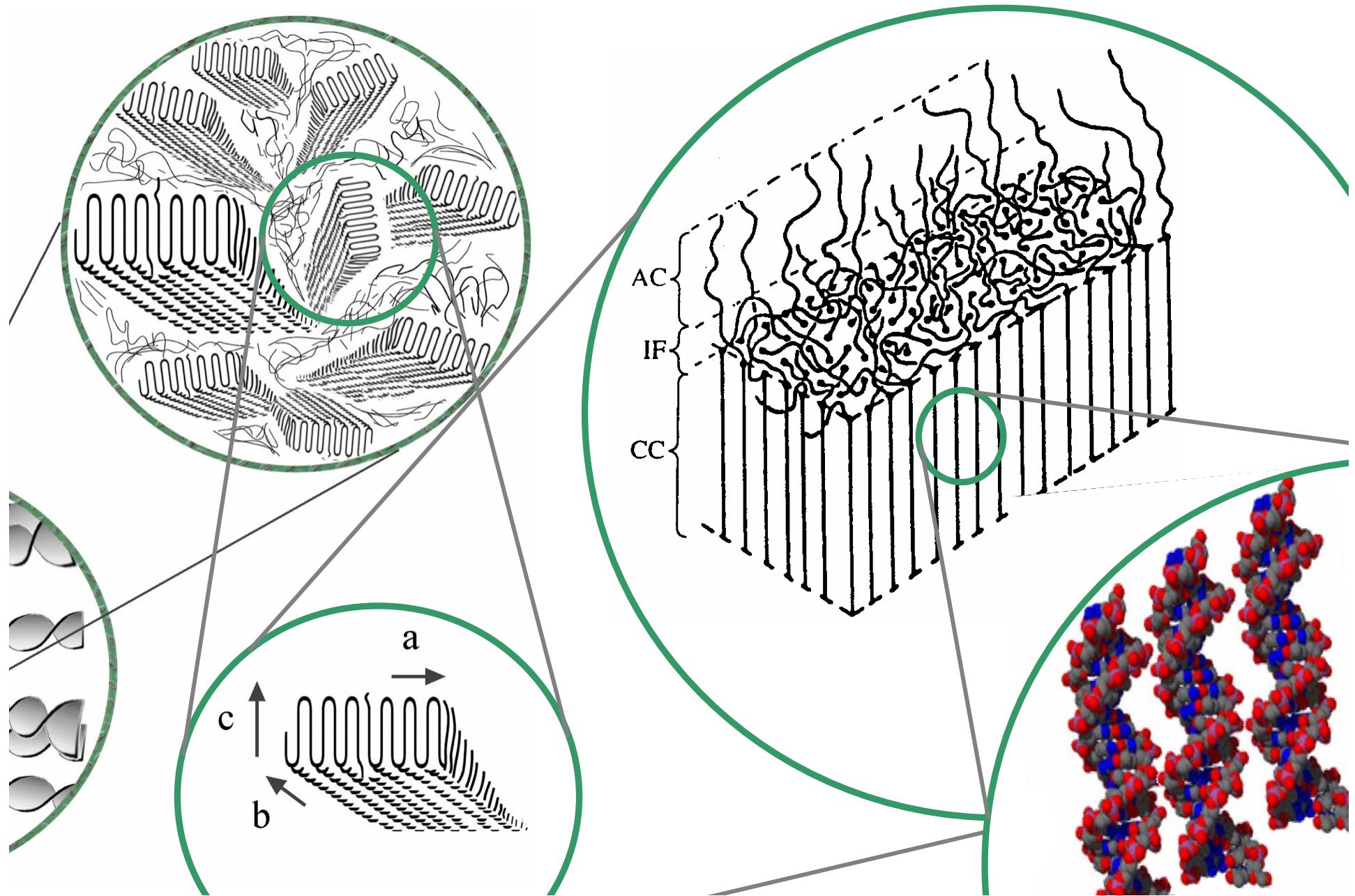
➤ Shish

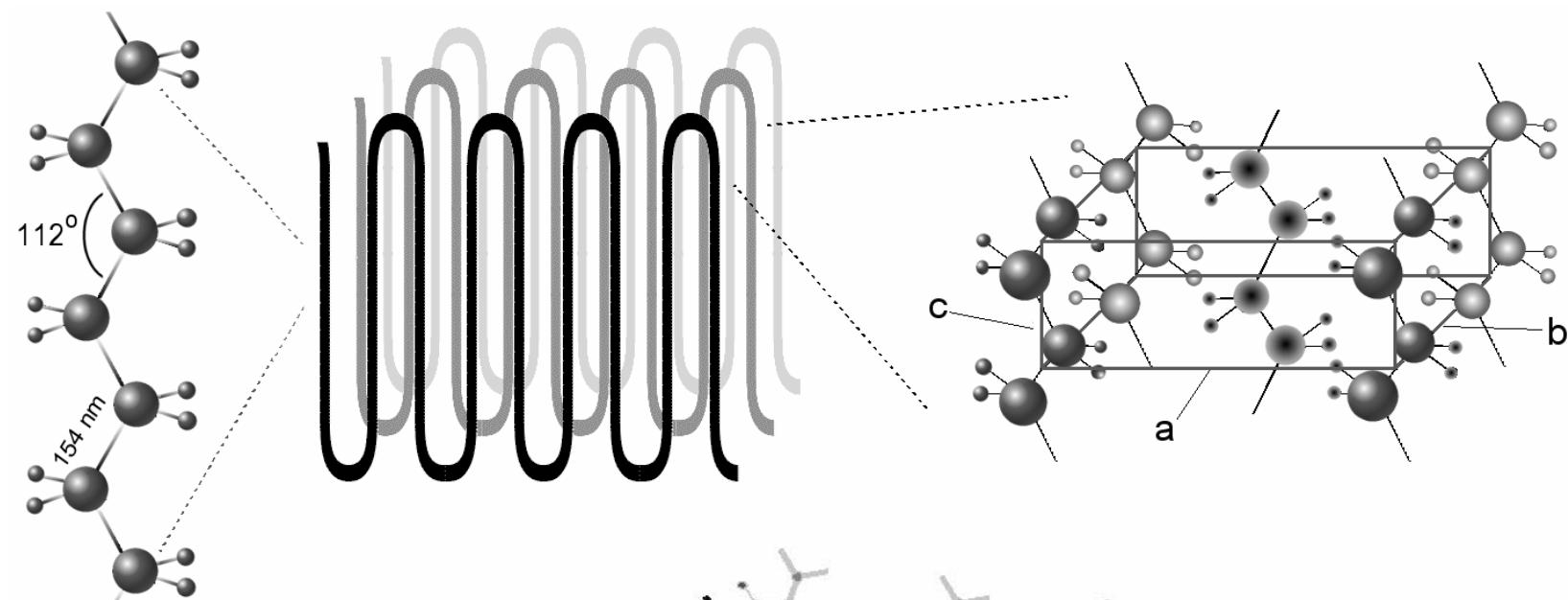
kebab

Structure, scales, partially crystalline polymers



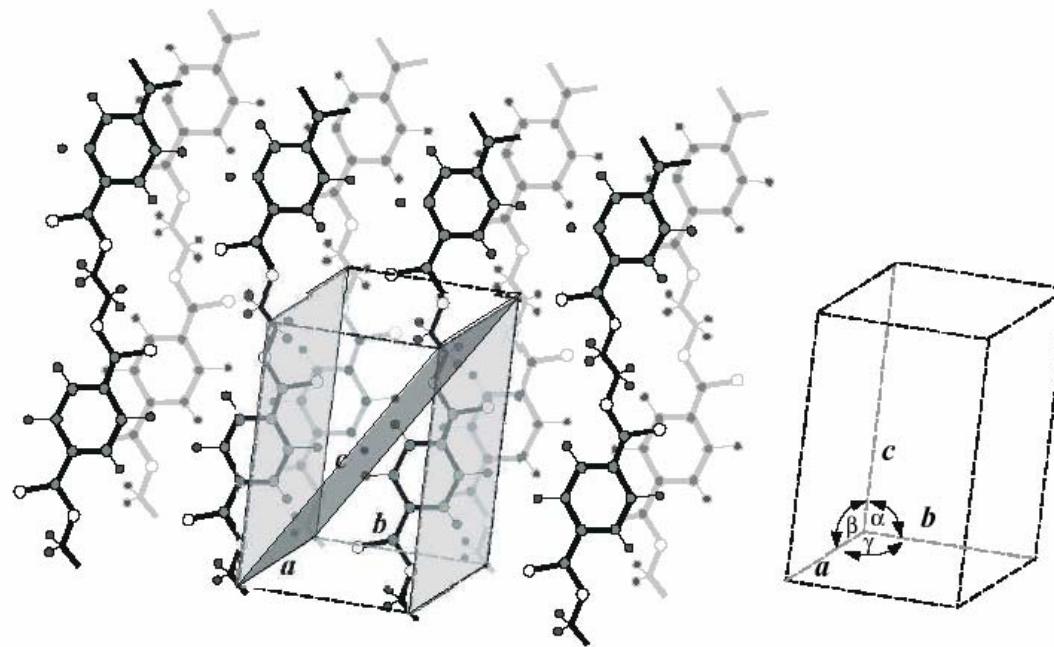
Structure, scales, partially crystalline polymers



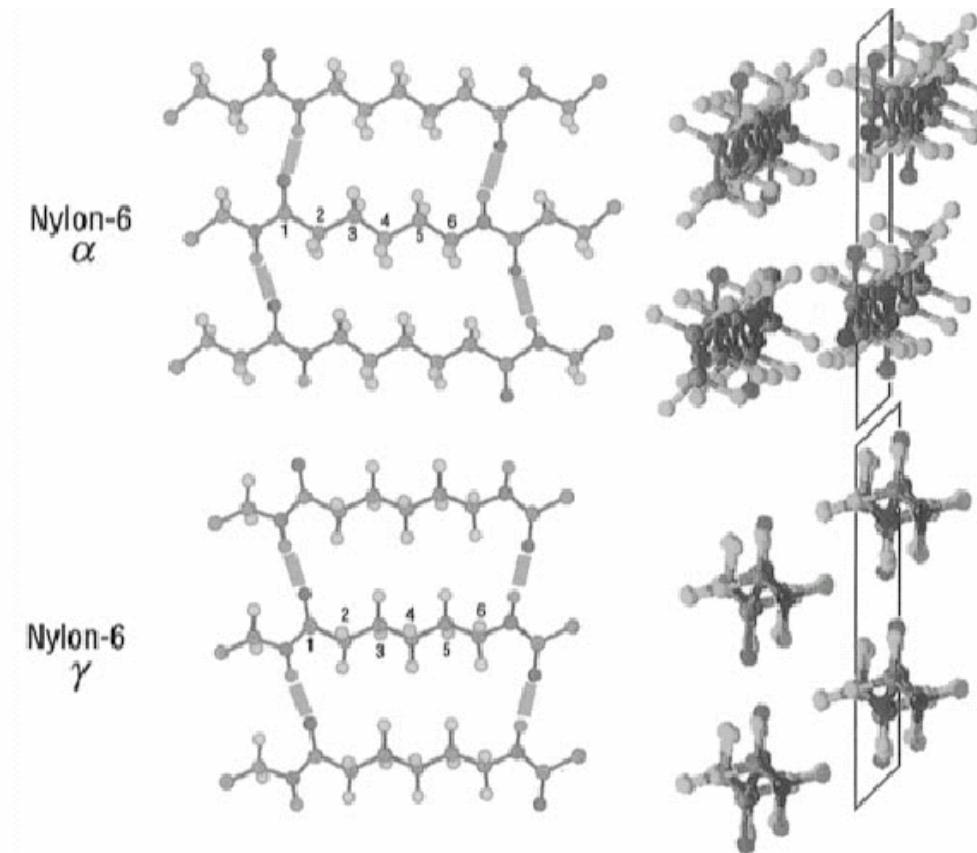


The crystalline unit cell for PET
 $(a = 4.56 \text{ \AA}, b = 5.94 \text{ \AA}, c = 10.75 \text{ \AA}; \alpha = 98.5^\circ, \beta = 118^\circ, \gamma = 112^\circ)$,

M Durell (2002)

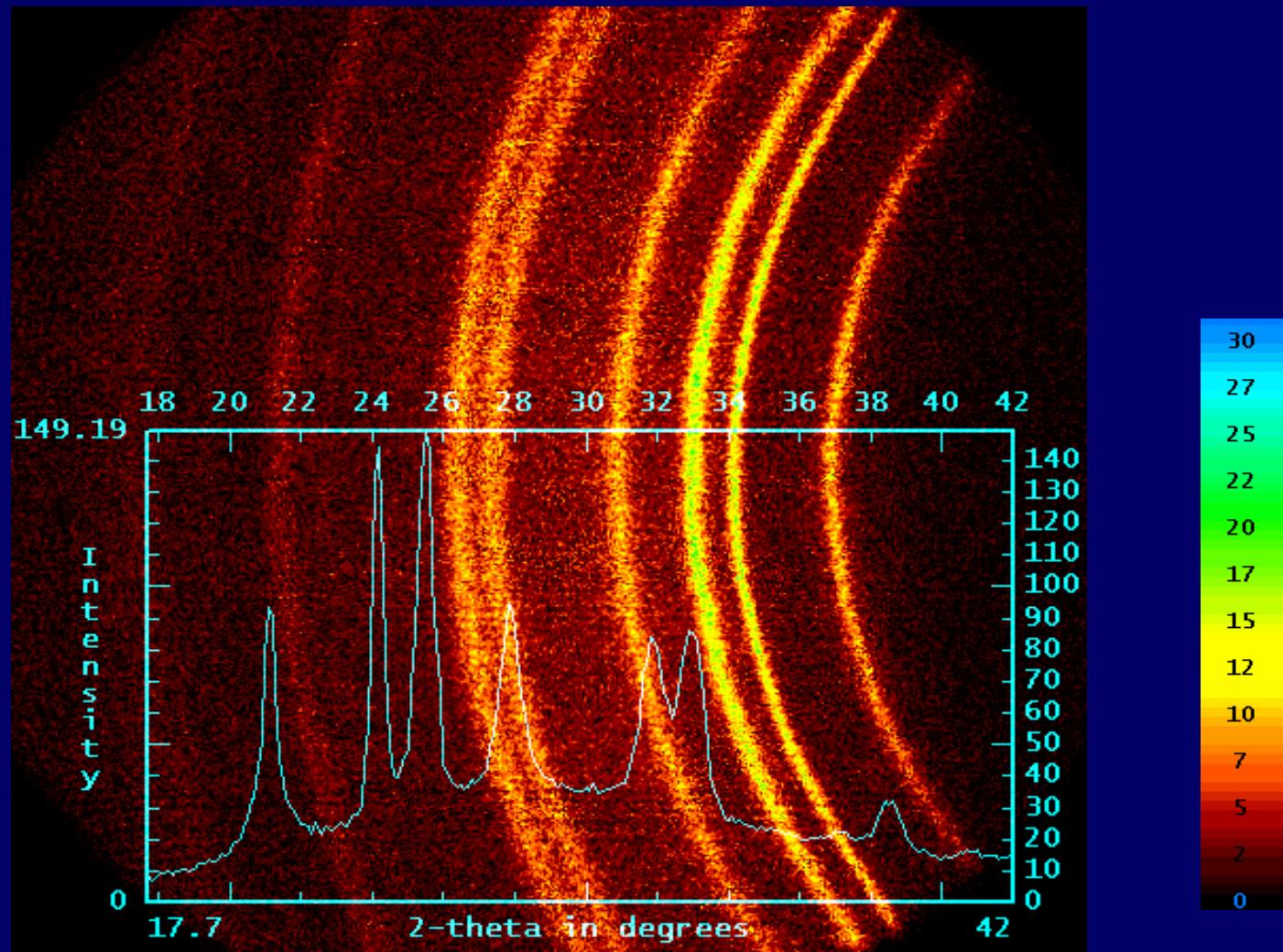


Nylon



Nylon exists in three forms, noncrystalline, α crystal and γ crystal. γ crystal is the most interested one. It is monoclinic, but could be more precisely described as pseudo-orthorhombic lattice with $a=0.482\text{nm}$, $b= 0.782\text{nm}$ and $c = 1.67\text{nm}$. And α crystal is monoclinic lattice with $a=0.478\text{nm}$, $b= 0.400\text{nm}$, $c=1.724\text{nm}$ and angle $\gamma = 67.5^\circ$.

PP



Polypropylen (PP), Cr-K α 1, 40KV 40mA

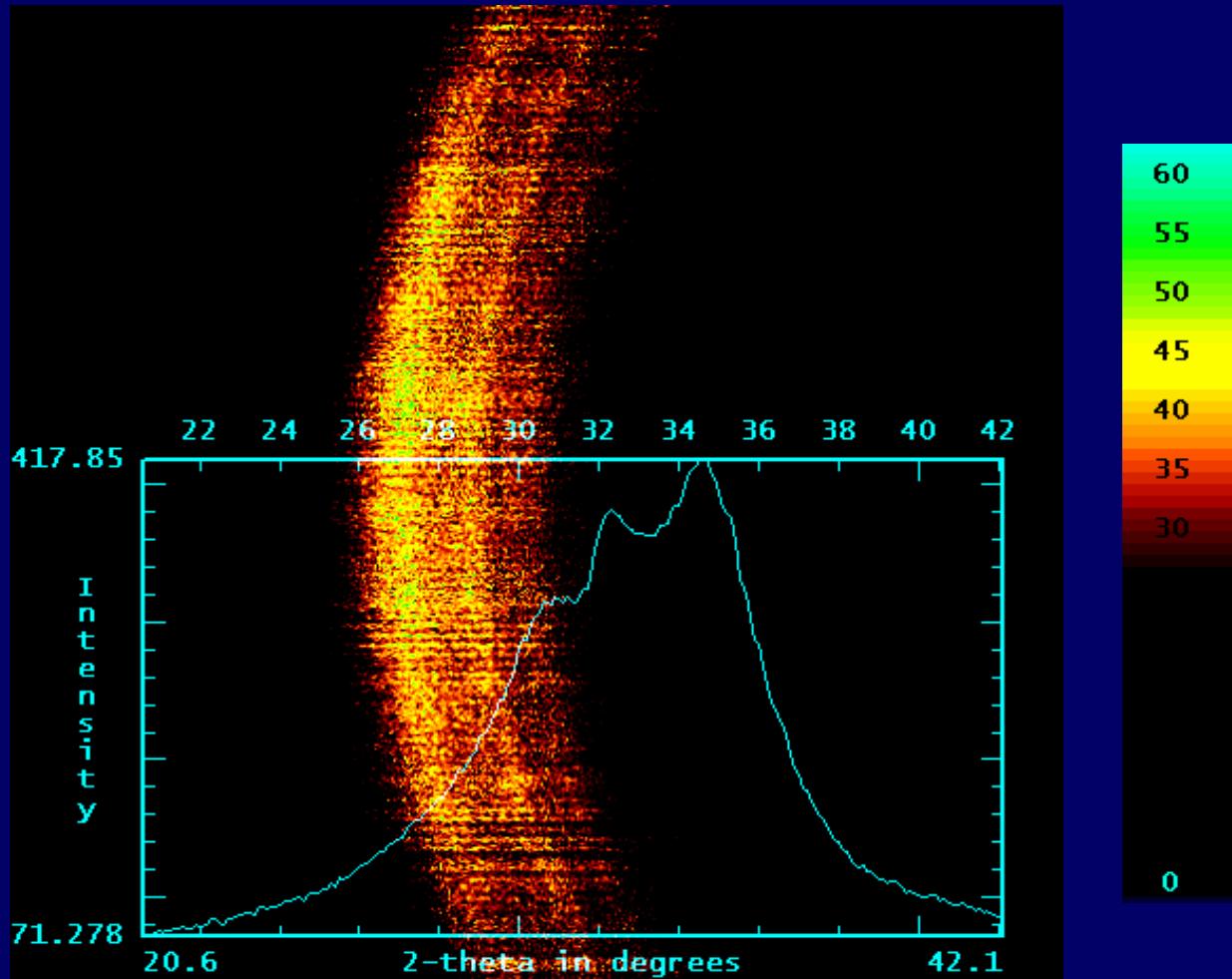
Nylon

α

4,489	(-131)	29,6
	(13-1)	29,6
4,416	(200)	30,1
4,278	(210)	31,1
	(2-10)	31,1
4,219	(-12-1)	31,54
	(121)	31,54
4,098	(-221)	32,5
	(22-1)	32,5
4,007	(-102)	33,26
3,93	(-220)	33,92
	(220)	33,92
3,903	(-1-12)	34,18
	(-112)	34,18
3,705	(002)	36,06

γ

4,599	(11-1)	28,88
	(-111)	28,88
4,159	(-121)	32,02
	(12-1)	32,02
4,109	(-102)	32,42
4,097	(100)	32,5
3,999	(002)	33,34
3,992	(-1-12)	33,38
	(-112)	33,38
3,982	(110)	33,48
	(-110)	33,48
3,891	(0-12)	34,28
	(012)	34,28
3,694	(12-2)	36,18



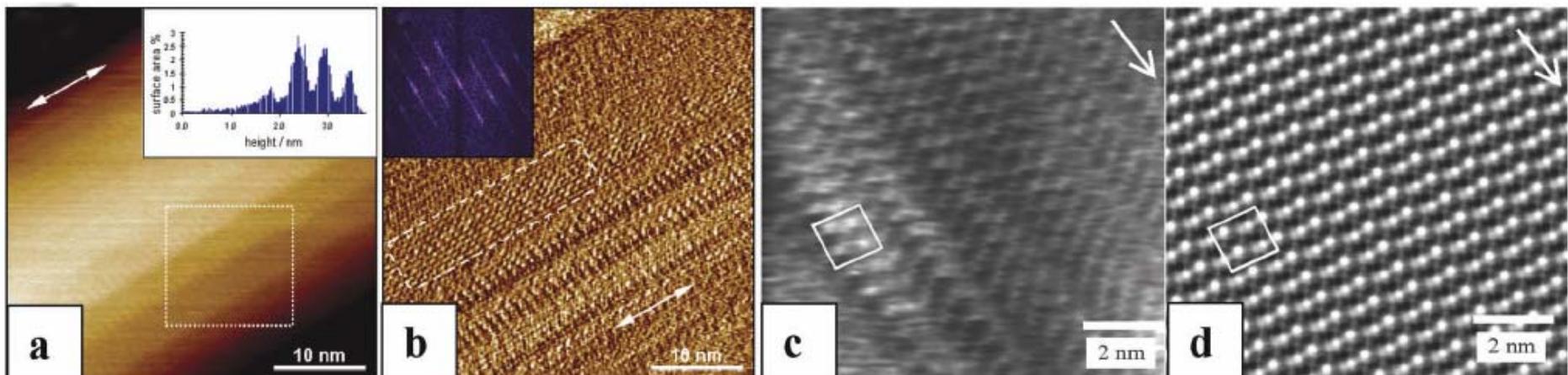
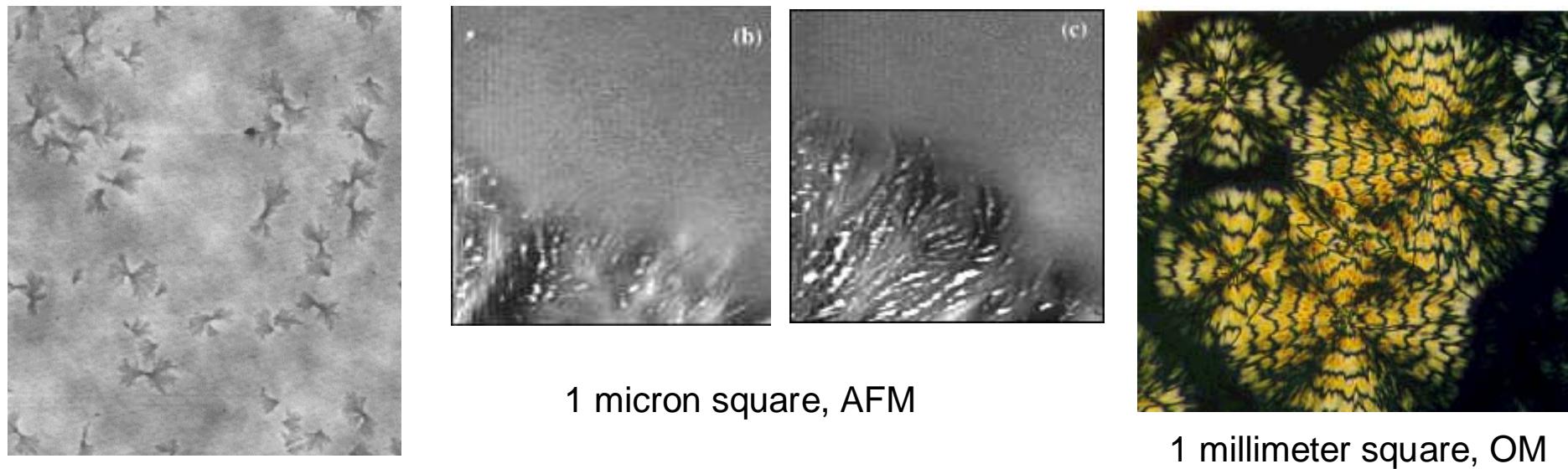
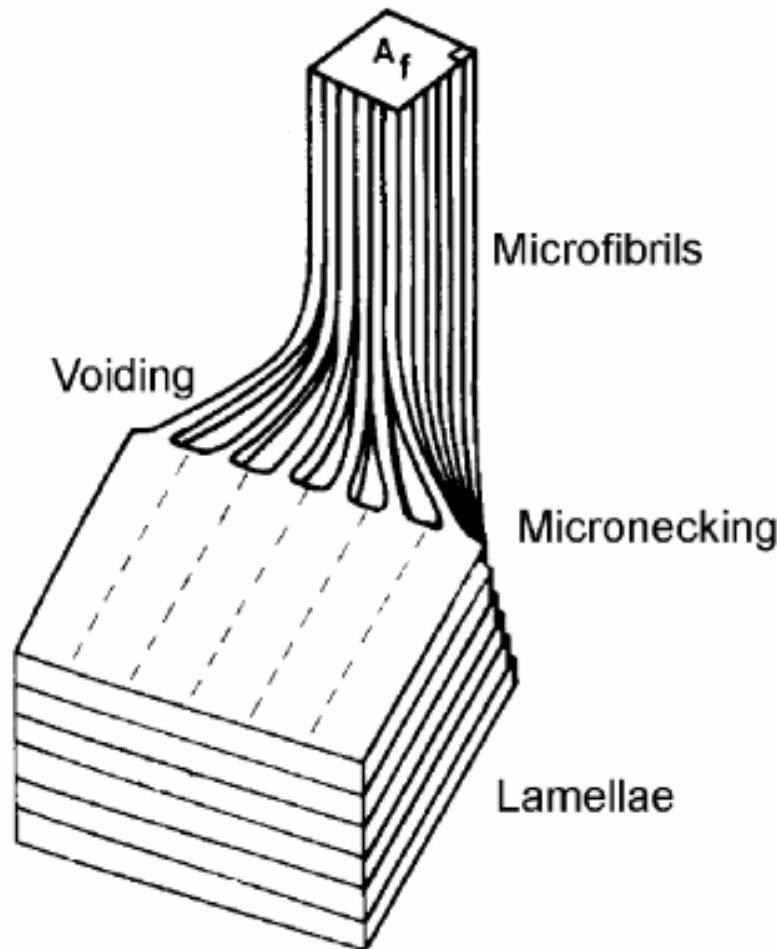


Fig. 2 AFM contact mode images. Cellulose molecules in micro-fibril crystal: (a) topographic image; and (b) error-signal image. Polybutene-1 molecules: (c) topographic image; and (d) simulated AFM image based on crystallographic data.

M. Miles, M. Antognazzi, H. Haschke, J. Hobbs, A. Humphris, T. McMaster
 H.H. Wills Physics Laboratory, University of Bristol, Materials Today, Feb. 2003

deformation

Tensile and drawing



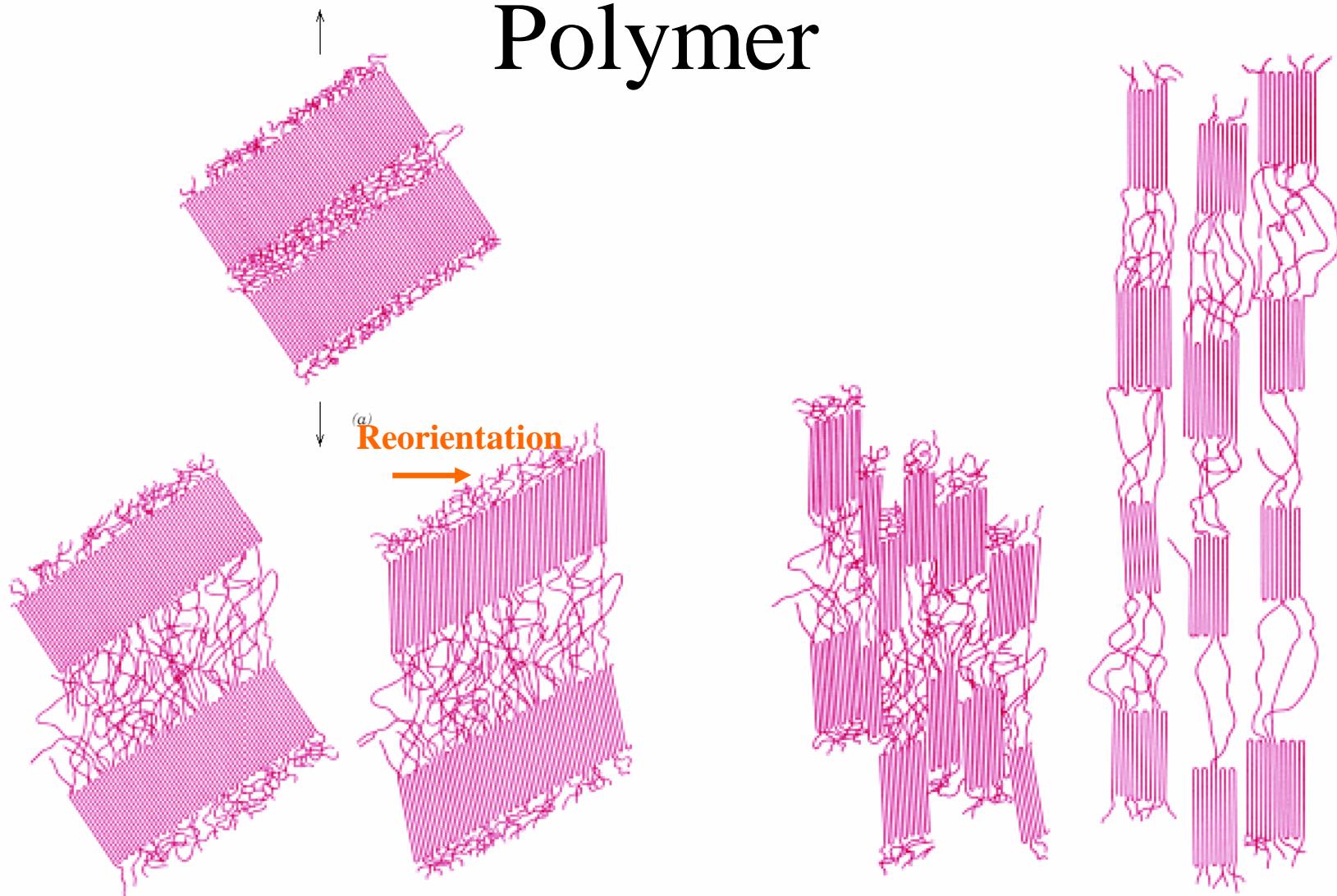
Schematic model of a cavitating semi-crystalline polymer under tensile deformation

A.Galeski, Progress in Polymer Science, (2003)

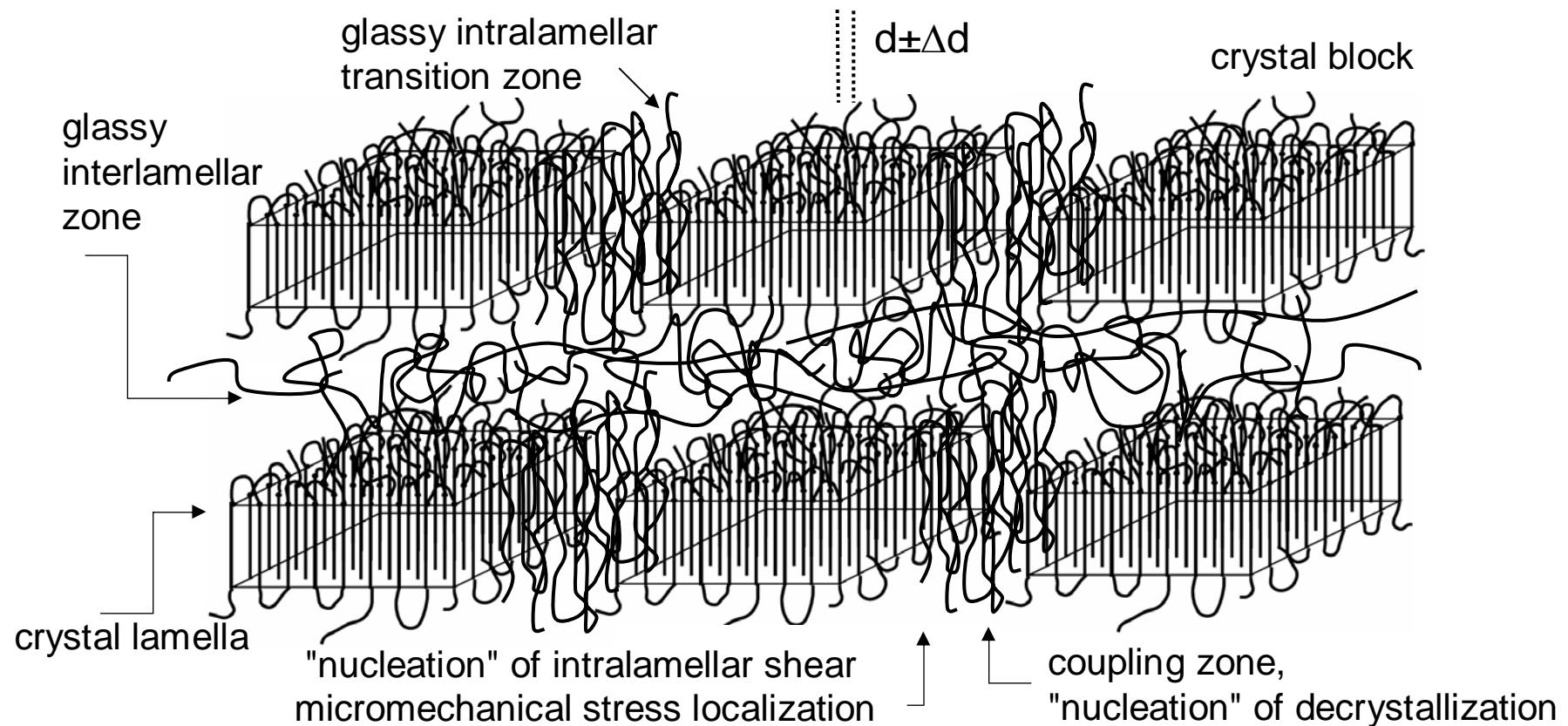
Peterlin: internal cavitation-micronecking

A. Peterlin, Journal of Materials Science , (1971)

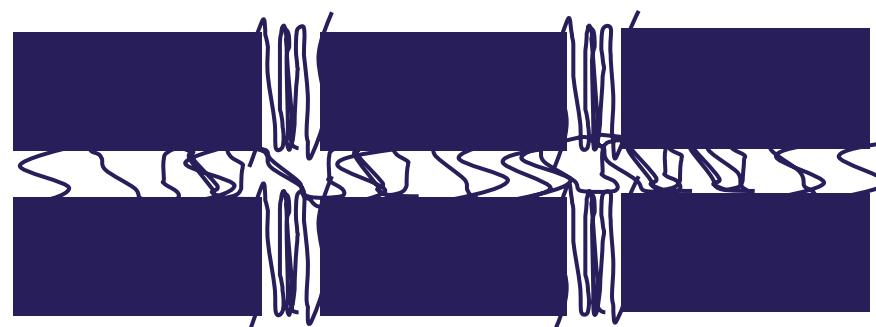
Polymer



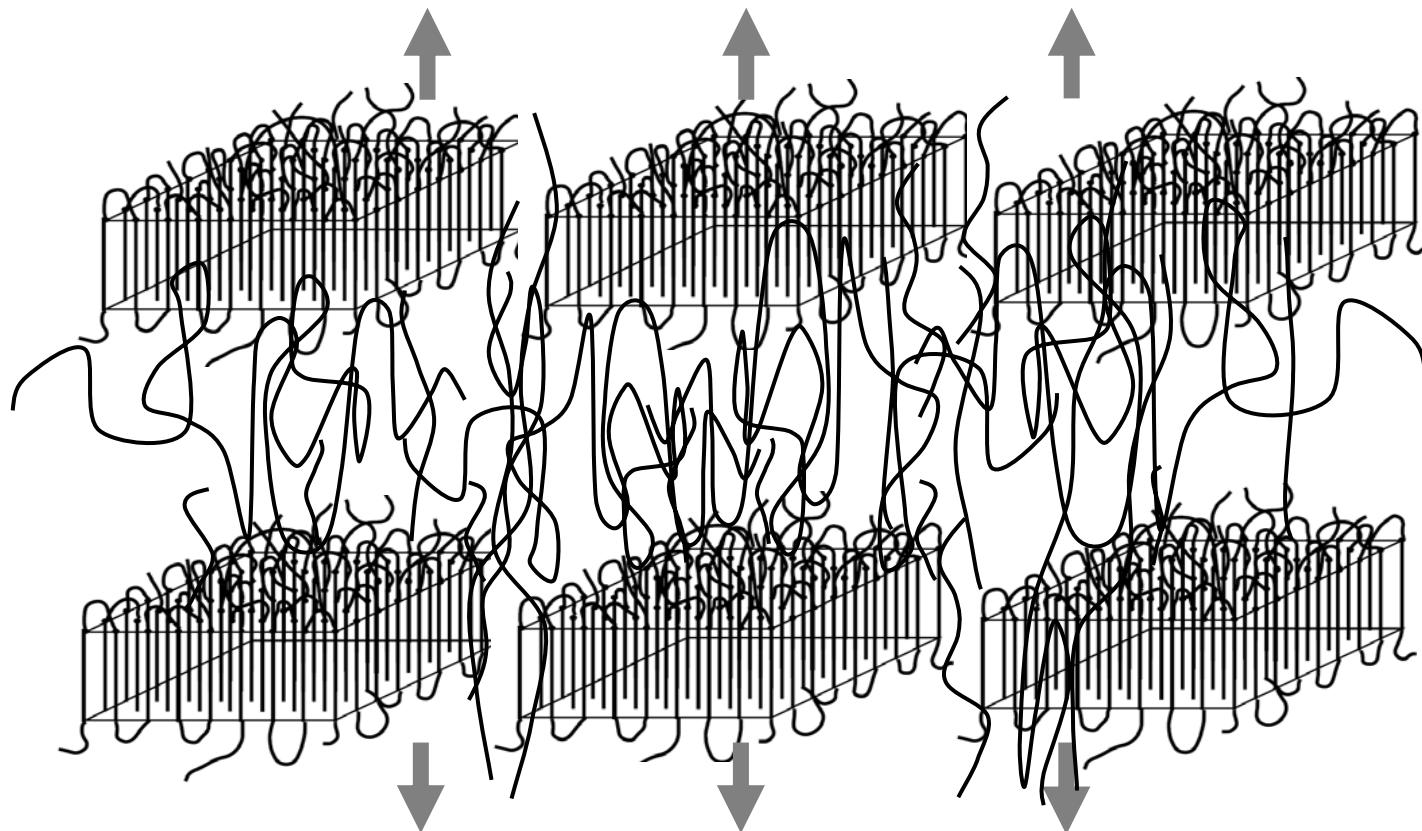
texture-relevant lattice defects, a tensorial view



**tensorial
nature
of polymer
deformation**

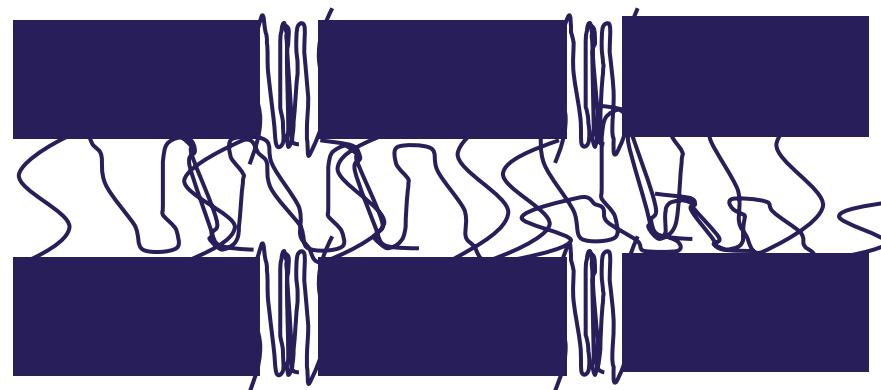


orientation dependence, elongation of soft phase

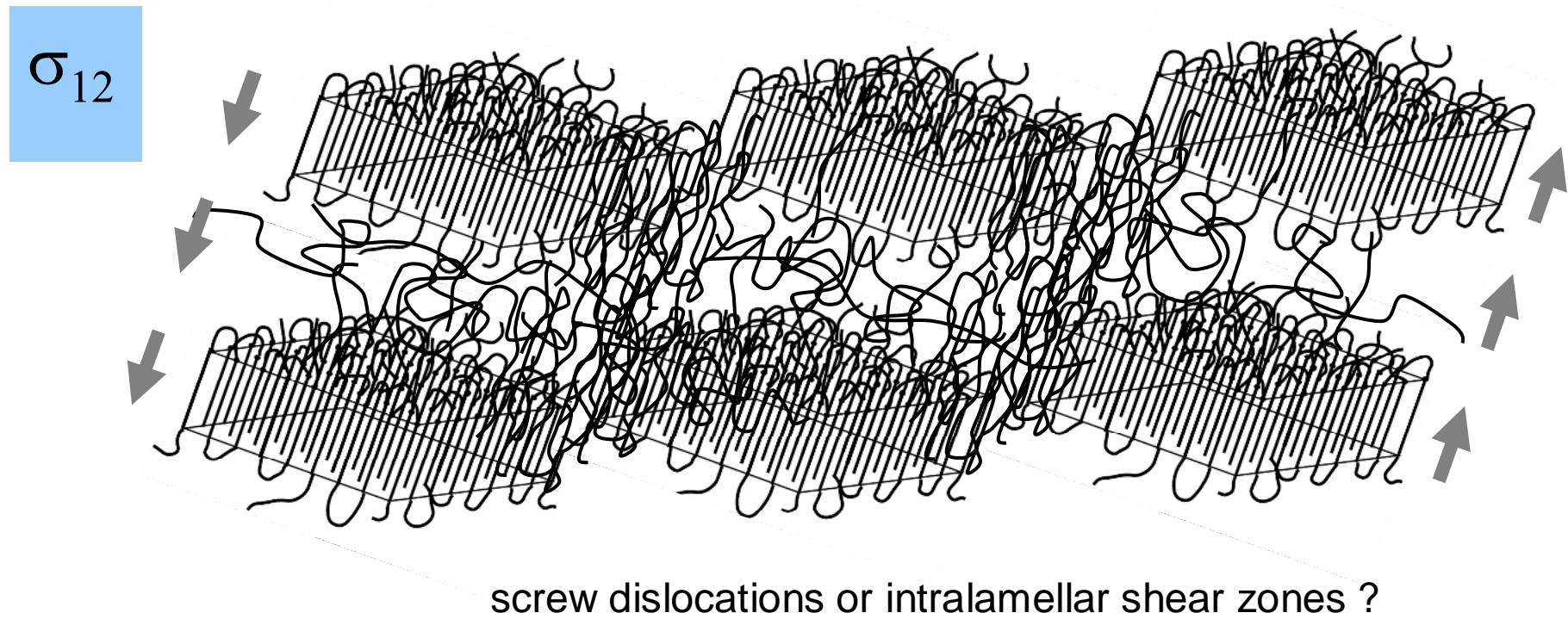


tensorial
nature
of polymer
deformation

$$\sigma_{11}$$



orientation dependence, shear of crystal phase #1



**tensorial
nature
of polymer
deformation**

$$\sigma_{12}$$

