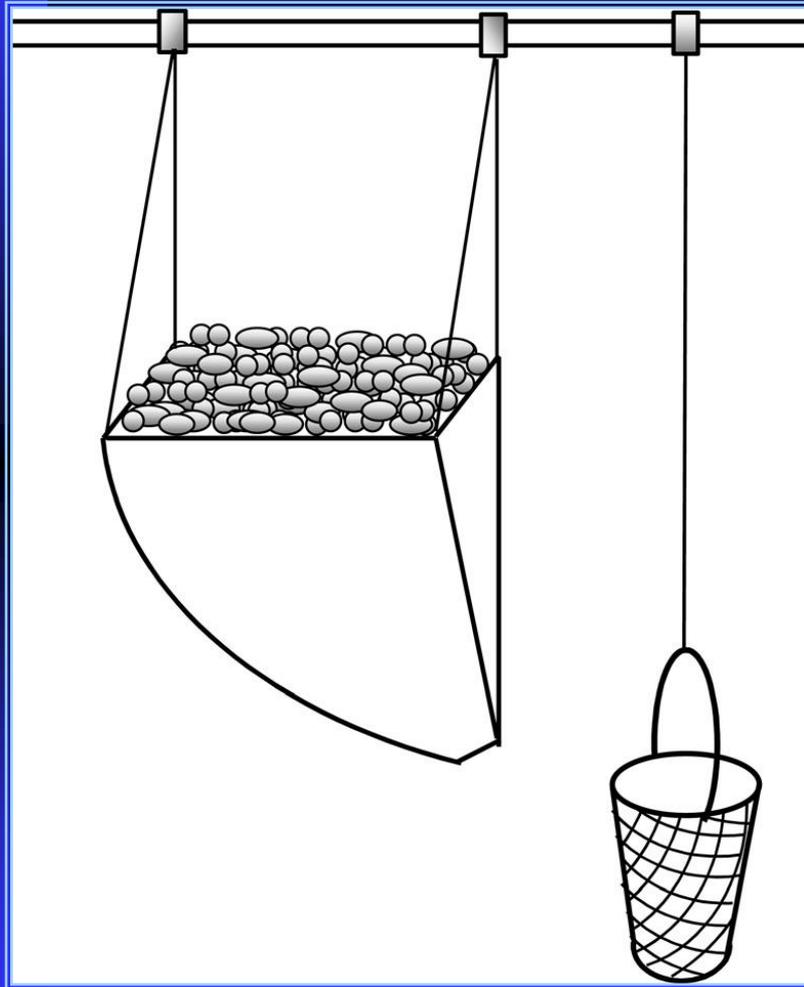


# Elasticity

- Elastic deformation (springs, spring washers)
- Structure rigidity (undesirable flexure, change of geometry)
- Structure vibrations dumping (eigenfrequency)

R. Hook: „ut tensio, sic vis“ (1676)

# Elasticity



- Leonardo da Vinci tensile test of wires
- 15<sup>th</sup> century !



## Tacoma Bridge Collapse



- 1) Elastic (Young's) modulus
- 2) Bonds between atoms
- 3) Atoms ordering in solids
- 4) Physical background of Youngs modulus

## Inelasticity/internal dumping

- 1) Parameters introduction and measurement
- 2) Physical and practical meaning

# Young's modulus

- ❑ Elastic characteristics of materials
- ❑ Measurements of elastic characteristics
- ❑ Values of Young's modulus

# Generalized Hooke's law

**Anisotropic material** (Anisotropy = dependence of physical properties of materials on the direction in which this property is determined)

Strain tensor

$$\begin{Bmatrix} \varepsilon_x \\ \varepsilon_y \\ \varepsilon_z \\ \gamma_{yz} \\ \gamma_{zx} \\ \gamma_{xy} \end{Bmatrix} = \begin{Bmatrix} S_{11} & S_{12} & S_{13} & S_{14} & S_{15} & S_{16} \\ S_{21} & S_{22} & S_{23} & S_{24} & S_{25} & S_{26} \\ S_{31} & S_{32} & S_{33} & S_{34} & S_{35} & S_{36} \\ S_{41} & S_{42} & S_{43} & S_{44} & S_{45} & S_{46} \\ S_{51} & S_{52} & S_{53} & S_{54} & S_{55} & S_{56} \\ S_{61} & S_{62} & S_{63} & S_{64} & S_{65} & S_{66} \end{Bmatrix} * \begin{Bmatrix} \sigma_x \\ \sigma_y \\ \sigma_z \\ \tau_{yz} \\ \tau_{zx} \\ \tau_{xy} \end{Bmatrix}$$

Elastic coefficients

Stress tensor

The matrix contains 36 elements, however, thanks to indexes symmetry only 21 are independent

# Generalized Hooke's law

(by introducing axis x,  
dependent elastic

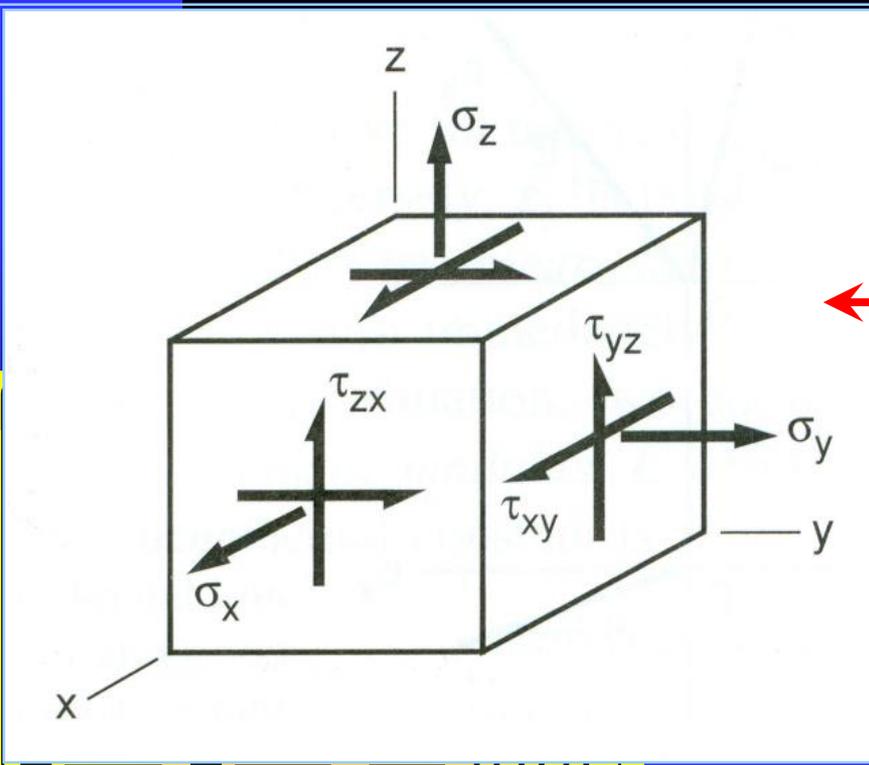
← **Orthotropic material**

Orthotropy = right angle anisotropy

Monocrystal, texture

Composite – matrix + fiber

(axis „z“ is neglected)

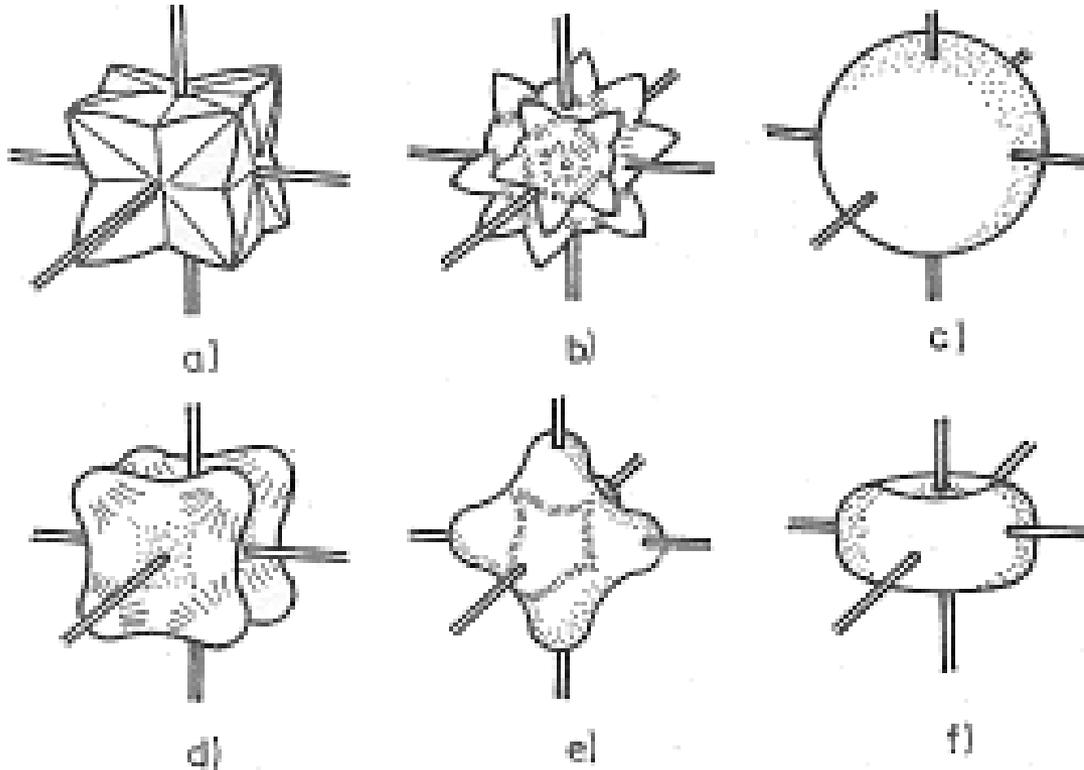


$$\left\{ \begin{array}{l} E_x, E_y, E_z, 0, 0, 0 \\ 0, 0, 0, \frac{1}{G_{yz}}, 0, 0 \\ 0, 0, 0, 0, \frac{1}{G_{zx}}, 0 \\ 0, 0, 0, 0, 0, \frac{1}{G_{xy}} \end{array} \right\}$$

$$\left\{ \begin{array}{l} \epsilon_x \\ \epsilon_y \\ \gamma_{xy} \end{array} \right\} = \left\{ \begin{array}{l} \frac{1}{E_x}, -\frac{\mu_{yx}}{E_y}, 0 \\ -\frac{\mu_{xy}}{E_x}, \frac{1}{E_y}, 0 \\ 0, 0, \frac{1}{G_{xy}} \end{array} \right\} * \left\{ \begin{array}{l} \sigma_x \\ \sigma_y \\ \tau_{xy} \end{array} \right\}$$

# Generalized Hooke's law

## End points of Young's modulus vectors



## Orthotropic material

monocrystals

a) Al (fract. stress)

b) Al (ductility)

c) Al (E)

d) Fe (E)

e) Fe (G)

f) Mg (E)

# Isotropic material

# Generalized Hooke's law

Isotropy = properties are independent of direction  
(contrary to anisotropy)

**Polycrystalline materials – metals,  
ceramics, polymers, partly also**

**... amorphous materials – glass**

**only two independent  
characteristics**

$$\mathbf{G} = \frac{E}{2(1 + \mu)}$$

$$\left\{ \begin{array}{l} \frac{1}{E}, -\frac{\mu}{E}, -\frac{\mu}{E}, 0, 0, 0 \\ -\frac{\mu}{E}, \frac{1}{E}, -\frac{\mu}{E}, 0, 0, 0 \\ -\frac{\mu}{E}, -\frac{\mu}{E}, \frac{1}{E}, 0, 0, 0 \\ 0, 0, 0, \frac{1}{\mathbf{G}}, 0, 0 \\ 0, 0, 0, 0, \frac{1}{\mathbf{G}}, 0 \\ 0, 0, 0, 0, 0, \frac{1}{\mathbf{G}} \end{array} \right\}$$

$$\varepsilon_x = \frac{1}{E} [\sigma_x - \mu(\sigma_y + \sigma_z)]$$

$$\varepsilon_y = \frac{1}{E} [\sigma_y - \mu(\sigma_x + \sigma_z)]$$

$$\varepsilon_z = \frac{1}{E} [\sigma_z - \mu(\sigma_x + \sigma_y)]$$

$$\gamma_{xy} = \frac{\tau_{xy}}{\mathbf{G}}$$

$$\gamma_{yz} = \frac{\tau_{yz}}{\mathbf{G}}$$

$$\gamma_{zx} = \frac{\tau_{zx}}{\mathbf{G}}$$

# Additional elastic characteristics, only two of them are independent

$$\sigma = E\varepsilon$$

$$\tau = G\gamma$$

$$p = -K \frac{\Delta V}{V}$$

$\mu$

standard form of Hooke's law for isotropic material loaded in tension and shear

modulus of elasticity of volume

Poisson ratio

**Important equations and quantities**

**Additional elastic characteristics, only two  
of them are independent**

$$E = 2G(1 + \mu)$$

$$K = \frac{E}{3(1 - 2\mu)}$$

$$\mu = \frac{E}{2G} - 1$$



**Important equations and quantities**

**Additional elastic characteristics, only two of them are independent**

**Isotropic material:  $\mu$  a  $E$**

**Poisson's ratio  $\mu$**  – ratio of relative transverse reduction in area to relative longitudinal elongation in field of elastic deformations.

The value  $\mu$  represents elastic compressibility of the body, i.e. ability of the material to decrease (at compression) and to increase (at tension) its volume in the course of elastic deformation.

# Definition of the $\mu$

relative transverse reduction in area

to

relative longitudinal elongation

$$\mu = \left| \frac{\Delta d / d_0}{\Delta l / l_0} \right| = - \frac{\Delta d / d_0}{\Delta l / l_0} = - \frac{\varepsilon_2}{\varepsilon_1} = - \frac{\varepsilon_3}{\varepsilon_1}$$

$$\varepsilon_3 = \varepsilon_2 = -\mu\varepsilon_1$$

# Definition of the $\mu$

What is the value of the  $\mu$  ?

$$(1 + \Delta V) = (1 + \varepsilon_1)(1 + \varepsilon_2)(1 + \varepsilon_3) =$$

$$(1 + \varepsilon_1)(1 - \mu\varepsilon_1)(1 - \mu\varepsilon_1) =$$

$$\left\{ \begin{array}{l} \text{neglecting the members in} \\ \text{second and third power } \varepsilon^2, \varepsilon^3 \end{array} \right\}$$

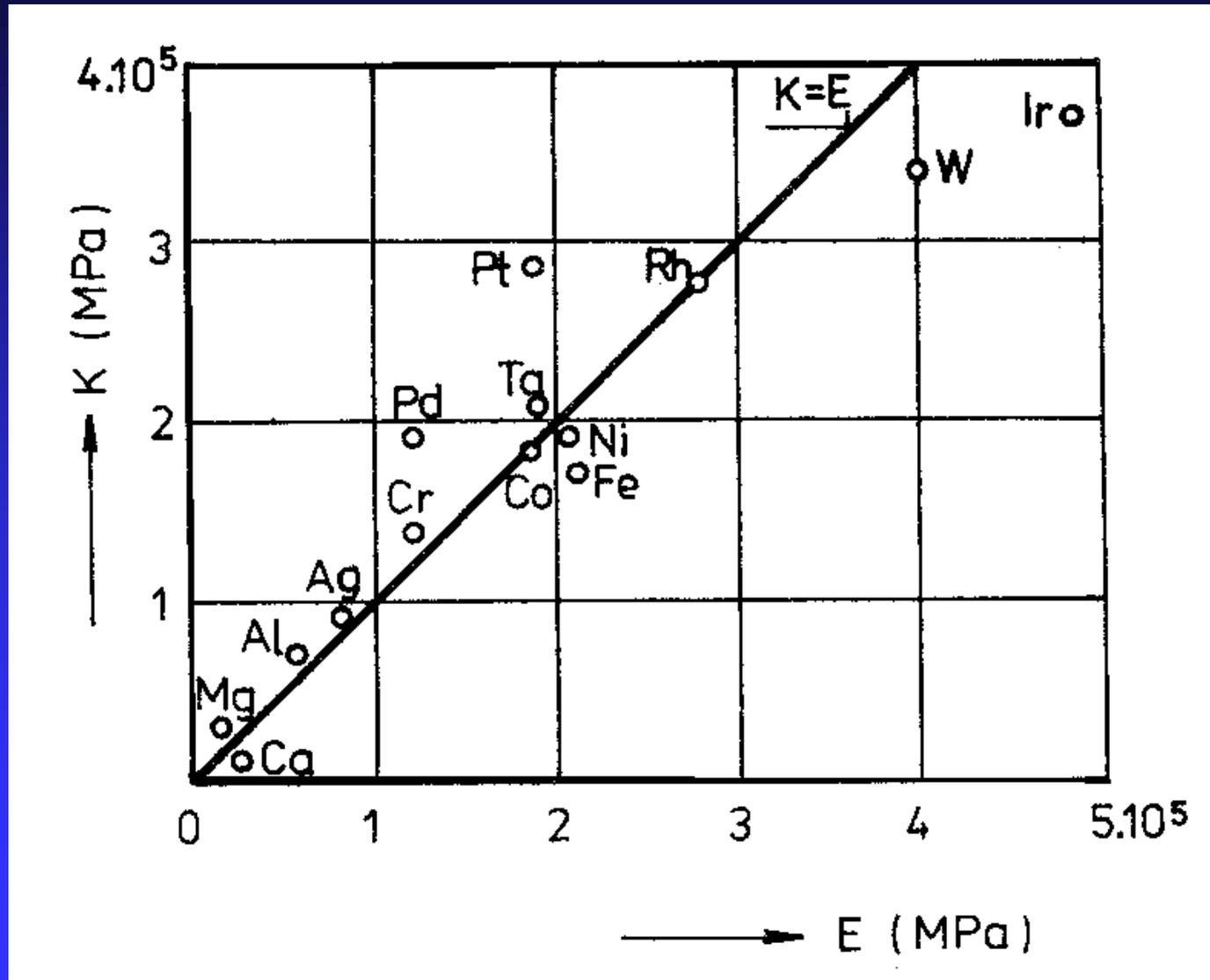
$$= 1 + \varepsilon_1(1 - 2\mu)$$

the volume increases  $\Rightarrow \Rightarrow 1 - 2\mu > 0 \Rightarrow \Rightarrow \mu < 0.5$

the volume is constant  $\mu = 0.5$

# Relation between $K$ and $E$

For metals,  $K$  is equal to  $E$



# What is the value of $\mu$ for metals ?

$$E \approx K$$

$$1 = 3(1 - 2\mu)$$

$$\mu = 0.333$$

$$G = \frac{3}{8} E$$

$$K = \frac{E}{3(1 - 2\mu)}$$

$$E = 2 * 10^5 \text{ MPa} = 200 \text{ GPa}$$
$$G = 7.5 * 10^4 \text{ MPa} = 75 \text{ GPa}$$

# Elastic characteristics - measurements

- Quasi-static methods
- Dynamic methods

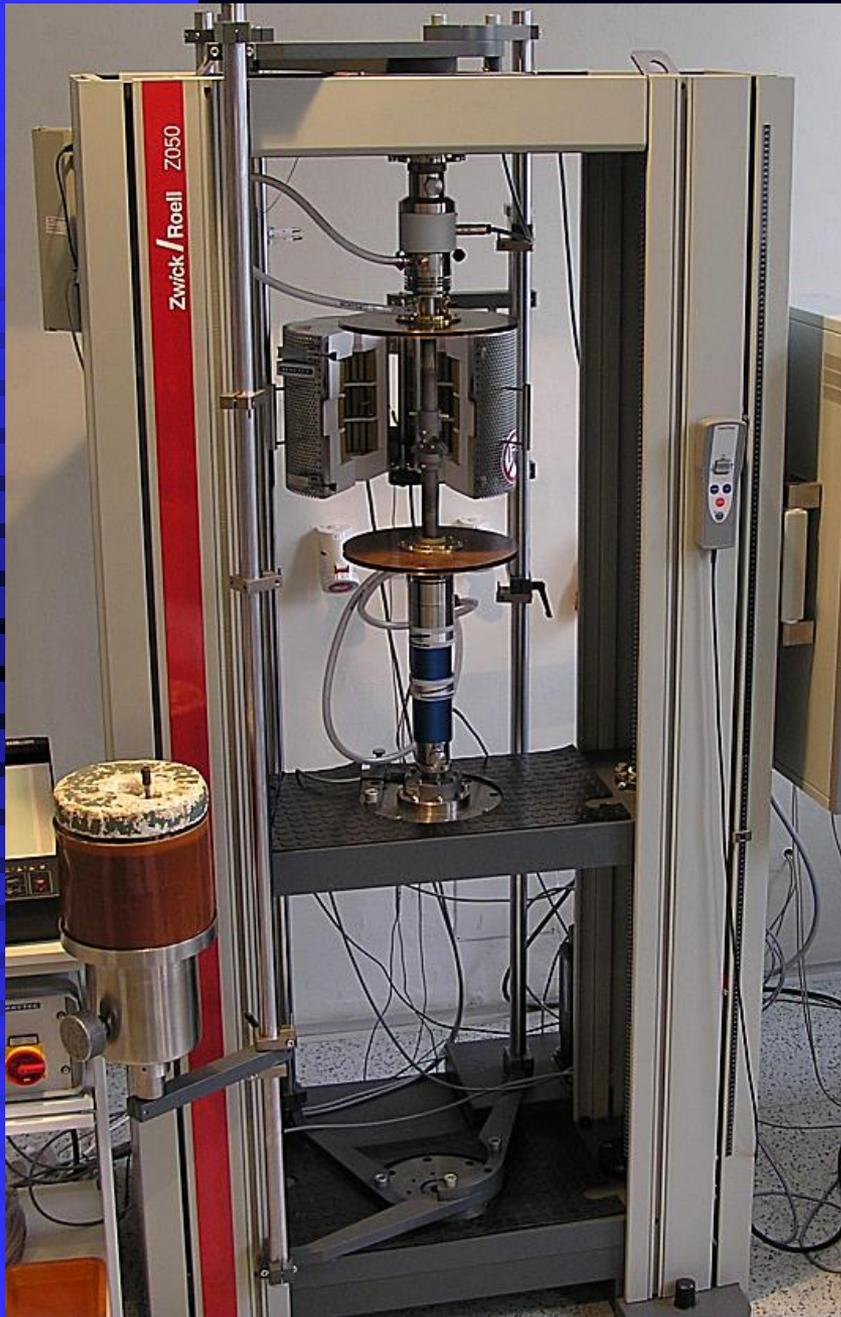
low frequency

high frequency

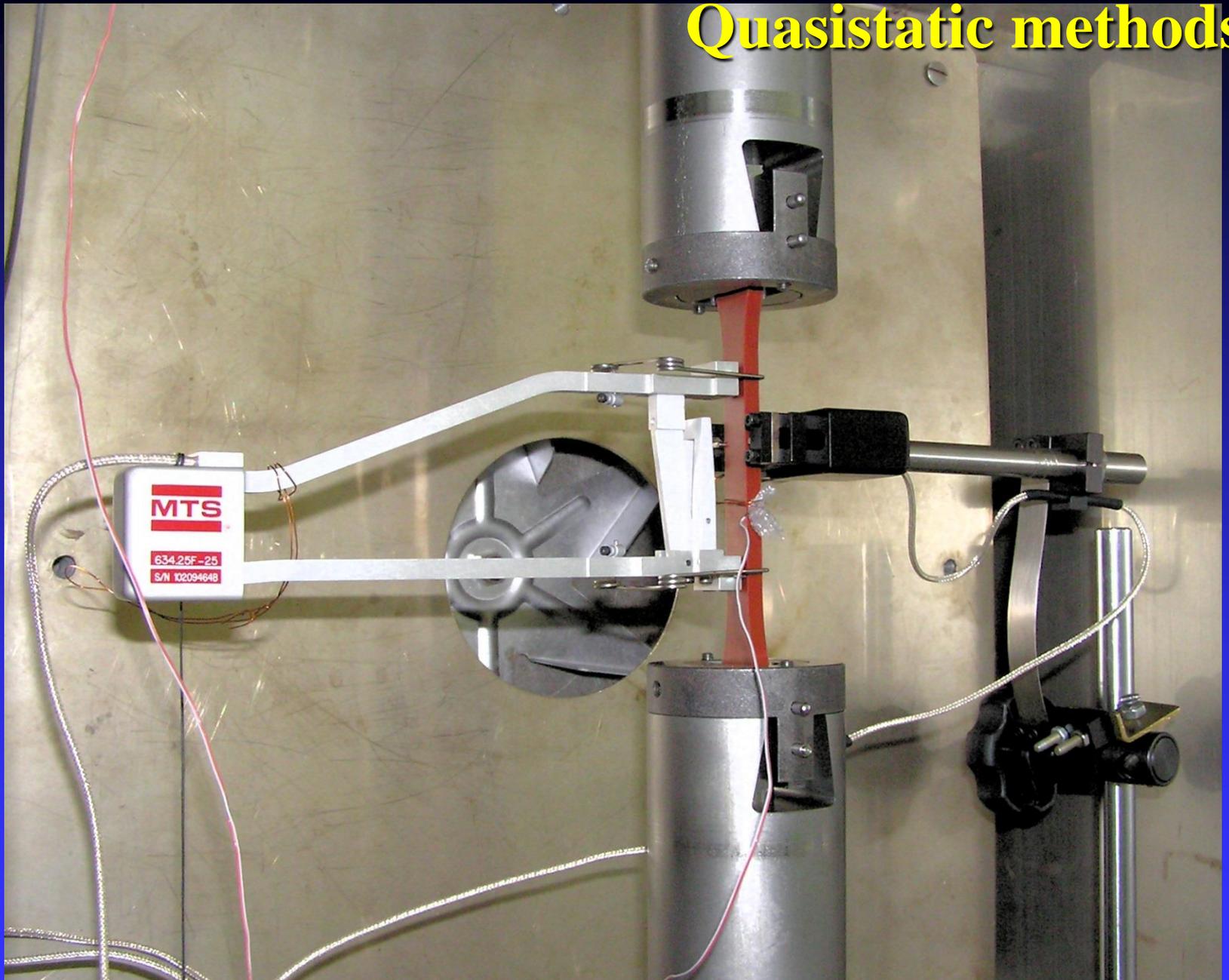
# Quasistatic methods

- Poisson ratio – from definition; using tensometers and/or transducers of longitudinal elongation and transversal contraction
- Modulus E – tensile test  
– flexural test
- Modulus G – torsion test

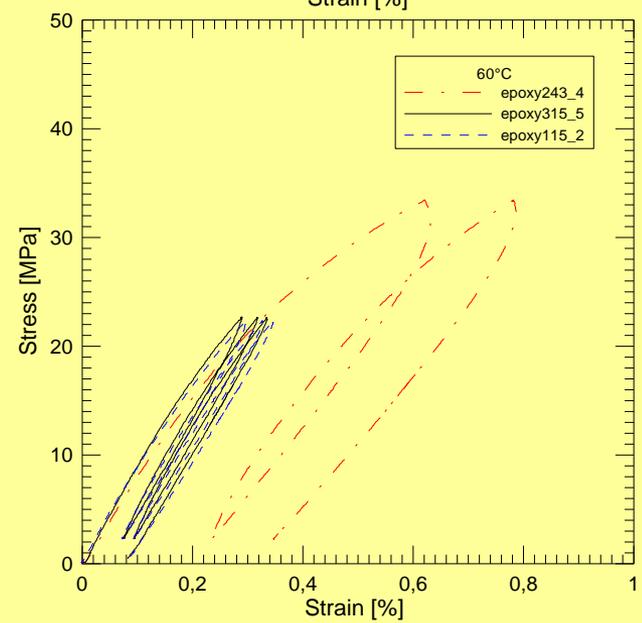
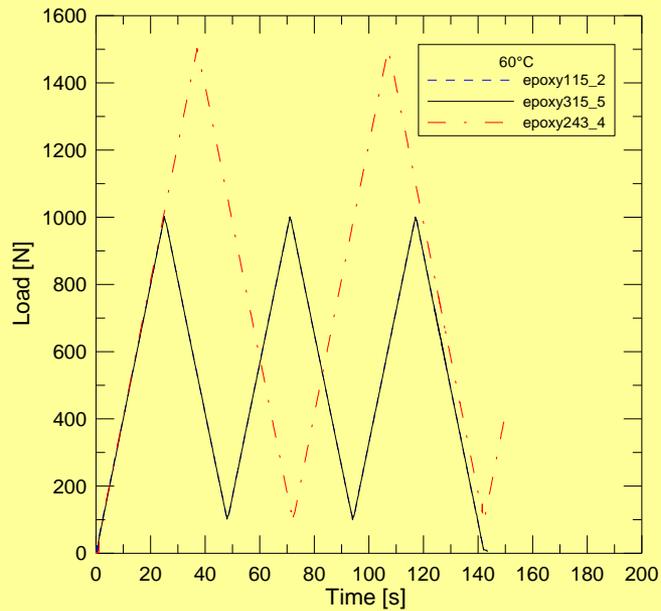
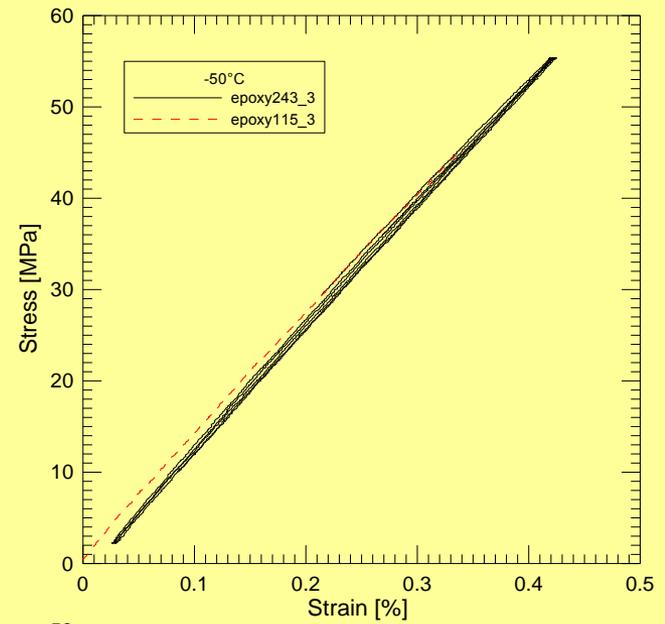
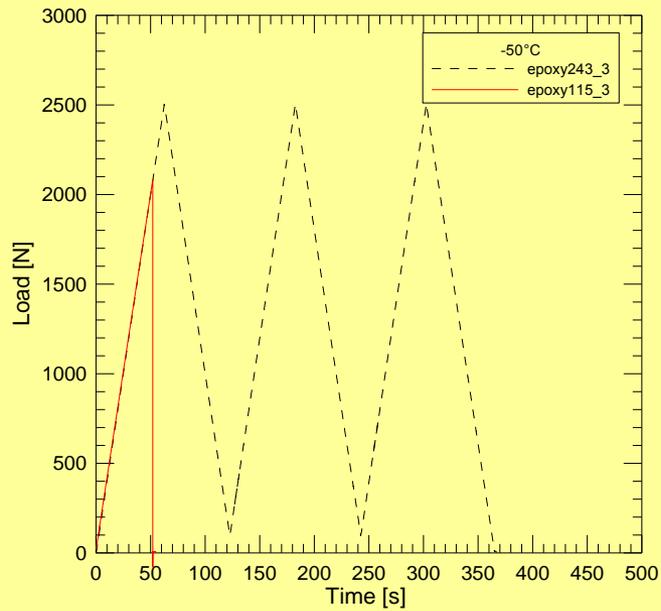
# Quasistatic methods



# Quasistatic methods

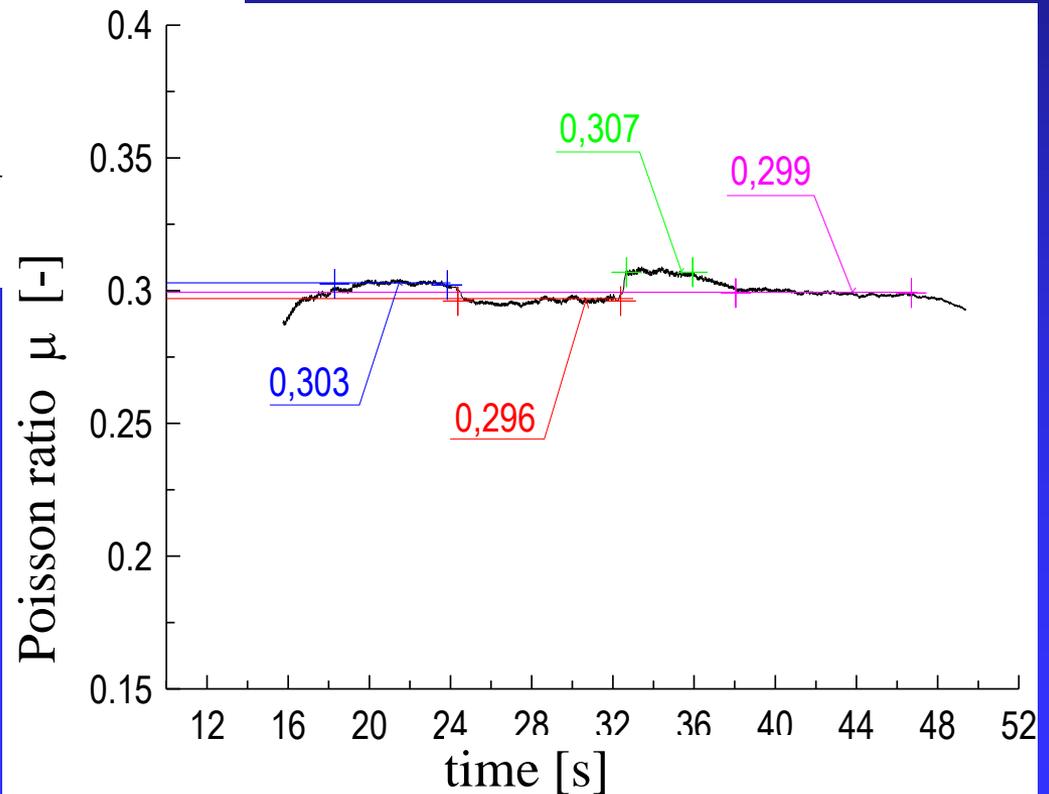
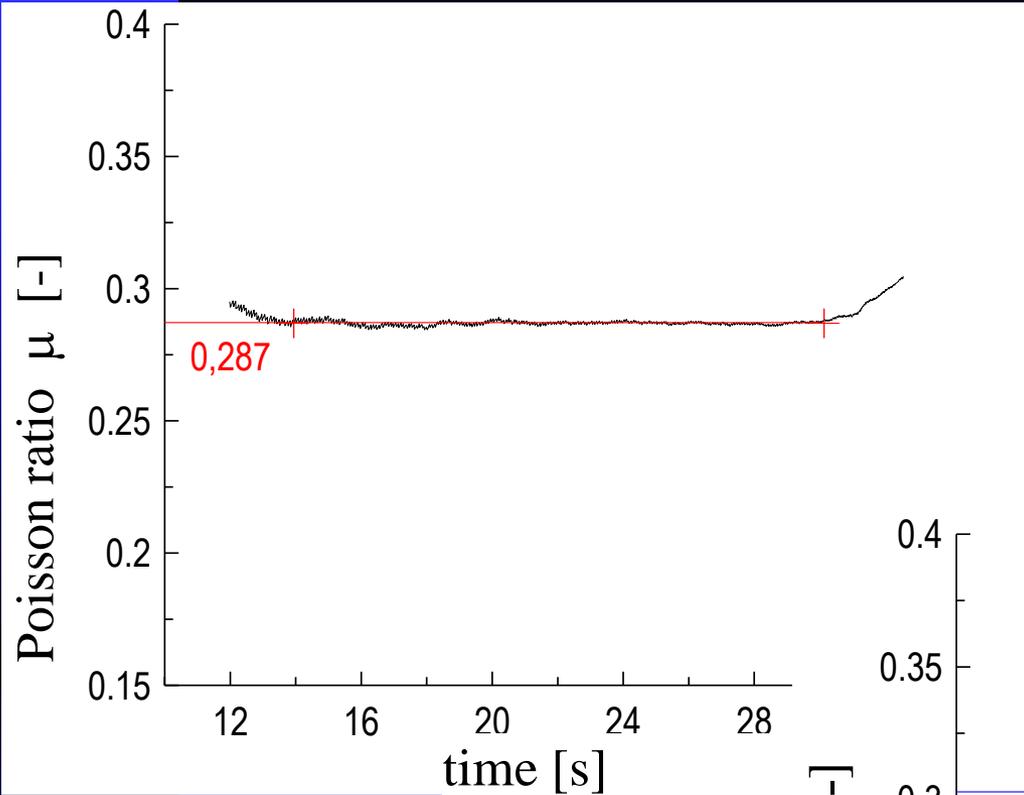


# Quasistatic methods

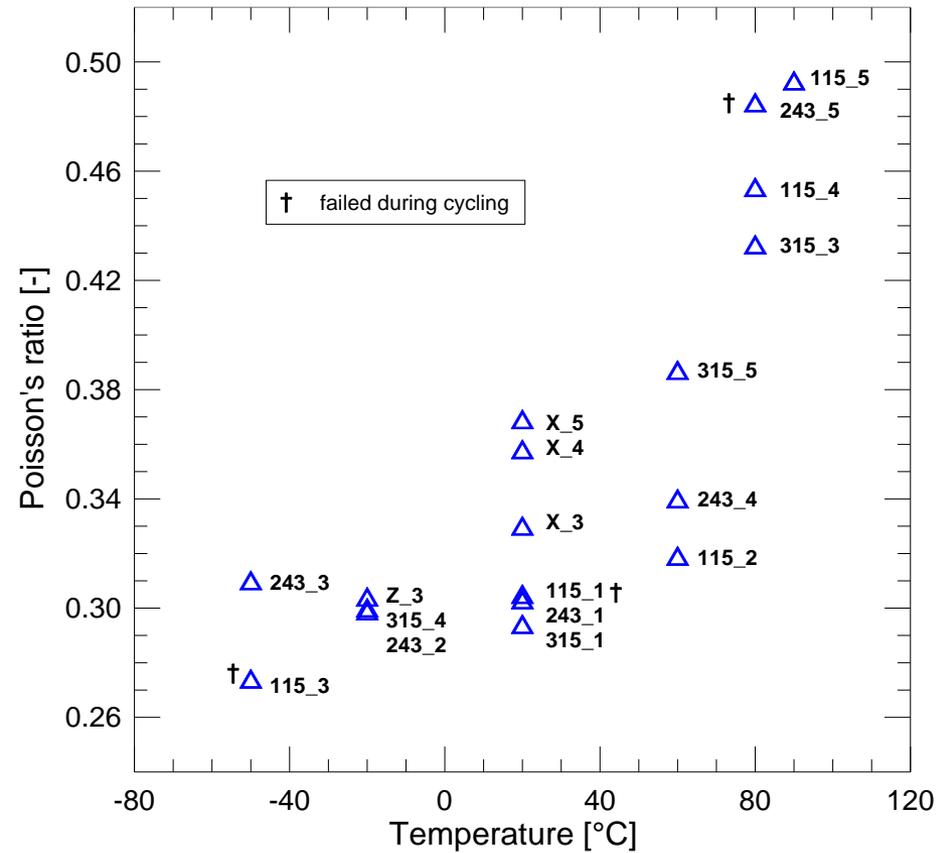
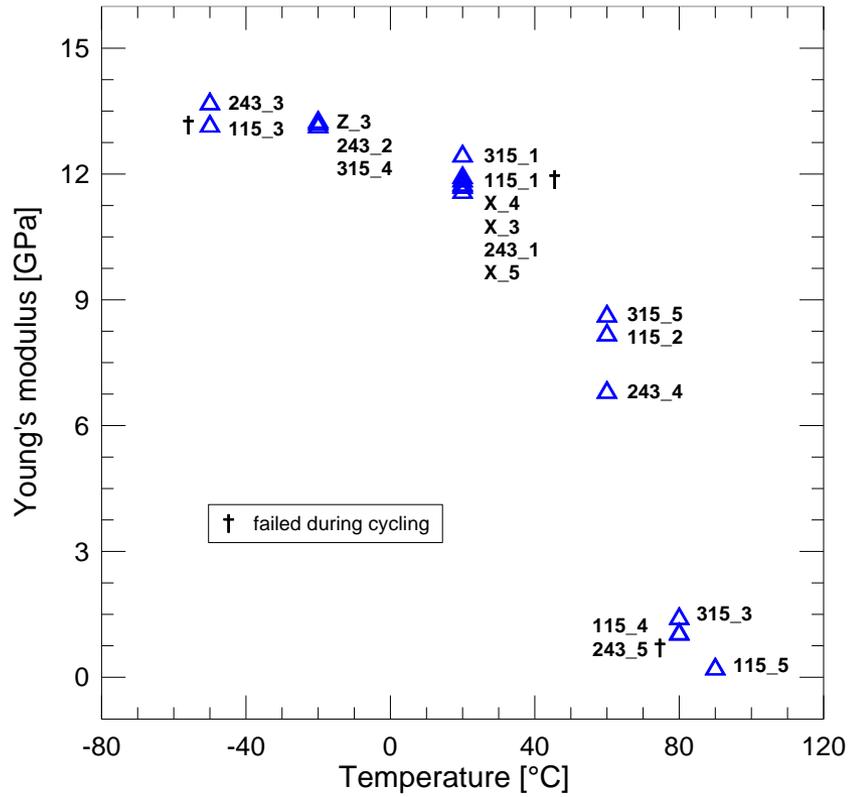


# Quasistatic methods

## Poisson's ratio

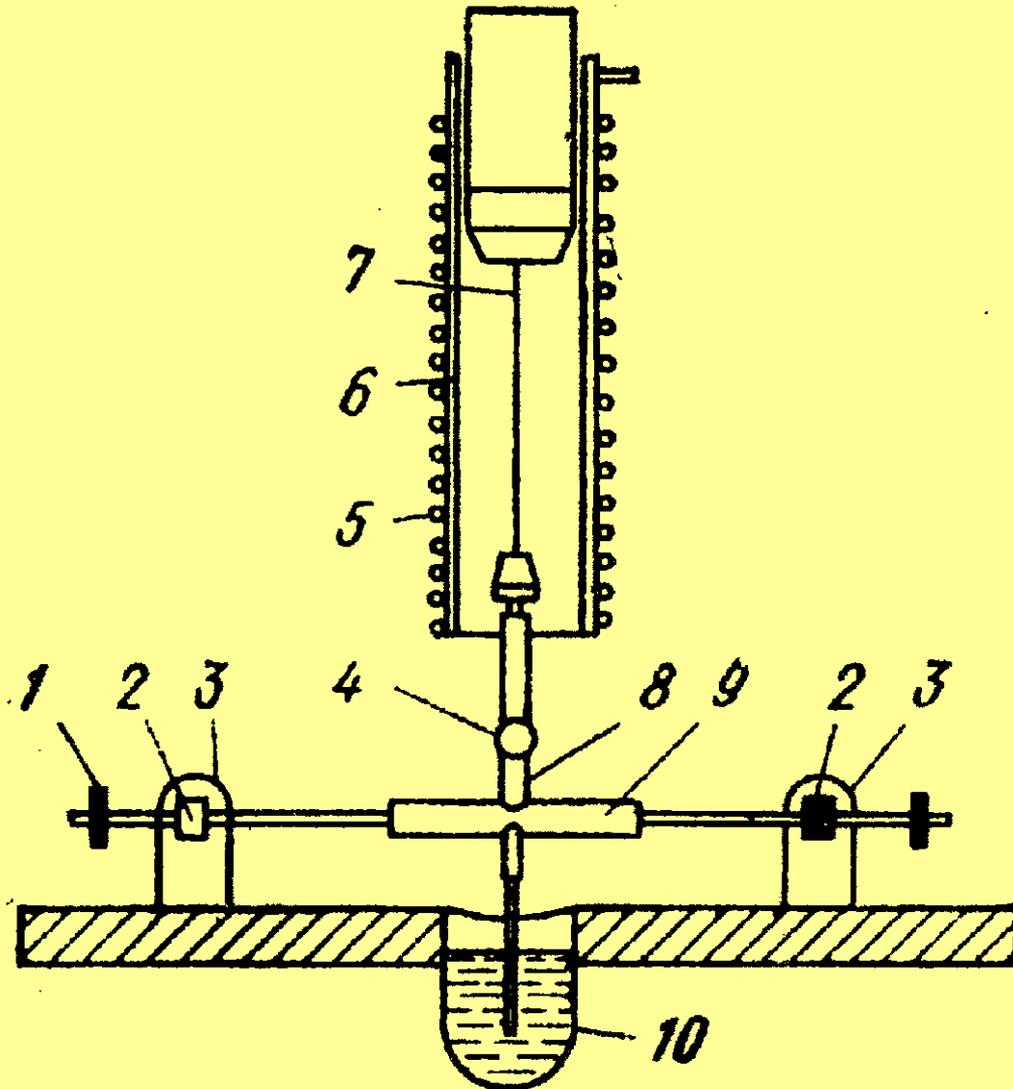


# Quasistatic methods



# torsional pendulum

## Dynamic methods



$$G = 128\pi.J_a.l.f^2/d^4$$

$d$  – sample diameter

$l$  – sample length

$J_a$  – axial moment of inertia

$f$  – eigenfrequency of pendulum vibrations

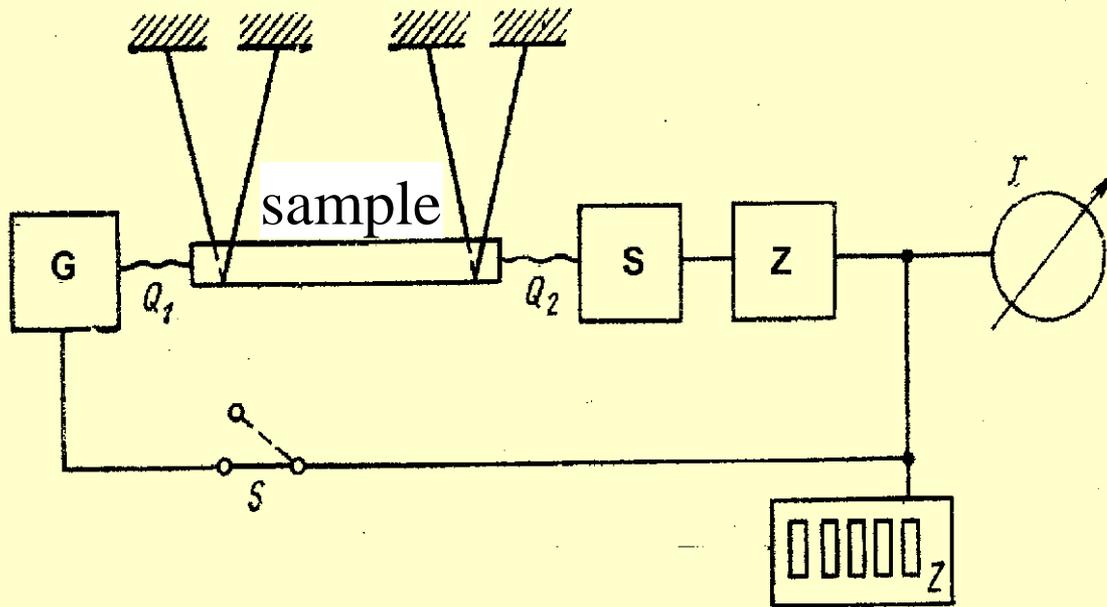
# Dynamic methods

- high frequency

$$v_D = (E/\rho)^{1/2} \quad \text{Speed of longitudinal wave propagation}$$

for sample of length  $l_0$   
eigenfrequency of vibrations  $f$ , and  
 $n$  harmonic wave

$$v_{Dn} = 2l_0 f_n / n, \quad (n = 2, 3, 4, \dots)$$



$$E = 4.l_0^2 . \rho . f_n^2 / n^2$$

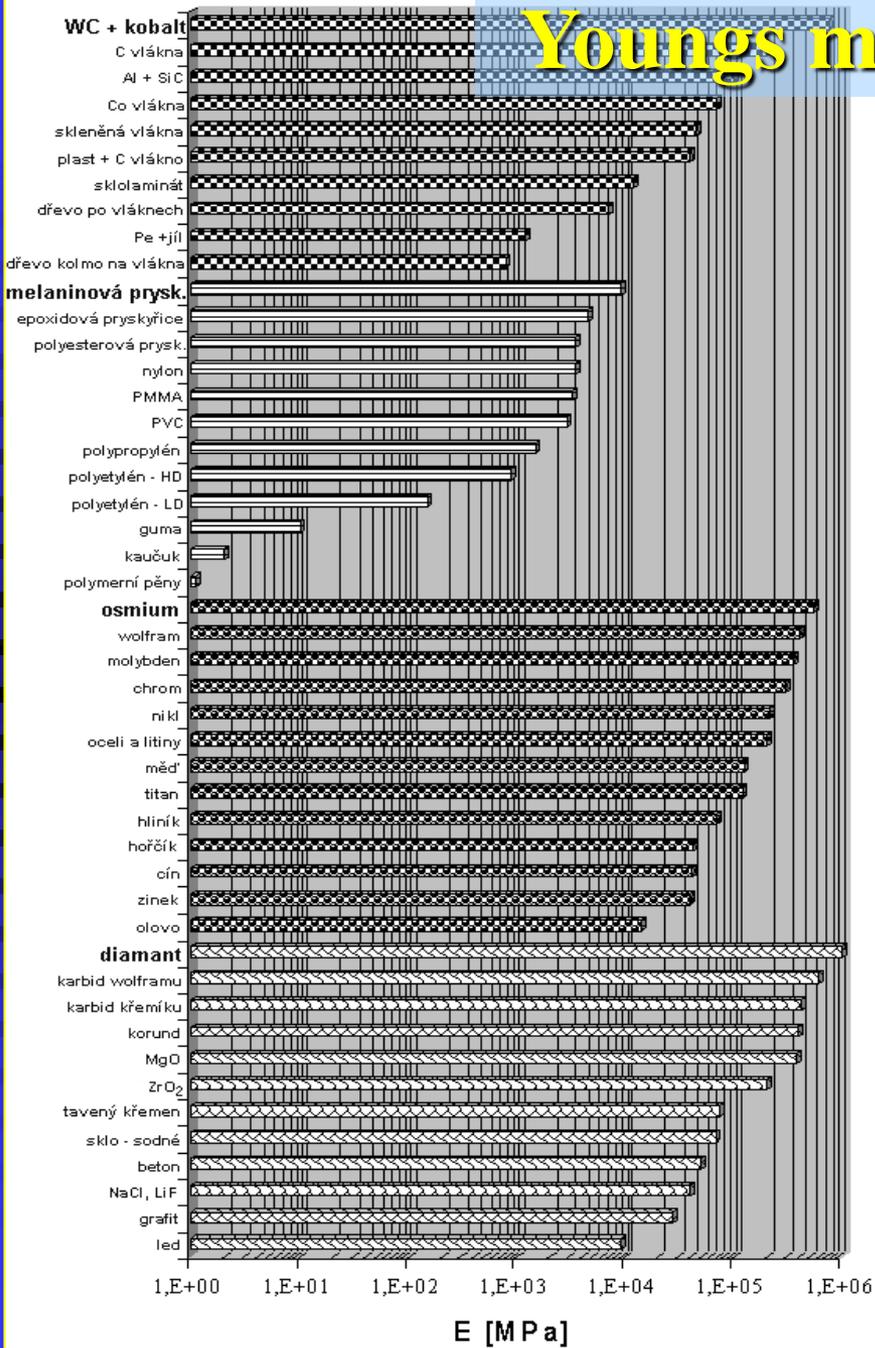
## Grindosonic

Vibrations introduced by  
small hammer blow  
(metallic, wood, plastic ..)

# Elastic properties measurements

- differences between E modulus values obtained at quasistatic and dynamic loading, (about 10% for metals, ceramics; for polymers of several orders)
- when measuring elastic properties for applications where vibrations are important, dynamic methods of measurements are preferred

# Youngs modulus - different materials



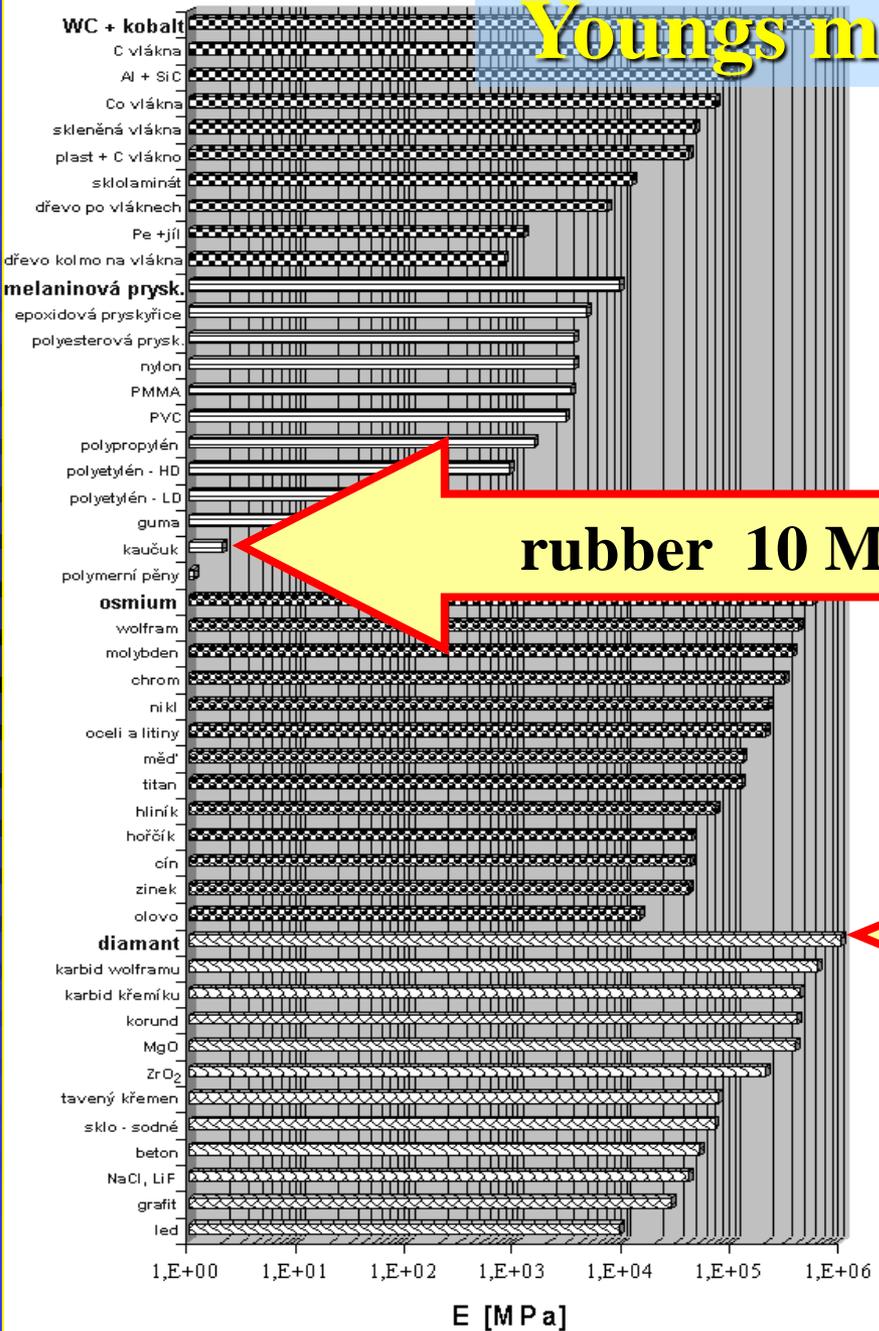
composites

polymers

metals

ceramics

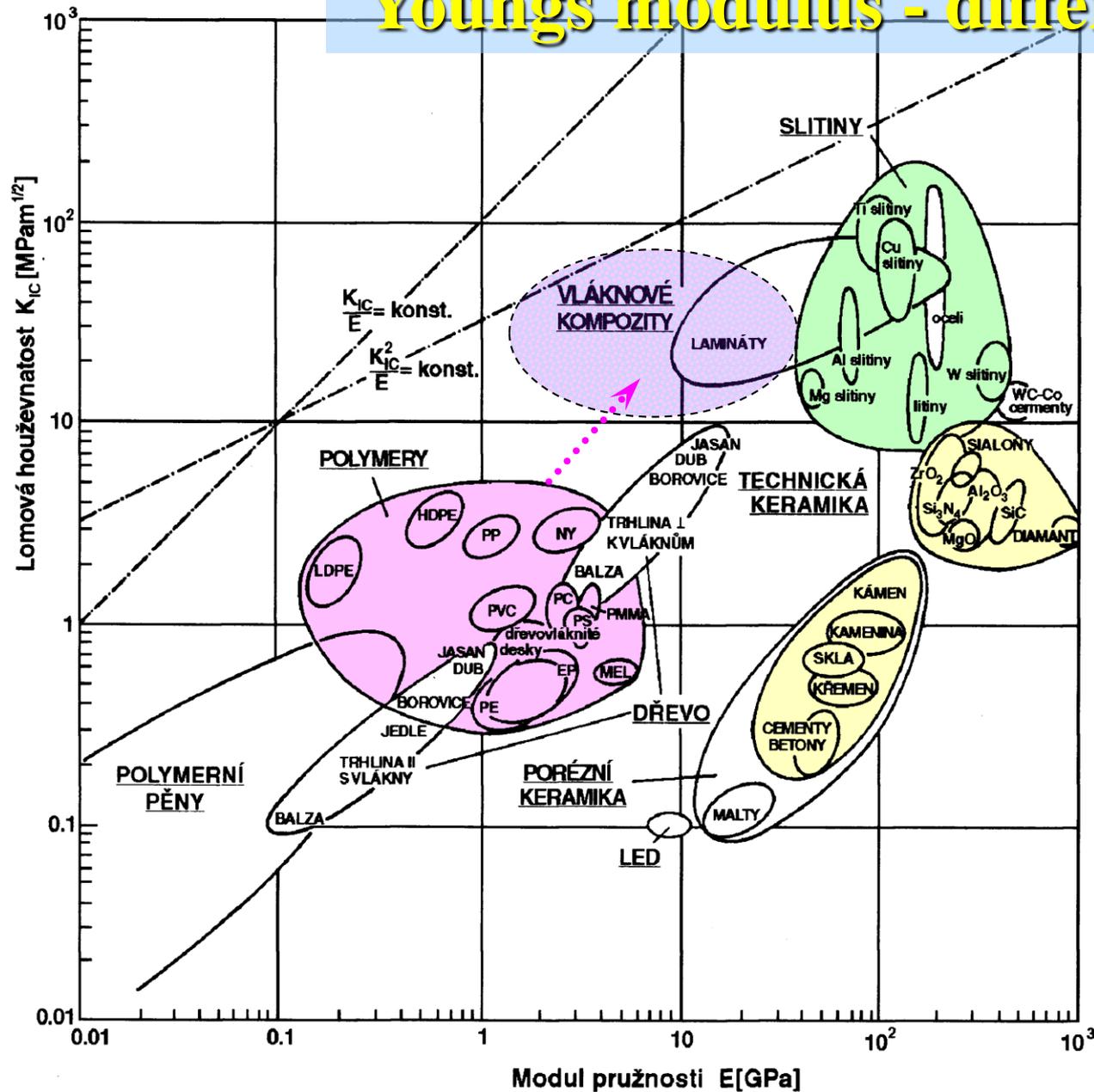
# Youngs modulus - different materials



**rubber 10 MPa**

**DIAMOND 1 000 000 MPa**

# Youngs modulus - different materials



# Young's modulus - different materials

metals, ceramics, glass  $(10^4 \text{ to } 10^6)$ MPa

polymers  $(1 \text{ to } 10^4)$ MPa

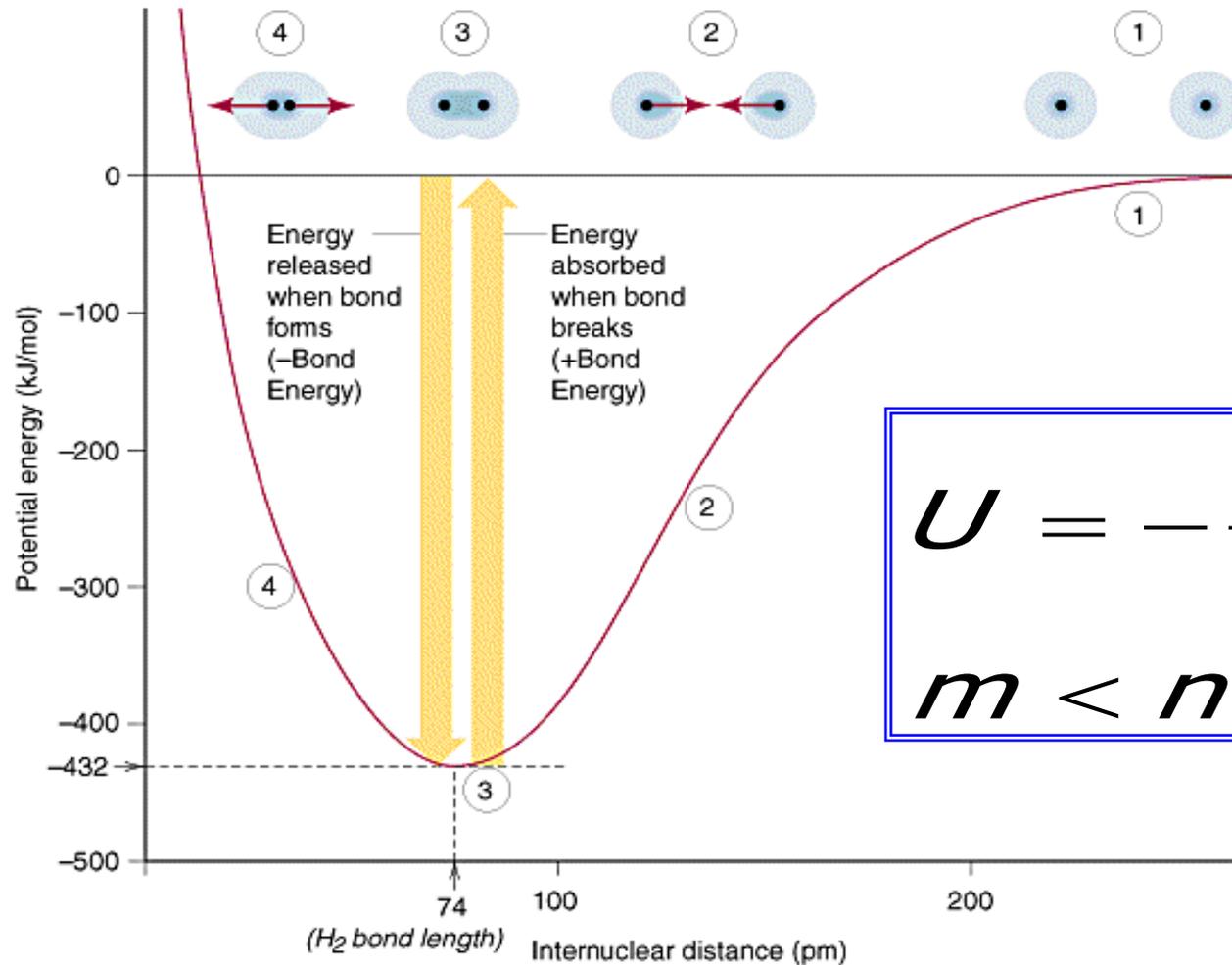
elastic deformation: dependence of Young's modulus and rigidity on

- interatomic bonding
- atoms ordering in space

- ✓ Young's modulus
- 2) Interatomic bonds
- 3) Atoms ordering in solids
- 4) Physical basis of Young's modulus

# Interatomic bonds

- Interatomic bonds – primary ( $T_m$  1000K to 5000K)
- secondary ( $T_m$  100K to 500K)



$$U = -\frac{A}{r^m} + \frac{B}{r^n}$$
$$m < n$$

# Interatomic bonds

Thanks to interatomic bonds condensed state of matter

**Metallic bond** - ions of metal atoms in cloud of electrons

>> close packed ordering = highest density

**Ion bond** - more heavily filled crystallic lattice – octaedric and tetraedric positions filled

**Covalent bond** – complicated lattice, highest density, highest strength

# Interatomic bonds

Matter performance  
controlled by  
secondary bonds

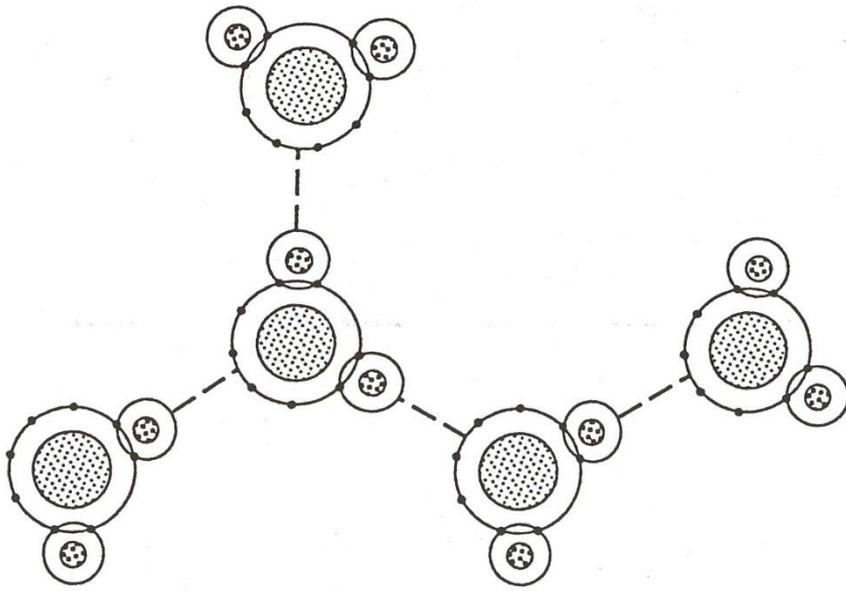
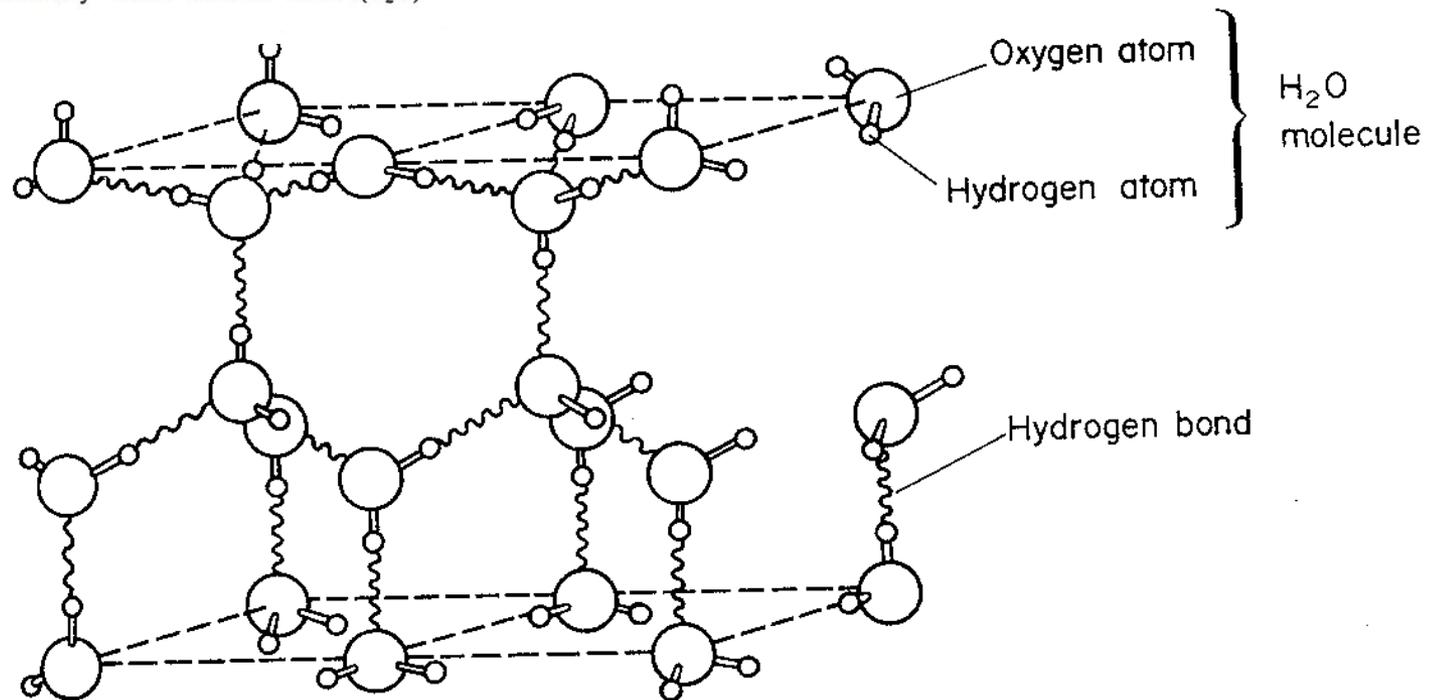
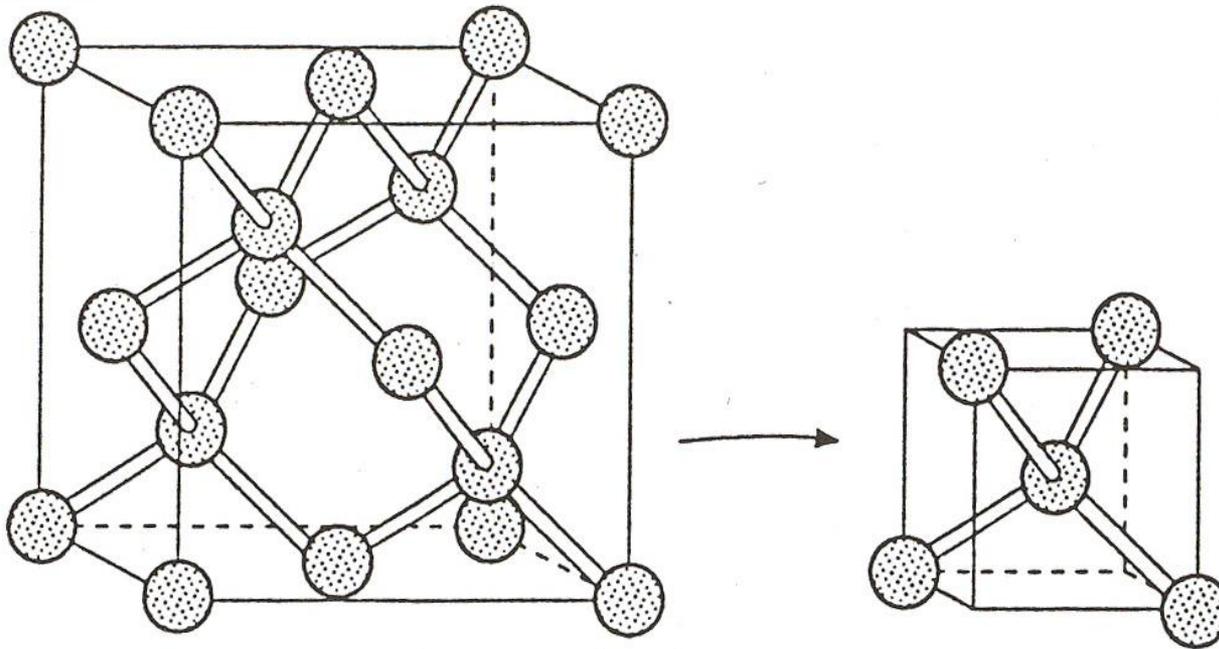


Figure 2.7 Oxygen-to-hydrogen secondary bonds between water ( $H_2O$ ) molecules.



## Matter performance controlled by primary bonds



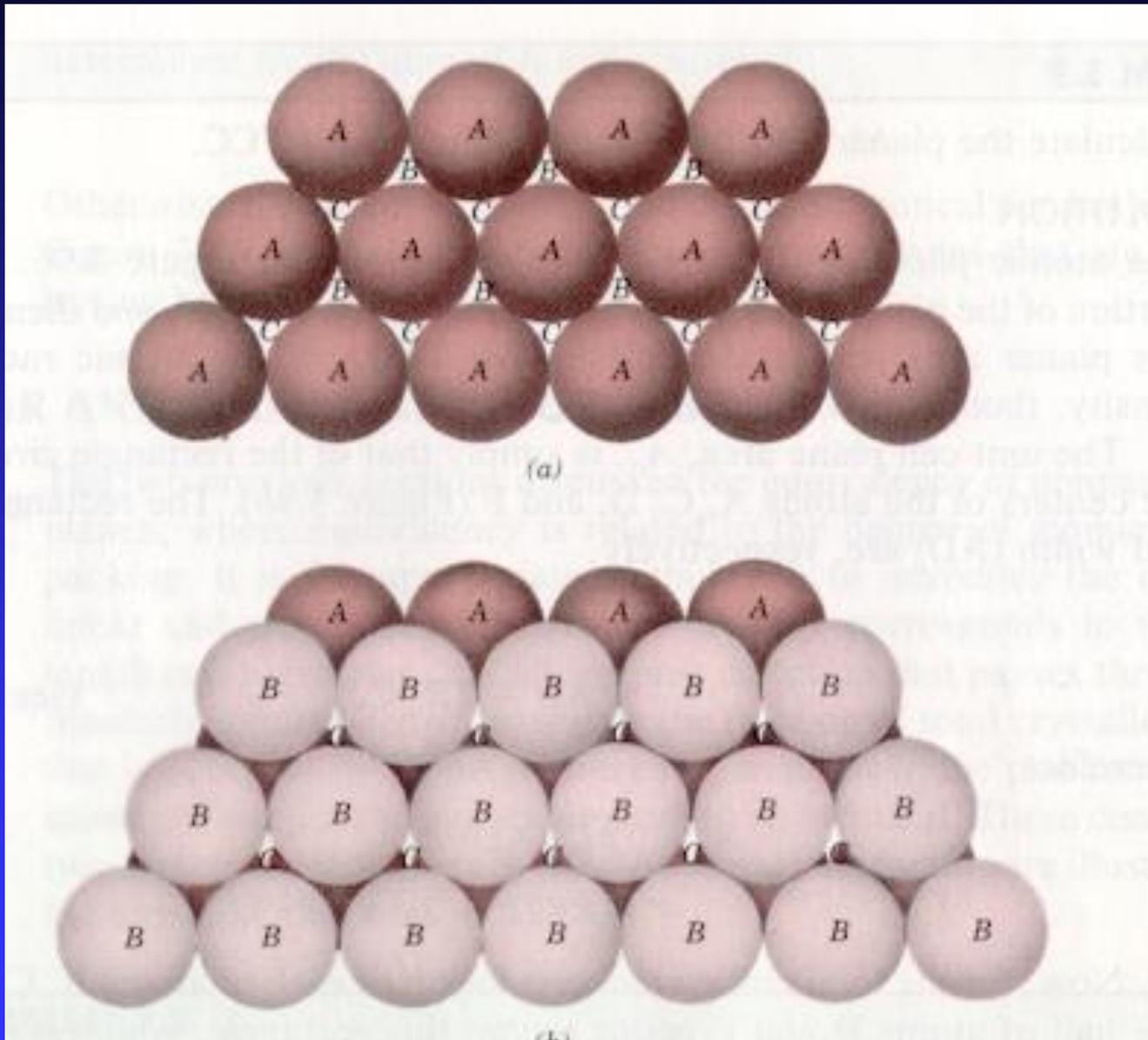
**Figure 2.5** Diamond cubic crystal structure of carbon. As a result of the strong and directional covalent bonds, diamond has the highest melting temperature, the highest hardness, and the highest elastic modulus  $E$ , of all known solids.

# condensed states

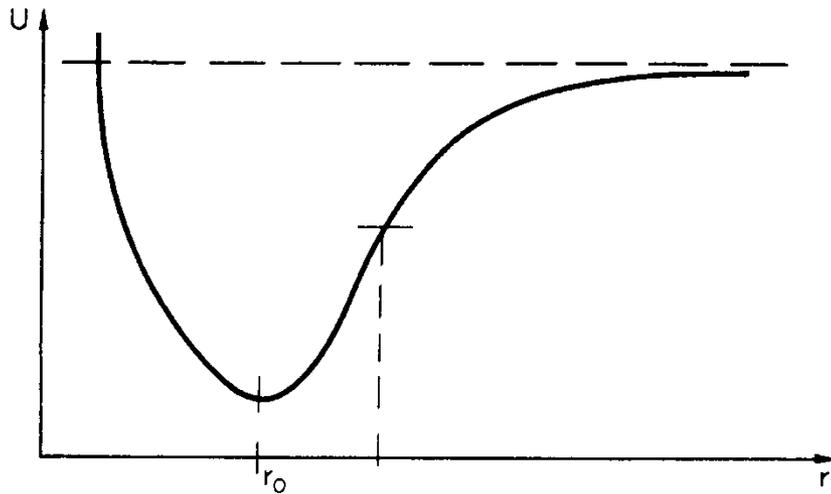
	state	bonds		K	E
		disordered	strong		
1	liquid	*		big	zero
2	Liquid crystal	*		big	small $\neq 0$
3	Rubber	Secondary	Primary	big	$E \ll K$
4	Glass		*	big	$E \approx K$
5	Crystal		*	big	$E \approx K$

- ✓ Young's modulus
- ✓ Interatomic bonds
- 2) Atoms ordering in solids
- 3) Physical basis of Young's modulus

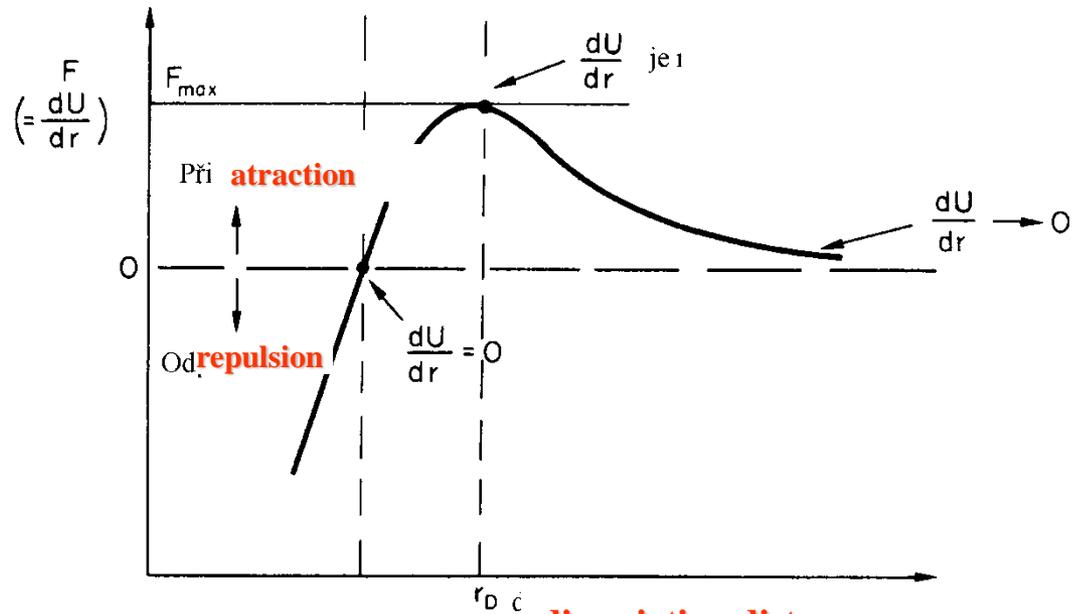
# Close packed lattice - metals



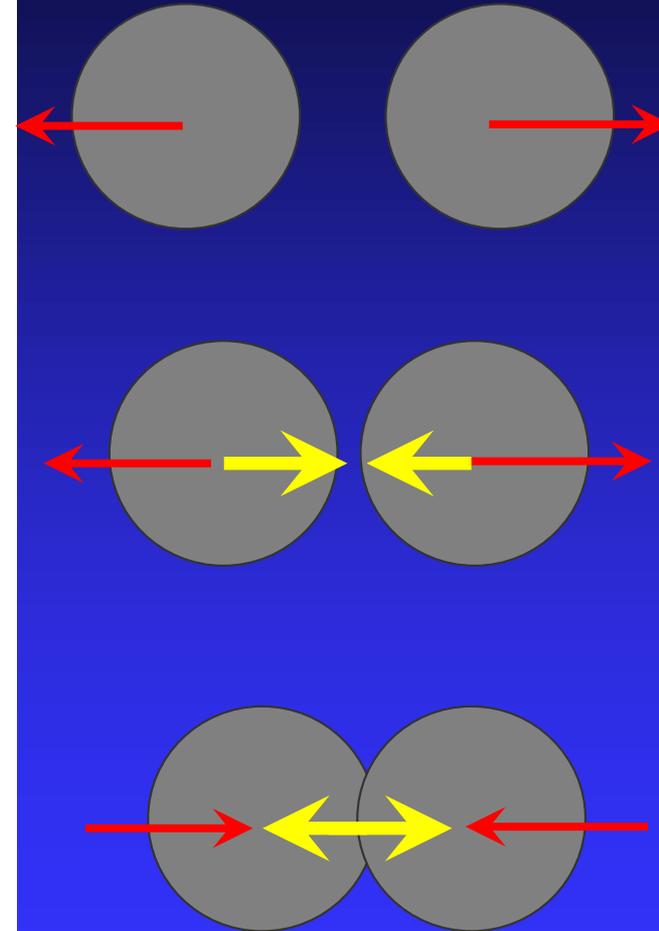
# Atoms ordering



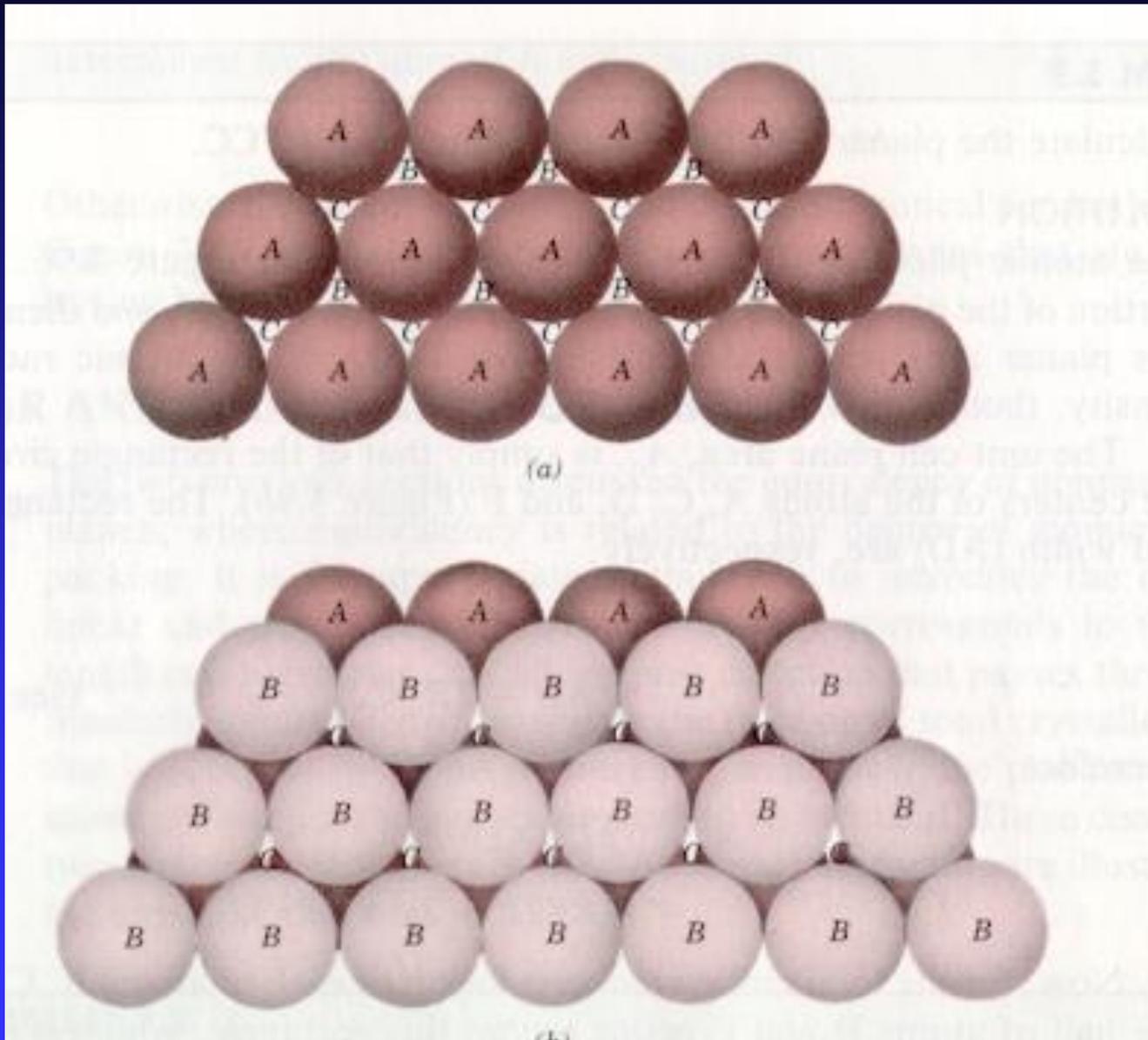
$r_0$  mřížkový parametr



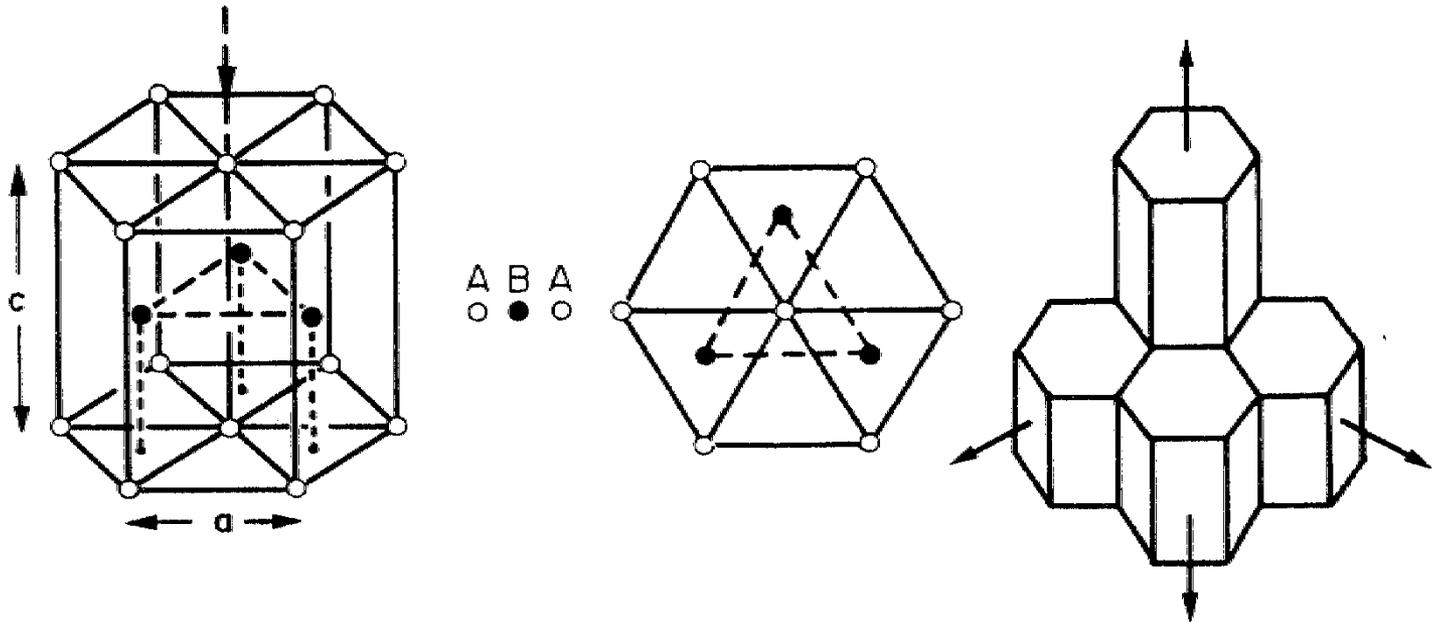
$r_{D c}$  dissociation distance



# Close packed lattice - metals

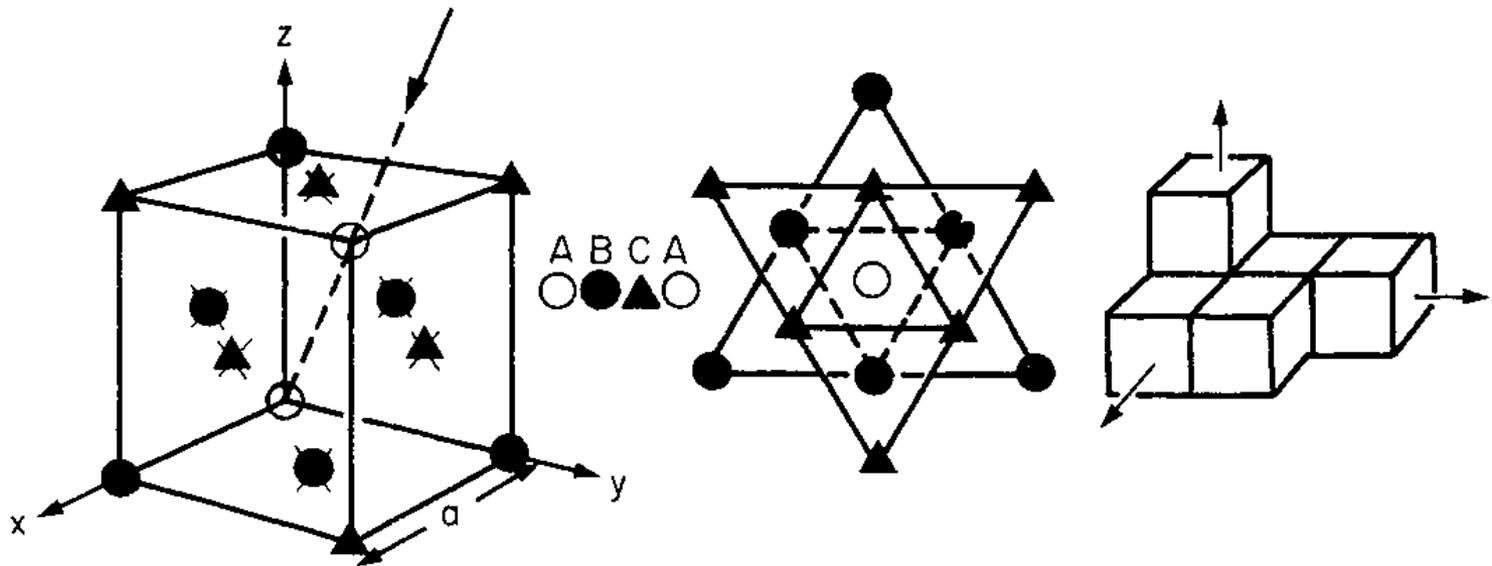


# Structure hcp –Mg, Zn, Co, a - Ti hexagonal - close – packed structure

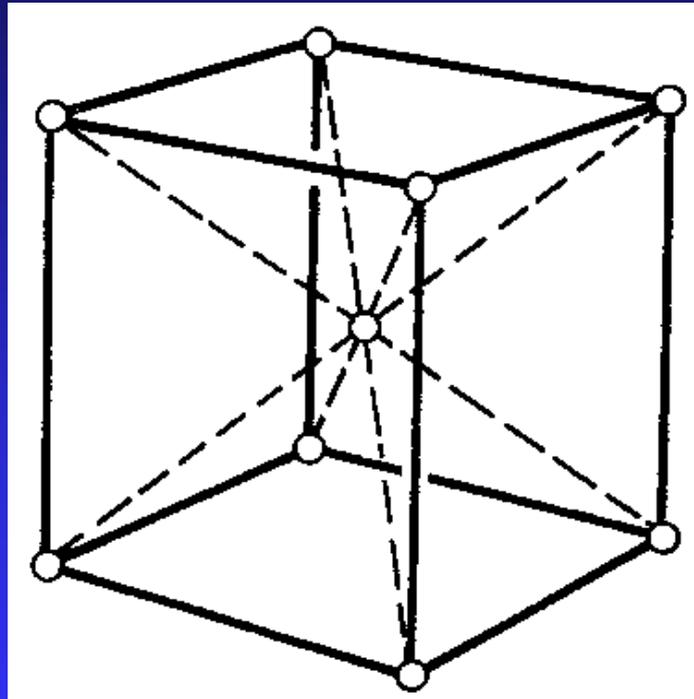


# Structure fcc – Al, Cu, Au, Ag, Pt, Ni, $\gamma$ -Fe

## face – centred – cubic structure



# Structure bcc – $\alpha$ -Fe, Mo, W, Ta body – centred - cubic structure

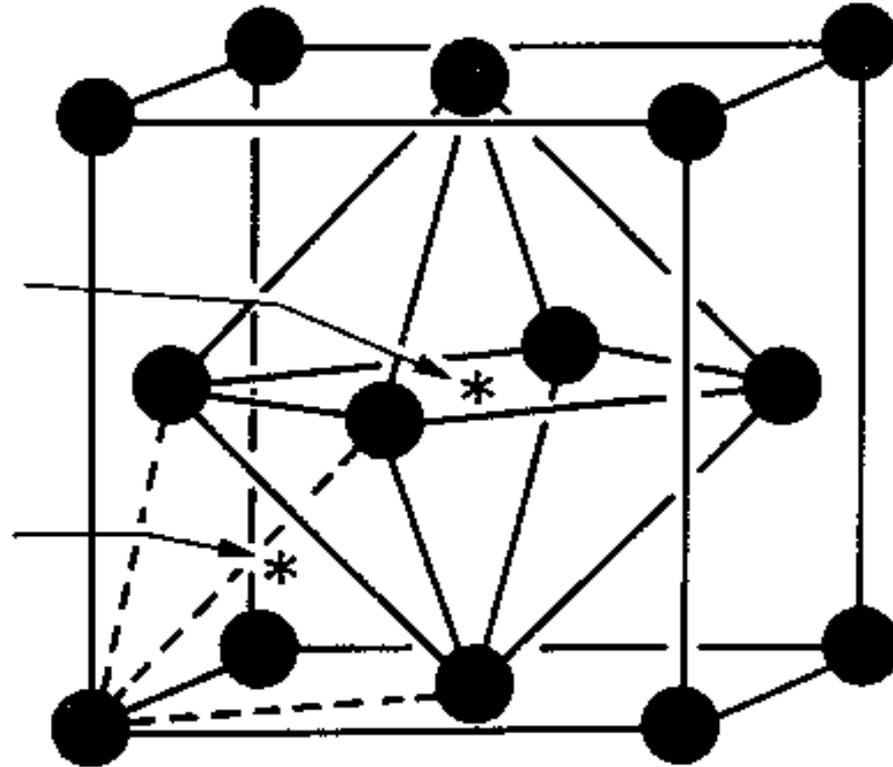


- Polymorphy – transformation of lattice with temperature,  $\alpha - \gamma - \delta$  (910/1400) $^{\circ}\text{C}$

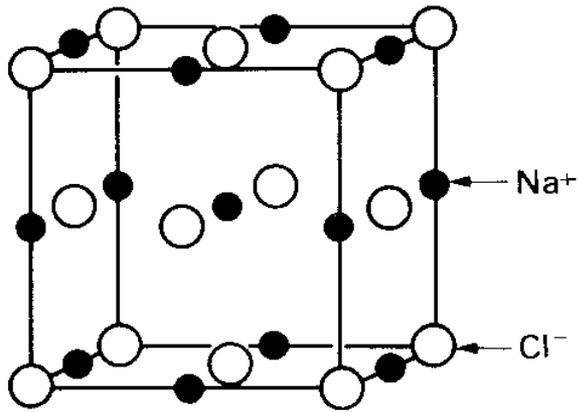
# Ion bonding - ceramics

octaedic  
location

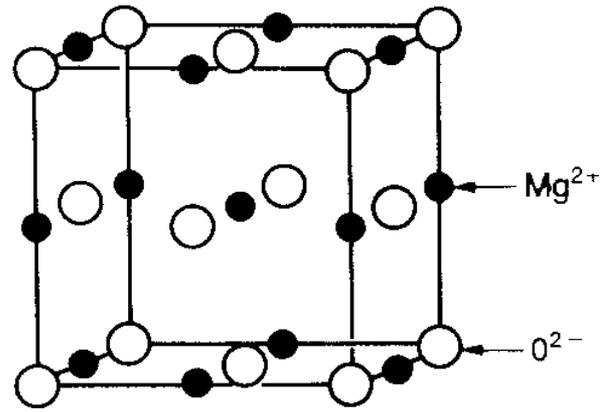
tetraedic  
location



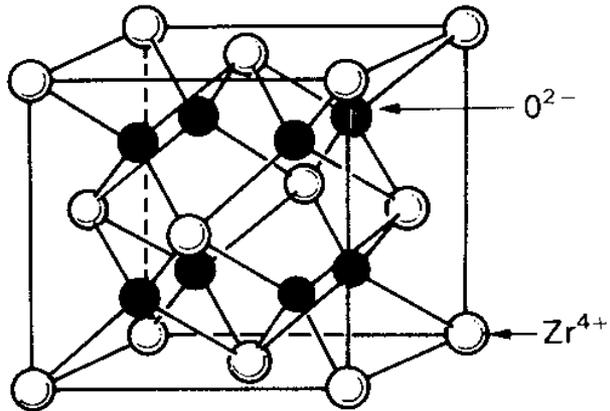
# Ion bonding - ceramics



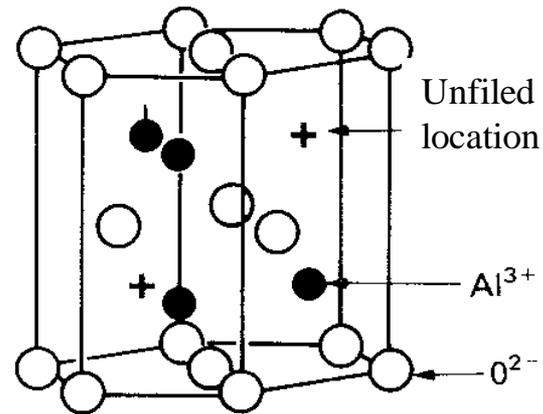
(a)



(b)



(c)



(d)

# Covalent bonding - ceramics

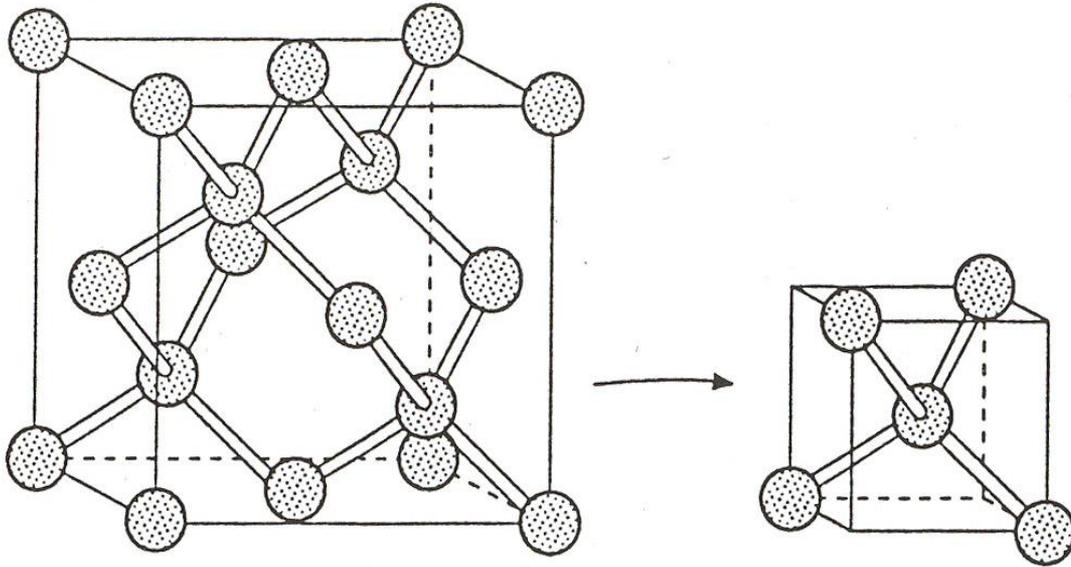
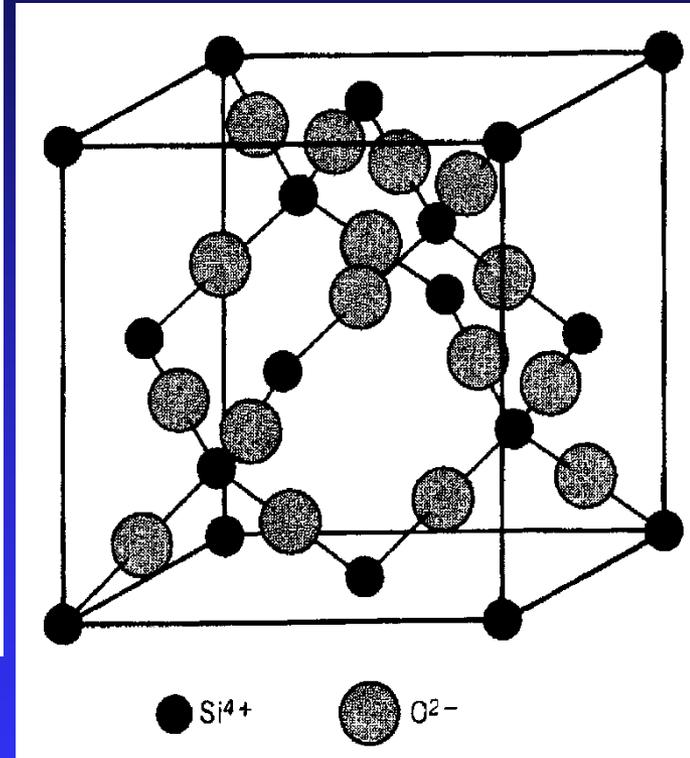


Figure 2.5 Diamond cubic crystal structure of carbon. As a result of the strong and directional covalent bonds, diamond has the highest melting temperature, the highest hardness, and the highest elastic modulus  $E$ , of all known solids.

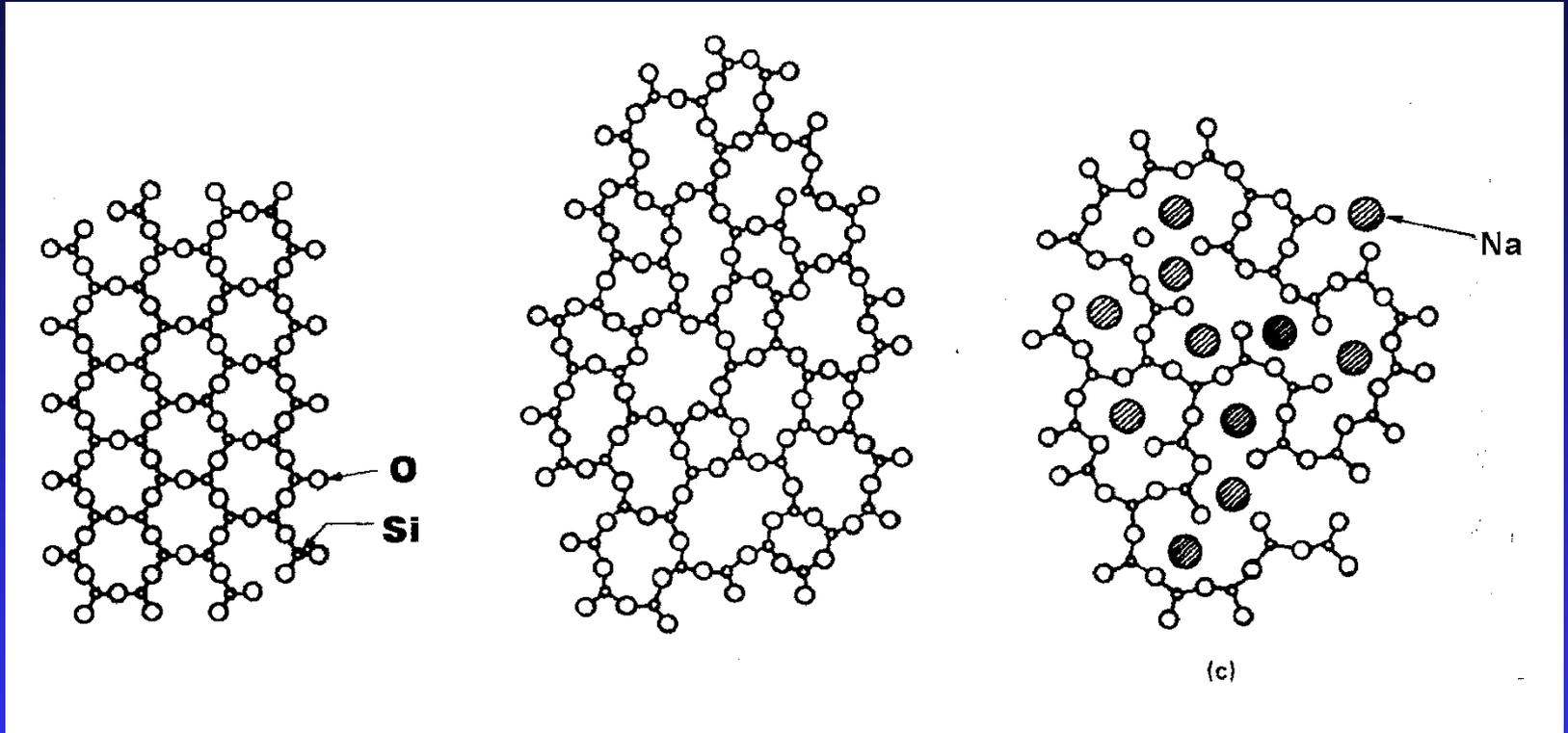
## Diamond



## Silica

Lattice is not exactly the close packed, but atoms fully filled space

# Covalent bonding - glass



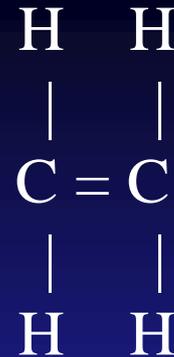
Silica glass – melting point  $1200^{\circ}\text{C}$

Na, Ca, Fe – terminators –  $700^{\circ}\text{C}$

# Covalent bonding – polymers (plastics)

- Thermoplastic PE, PP, PS, PMMA
- Thermosets (reactoplasts)
- Elastomers (rubbers, elastic)

**Etylen**



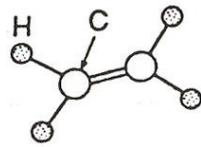
**monomer**

<b>Number of monomers (n)</b>	<b>Temperature of softening</b>	<b>State at 20°C</b>
<b>1</b>	<b>-167</b>	<b>gas</b>
<b>6</b>	<b>-12</b>	<b>liquid</b>
<b>35</b>	<b>37</b>	<b>oil</b>
<b>140</b>	<b>93</b>	<b>wax</b>
<b>430</b>	<b>109</b>	<b>solids</b>

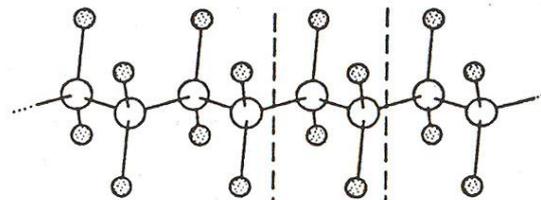
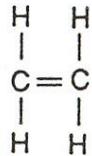
▼  
Polymerization degree

$$10^3 \approx 10^5$$

# Polymers (plastics)



Ethylene



Polyethylene

repeating unit

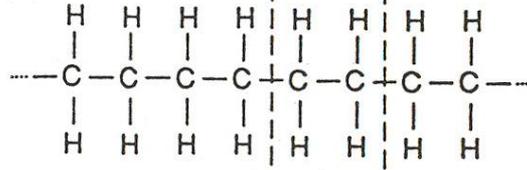
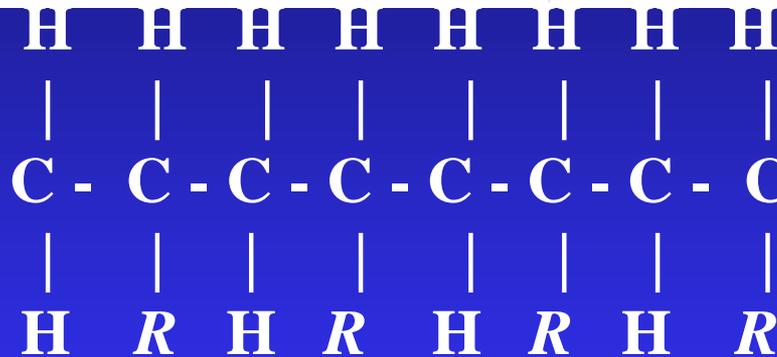
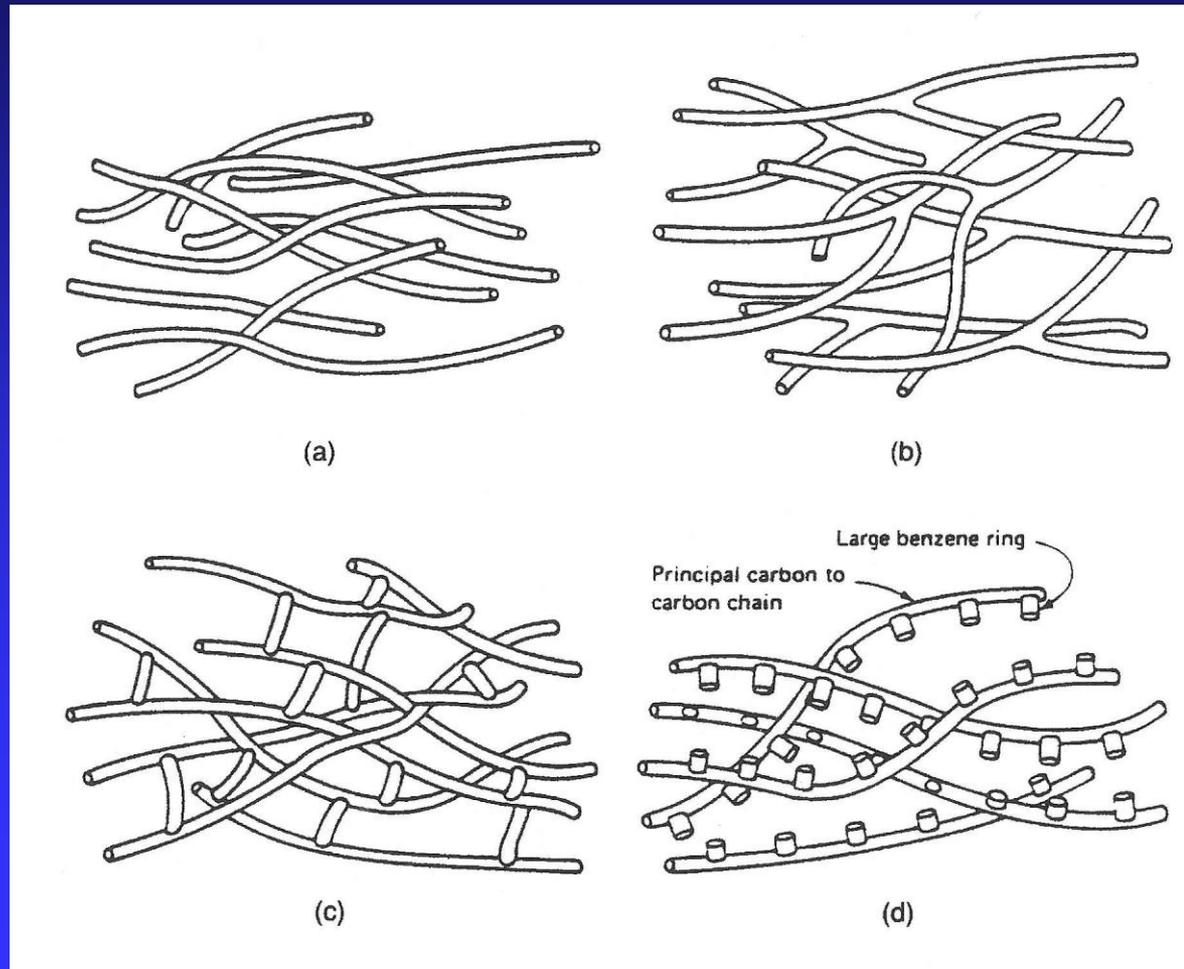


Figure 2.6 Molecular structures of ethylene gas ( $\text{C}_2\text{H}_4$ ) and polyethylene polymer. The double bond in ethylene is replaced by two single bonds in polyethylene, permitting formation of the chain molecule.



- R ---- H** PE / polyethylene (tubes, isolators)
- R ---- CH<sub>3</sub>** polypropylen (better resistance against light)
- R ---- Cl** polyvinylchlorid (roofing and floor covering)
- R ---- C<sub>6</sub>H<sub>5</sub>** polystyrén (gramophone disks, cutlery, příbory, izolation)

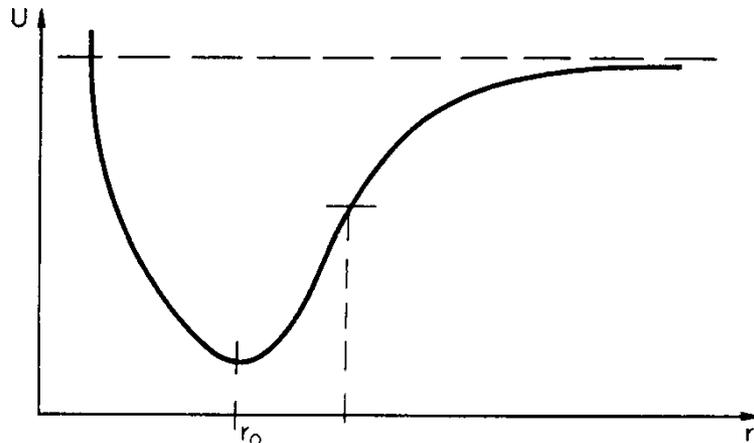
## Shape of plastics molecule



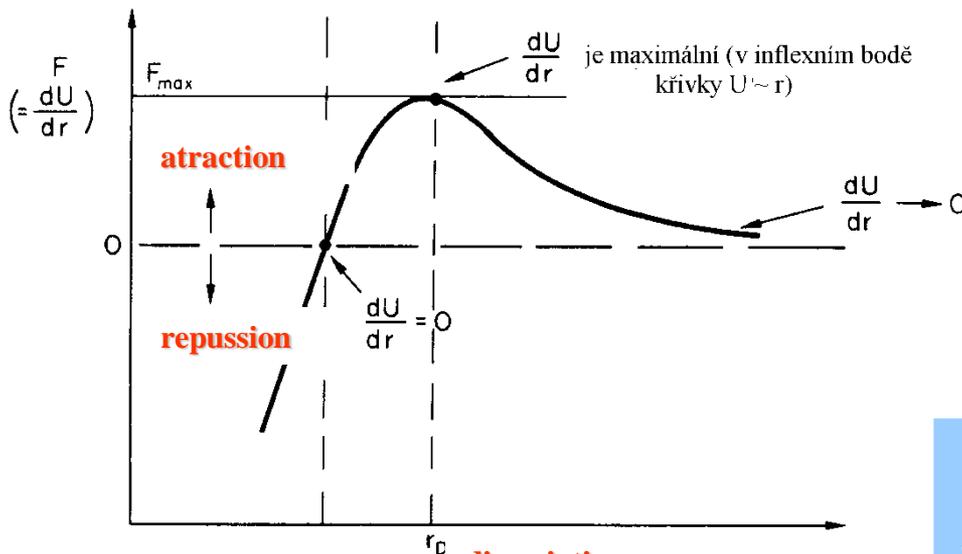
- ✓ Young's modulus
- ✓ Interatomic bonds
- ✓ Atoms ordering in solids

Physical basis of Young's modulus

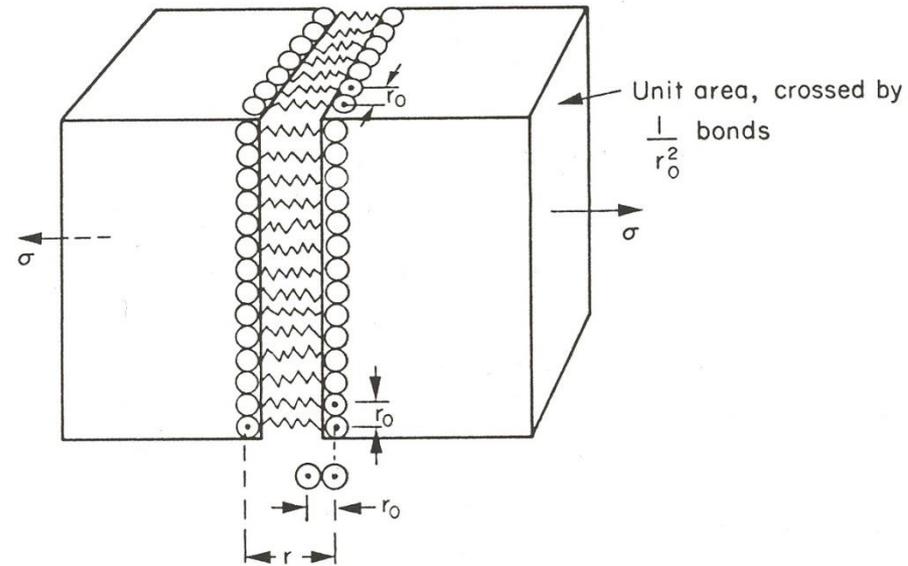
# Physical nature of elastic modulus



$r_0$  mřížkový parametr



**dissociation distance**



**rigidity of the bond**

(second order derivative of the potential energy)

$$S_0 = \left( \frac{d^2U}{dr^2} \right)_{r=r_0}$$

## bond rigidity

## Young's modulus

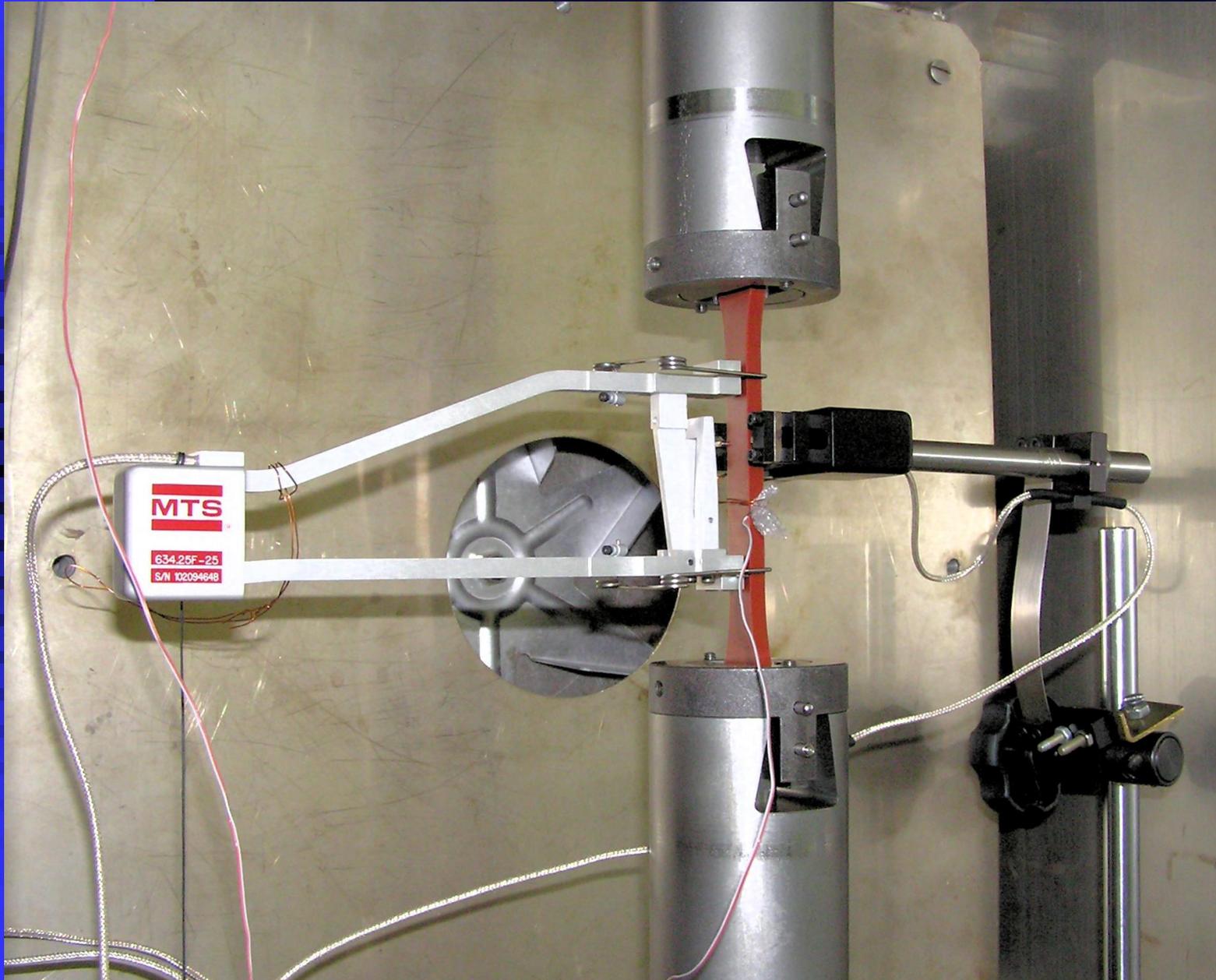
bond type	$S_0$	$E \approx S_0/r_0$ [MPa]
C-C	180	1 000 000
ionic Na-Cl	9 - 21	(0.3 – 0.7) 100 000
metallic Cu-Cu	15 - 40	(0.3 – 1.5) 100 000
H - bond	2	8 000
Van der Waals	1	2 000

Valid for metals and ceramics; for selected plastics (rubber, caoouchouck) by 3 order lower comparing to theory.

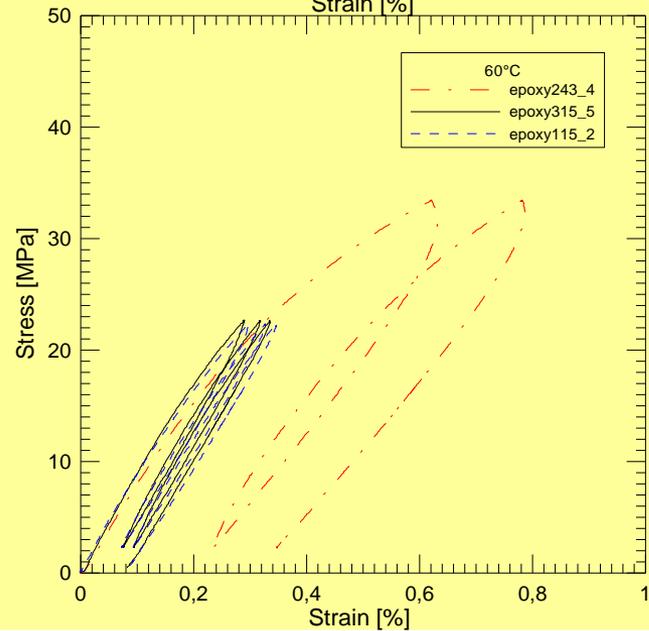
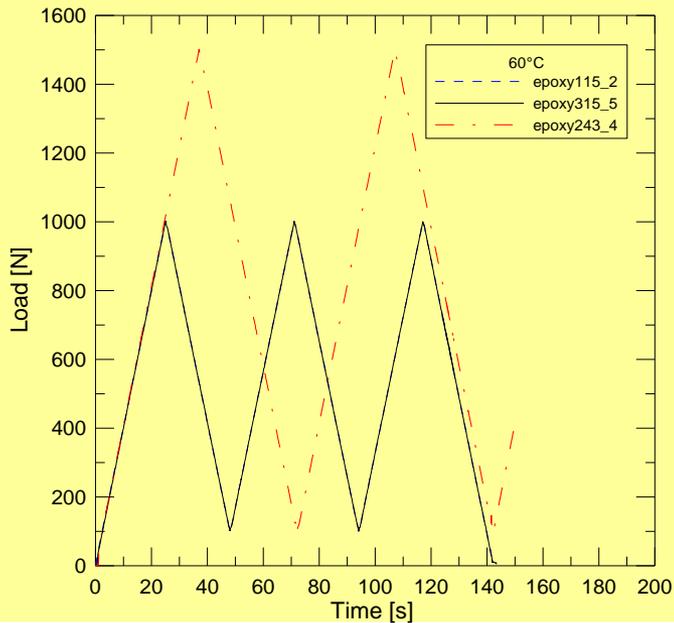
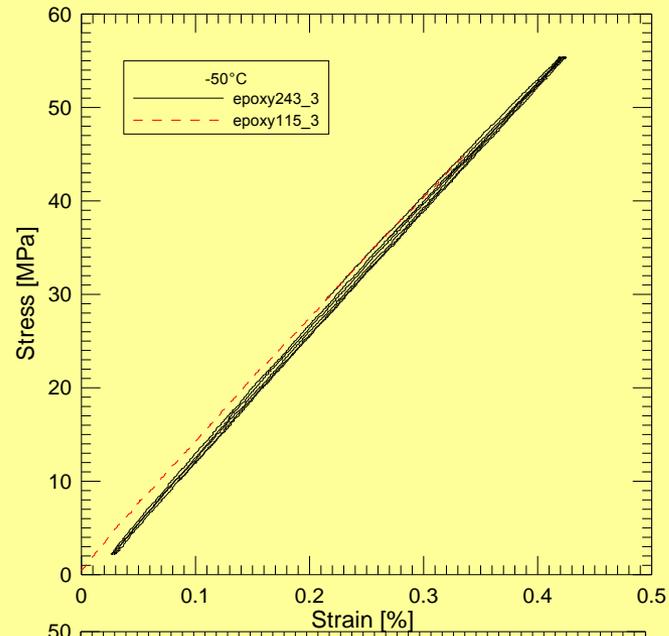
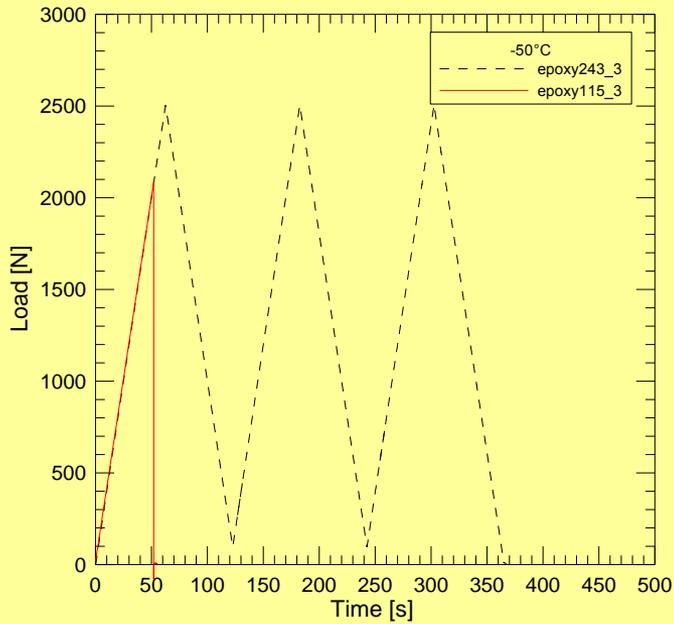
# Inelasticity – internal dumping

- Deformation energy losses during vibrations
- Noise dumping

# Určování modulu pružnosti

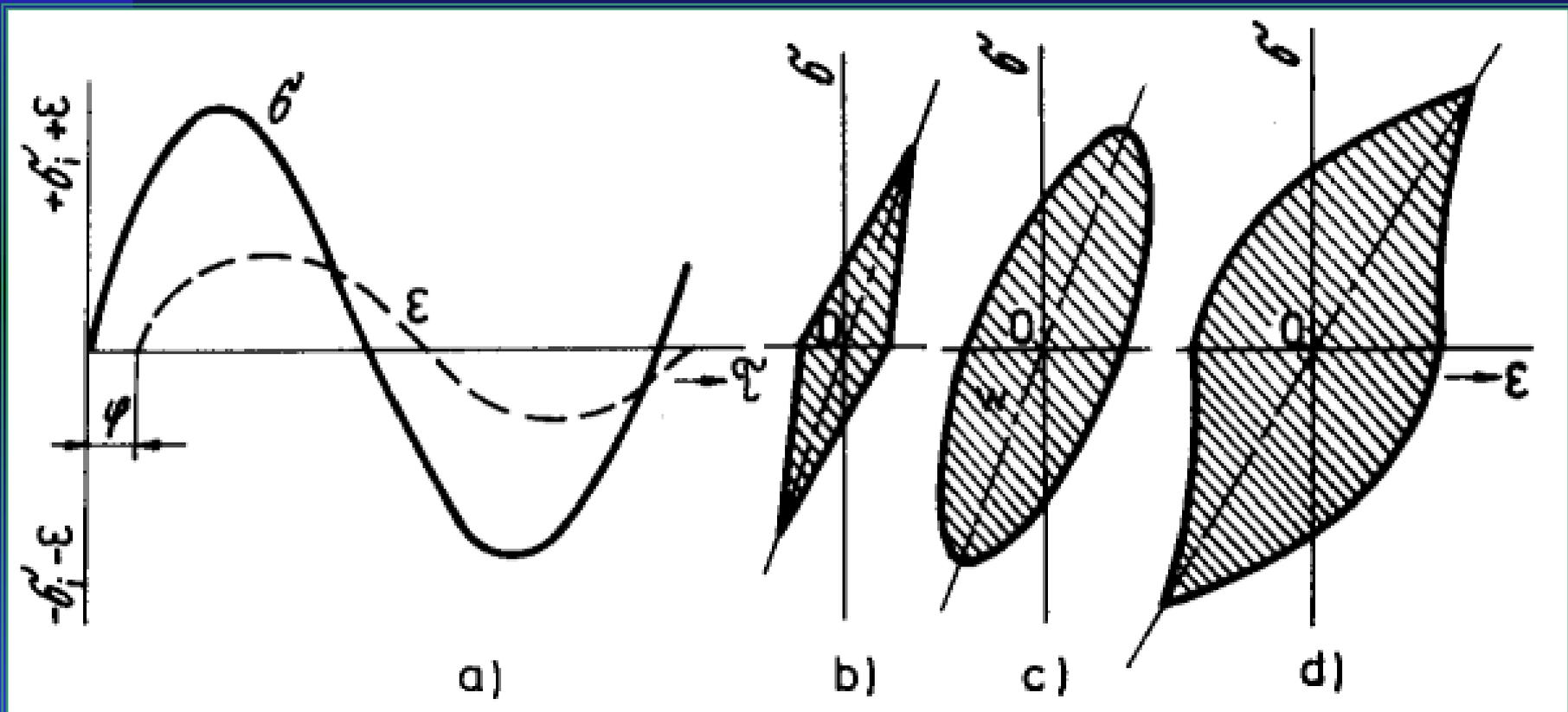


# Quasistatic methods



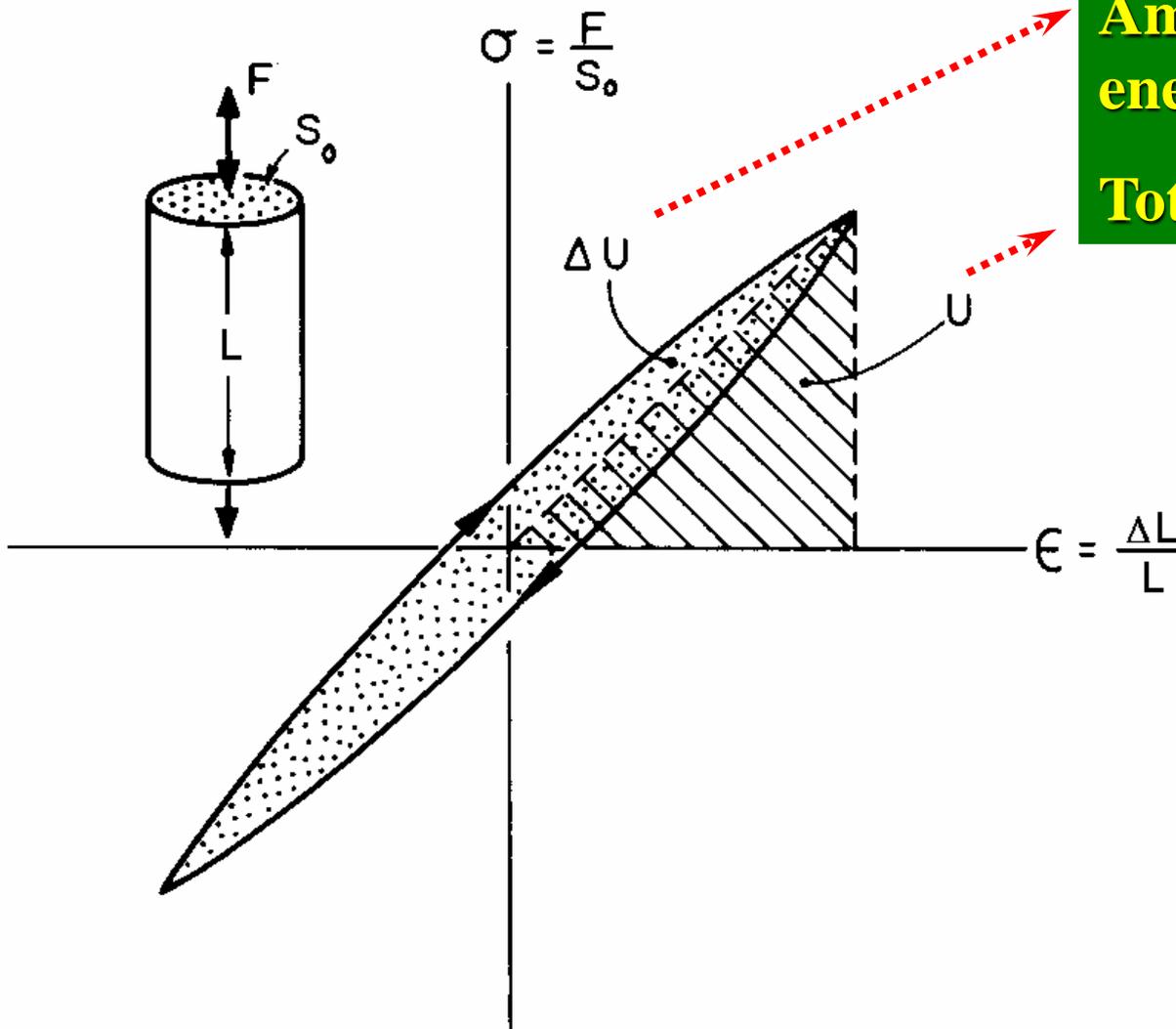
# Hysteresis - inelasticity

$Q^{-1} = \operatorname{tg} \alpha$  (internal damping – coefficient of internal losses)



**Ability of the materials to dissipate elastic energy during vibrations**

# Inelasticity / Internal damping



Amount of dissipated energy per one cycle

Total energy input

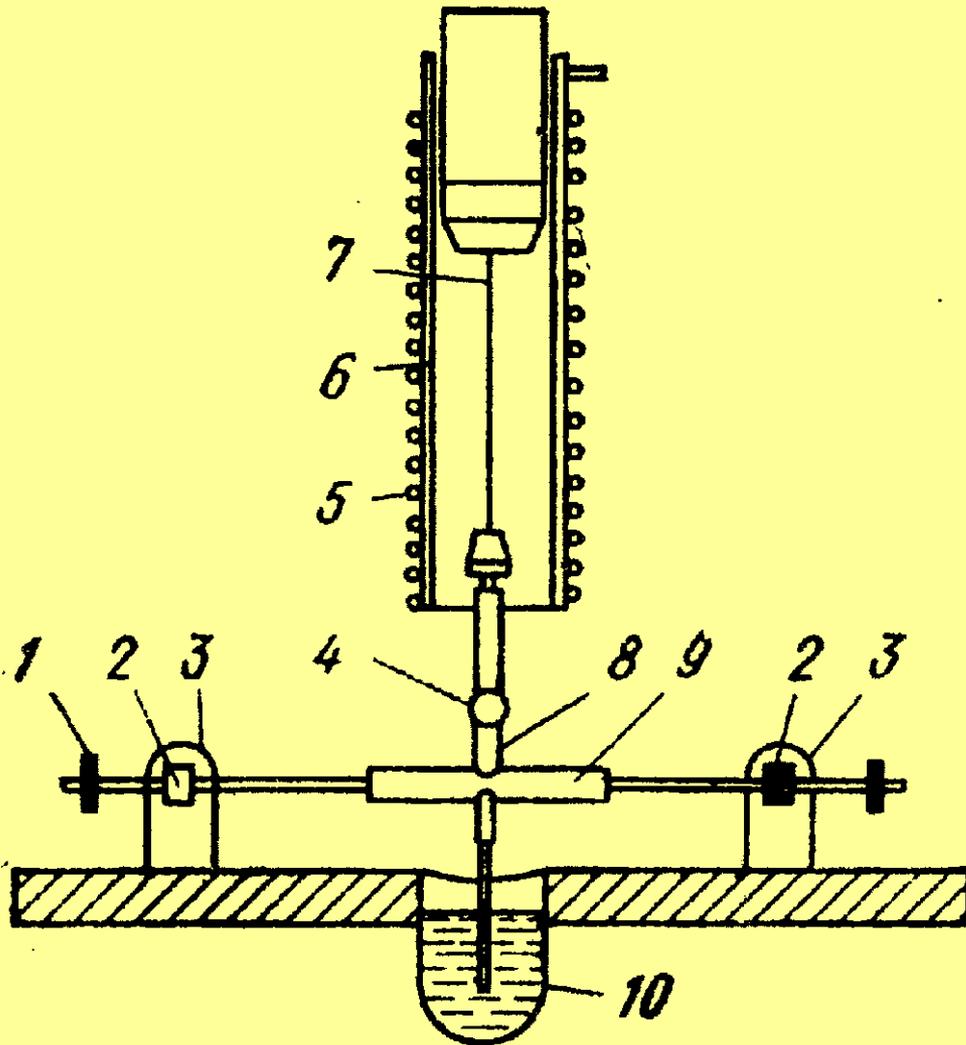
$$\eta = \frac{\Delta U}{2\pi U}$$

# Measurements of internal damping parameters

## Torzional pendulum

$$G = 128\pi.Ja.l.f^2 / d^4$$

$\Delta\omega$  – half of width of the resonance peak in the half of its height  
 $f$  – eigenfrequency pendulum  
 $\omega_0$  – resonance circular frequency of sample



$$\eta = \frac{\Delta U}{2\pi U} = \frac{\Delta\omega}{\sqrt{3}\omega_0}$$

# Inelasticity – internal dumping

- Ability of the material dissipate elastic energy during vibrations

- Internal dumping quantification

- coefficient of internal losses

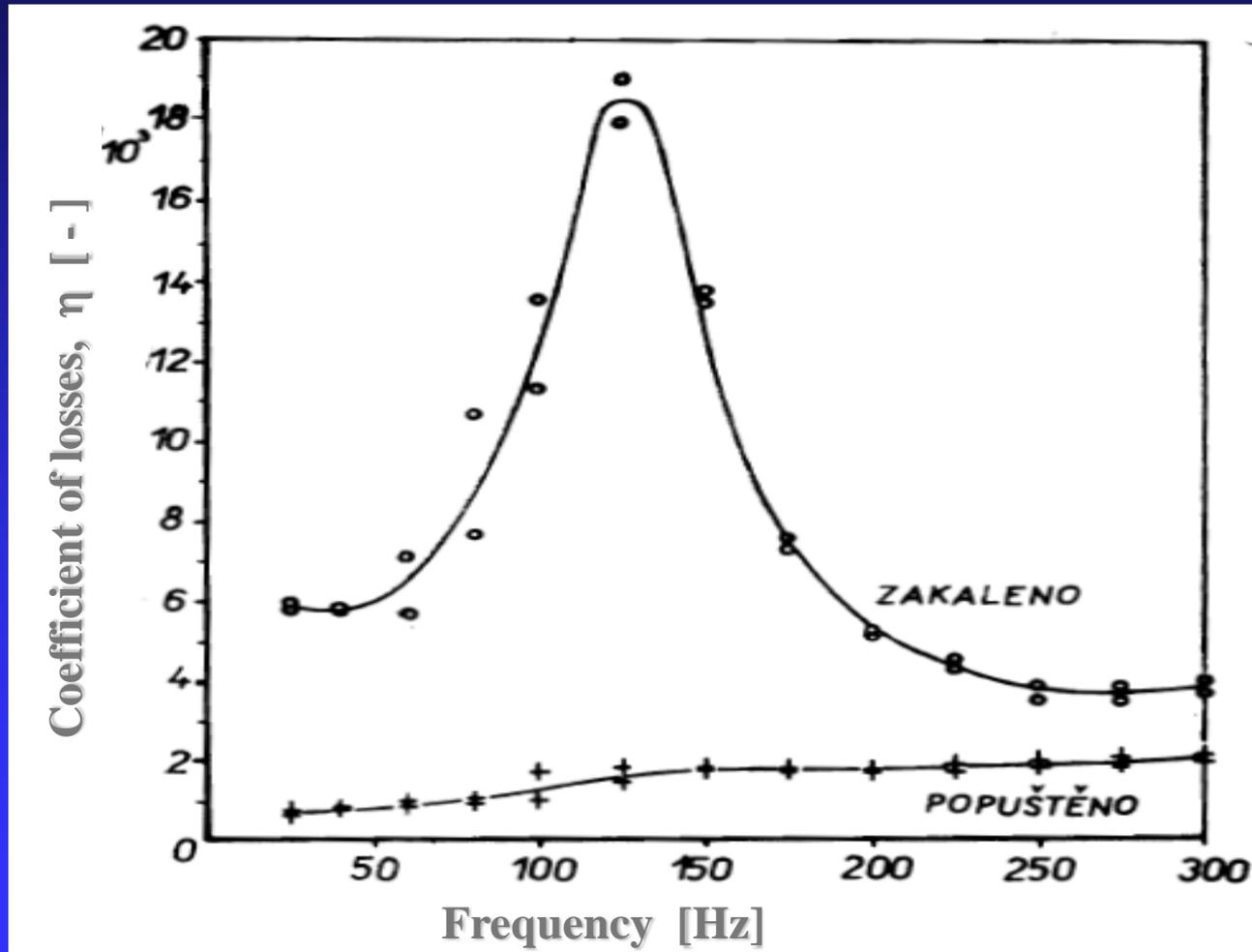
$$\eta = \frac{\Delta U}{2\pi U} \quad (Q^{-1})$$

- logarithmic decrement of dumping

$$\mathcal{D} = \ln \left( \frac{a_n}{a_{n+1}} \right)$$

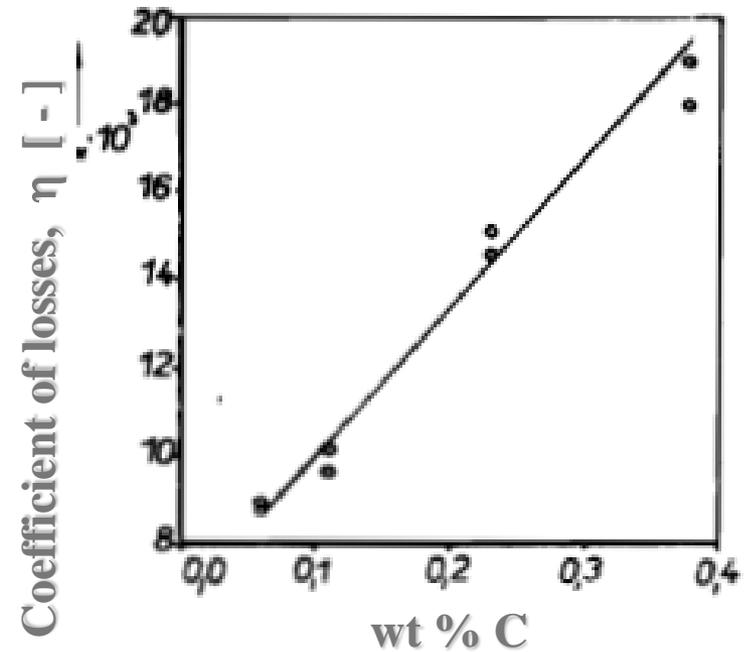
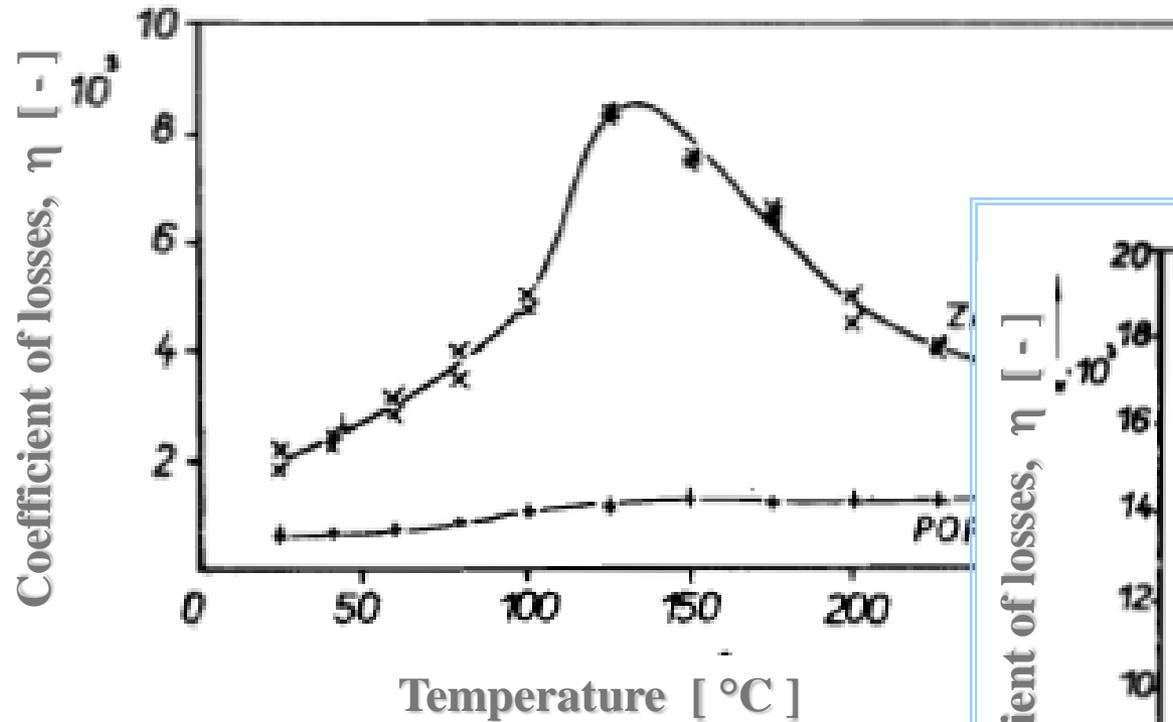
# Measurements of internal damping parameters

Frequency dependence of coefficient of losses



# Measurements of internal damping parameters

Temperature dependence of coefficient of losses



# Measurements of internal damping parameters

- mechanical spectroscopy



# Inelasticity – internal dumping

- Ability of the material dissipate elastic energy during vibrations
- Internal dumping quantification
  - coefficient of internal losses
  - logarithmic decrement of dumping
- *Practical meaning:*
  - gear box, engine block, railway wheels etc.*

# Inelasticity – internal dumping

## Physical background

### Frequency dependent internal dumping

- thermoelastic phenomenon (elongation – cooling, compression - heating)
- motion of valency electrons (at several Kelvins)
- viscous flow - grain boundaries (K<sub>e</sub>, self diffusion)
- carbon diffusion in solid solution of  $\alpha$ -iron (Snoek)
- change of orientation of bi-vacancies etc.
- relaxation of dislocations – Bordoni peak
- dynamic properties of dislocations
- relaxation processes in plastics

# Inelasticity – internal dumping

## Physical background

### Frequency independent internal dumping

- behaviour of dislocations segments (dislocations between anchoring points)
- magnetic-mechanic dumping (magnetic domains rotation)

# Inelasticity – internal dumping

## Importance

## Frequency dependent internal dumping

Interesting for materials scientists rather than mechanical engineers

(phenomena conditioned by some relaxation process):

- carbon diffusion in solid solution of  $\alpha$ -iron

(Snoek)

- relaxation in dislocations (Bordoni)

- dynamic properties of dislocations and interactions of dislocations with obstacles – dislocation, atmospheres, precipitates, stacking faults)

# Inelasticity – internal dumping

## Importance

## Frequency independent internal dumping

Key property for designers

- railway cars (wheels)
- gear box
- noise dumping etc.

## Suitable materials:

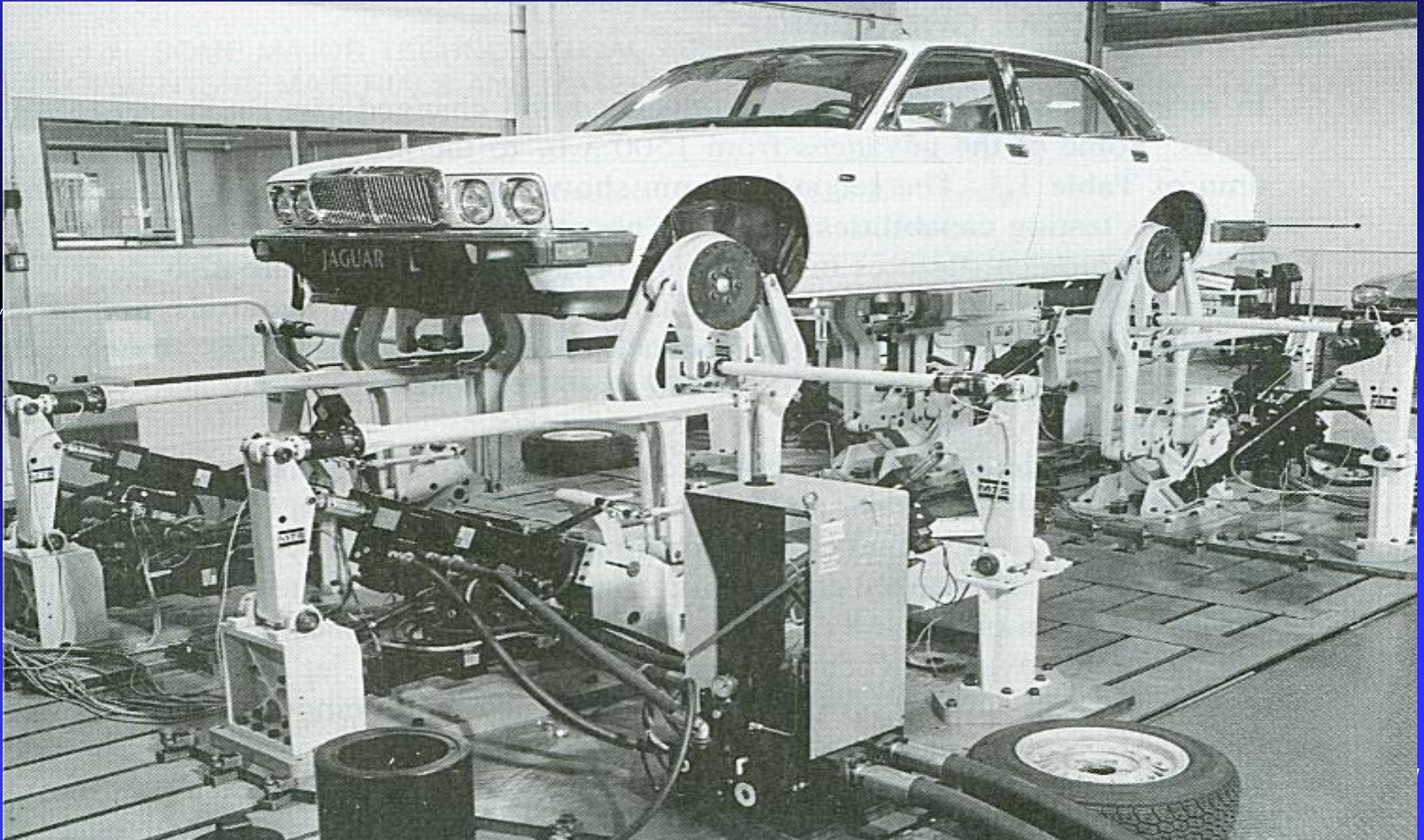
- magnesium alloys – reversible twin deformation
- Mn-Cu alloy – reversible martensitic transformation
- cast iron – energy dissipation on graphite particles
- concrete – energy diss. on inhomogeneous structure

# Inelasticity – internal dumping

## Material selection of vibration table

(vibration table – structures testing for cars, aircrafts)

Material demands:



# Inelasticity – internal dumping

## Material selection of vibration table

(vibration table – structures testing for cars, aircrafts)

Material demands:

low specific weight (density)  $\rho$

high rigidity  $E > 30$  GPa

high eigenfrequency  $(E/\rho)^{1/2}$

high coefficient of internal losses  $\eta \approx 10^{-2}$

Material	$\eta$	$(E/\rho)^{1/2}$	$\rho$ [Mg/m <sup>3</sup> ]	Evaluation
Mg alloys	$10^{-2} - 10^{-1}$	$5 \cdot 10^3$	1,75	best solution
alloy Mn-Cu	$10^{-1}$	$3,5 \cdot 10^3$	8	damping OK, but heavy
KFRP/GFRP	$2 \cdot 10^{-2}$	$6 \cdot 10^3$	1,8	acceptable
cast iron	$2 \cdot 10^{-2}$	$6 \cdot 10^3$	7,8	acceptable/heavy
concrete	$2 \cdot 10^{-2}$	$3,5 \cdot 10^3$	2,5	acceptable

# Inelasticity – internal damping

