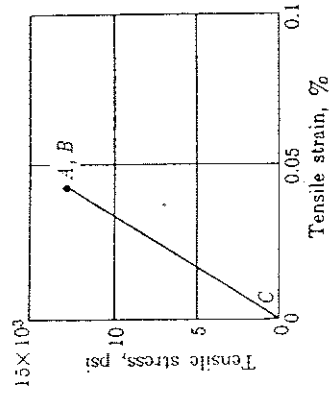


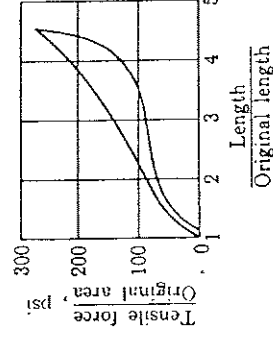
Introduction to Elastic Behavior

- Elastic behavior does not result in permanent deformation upon removal of applied loads
- It is induced primarily by the stretching and rotation of bonds in crystalline and non-crystalline solids
- It may be either linear or non-linear
- It may also be instantaneous or time-dependent

Elastic Deformation of 1020 Steel



Non-linear Deformation of Rubber



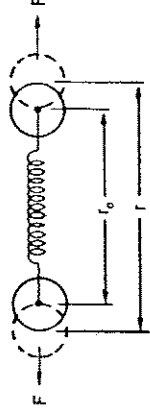
Objectives of this Lecture

- To present an introduction to the basic concepts of elasticity
 - physical basis
 - theoretical basis
- To introduce the anisotropic nature of elastic deformation of crystals
 - single crystals
 - polycrystals

Reasons for Elastic Behavior

- Consider two atoms A and B that are chemically bonded
- The bonds may be either:
 - strong primary bonds (ionic, covalent, metallic)
 - weak secondary bonds (Van der Waal's forces or hydrogen bonds)
- Deformation in early stages controlled by the stretching or rotation of bonds

Stretching of Chemical Bonds Between A and B



Potential Energy and Force Dependence on Atomic Separation

- The potential energy of a bonded system has two components, U_a and U_R :

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

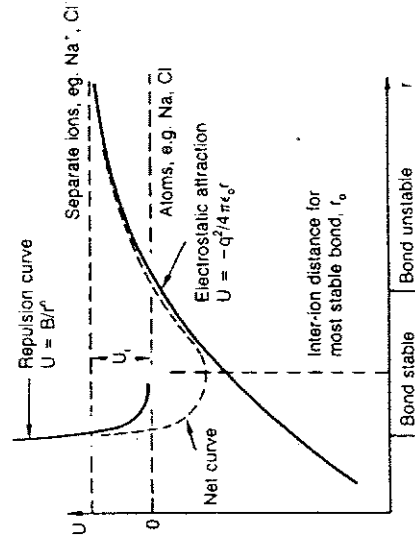
The force is given by the derivative of U

$$F = -\frac{dU}{dr} = -\frac{Am}{r^{m+1}} + \frac{Bn}{r^{n+1}}$$

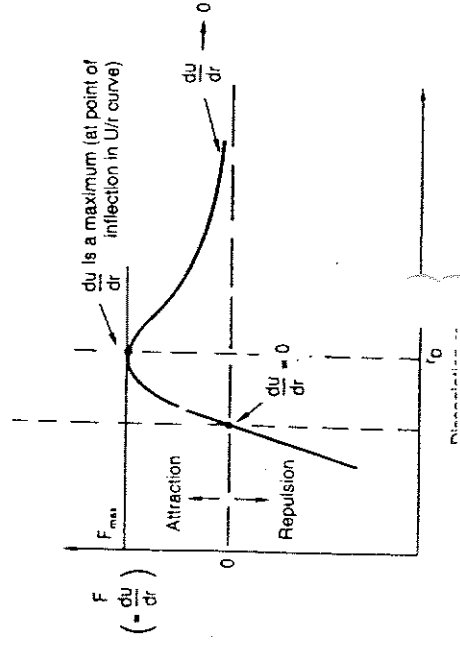
- $F = 0$ corresponds to the equilibrium separation r_0 and binding energy U_0

Schematic of Energy/Force Versus Separation Functions

Relationship Between U and r



Relationship Between F and r



Introduction to Linear Elasticity

- The simple relation between stress and strain was first proposed by Robert Hooke in 1678 (Hooke's law)

$$\sigma = E\varepsilon$$

and

$$\tau = G\gamma.$$

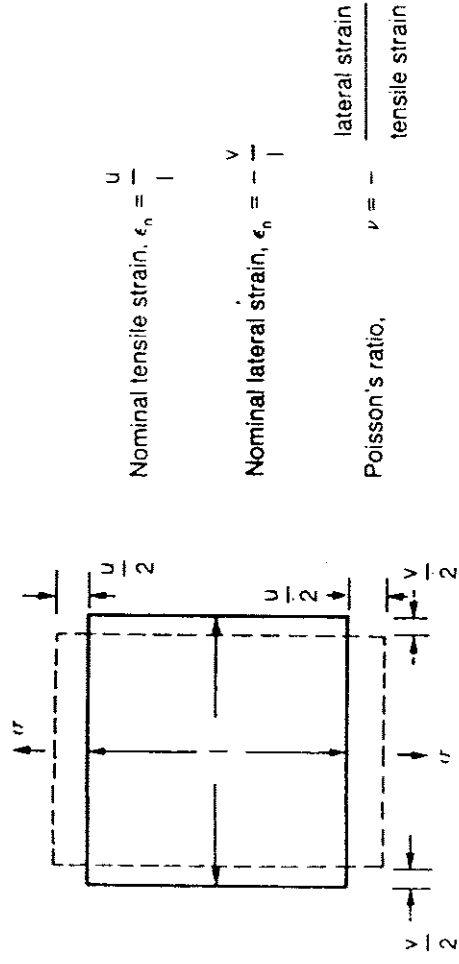
and

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

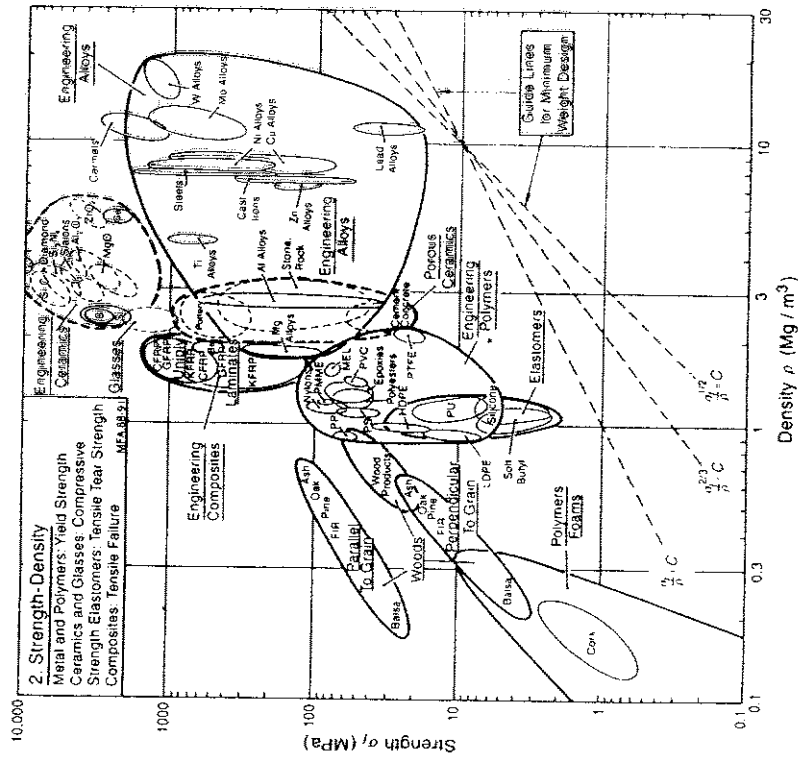
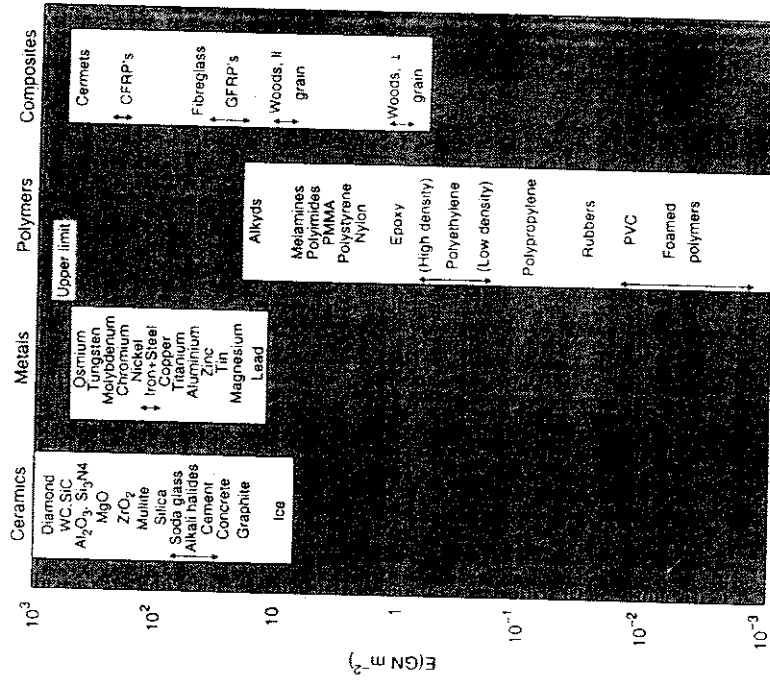
Introduction to Poisson's Ratio

- Lateral contract generally associated with axial deformation and vice-versa
- This was first recognized by Poisson in the 19th century
- Poisson's ratio is defined as

Schematic of Lateral Contraction



Typical Elastic Properties for Different Materials



Crystal Symmetry and Anisotropic Deformation

- There appear to be 36 C_{ij} or S_{ij} terms – at first glance
- However, only 21 of these terms are independent
- Number of independent elastic terms decreases with increasing crystal symmetry
 - least symmetric crystals (triclinic crystals) have 21 independent terms
 - orthorhombic have 9 and hexagonal have 5
 - cubic crystals have 3 independent terms

Hooke's Law for Cubic Crystals

$$\begin{bmatrix} \epsilon_{11} \\ \epsilon_{22} \\ \epsilon_{33} \\ \gamma_{23} \\ \gamma_{13} \\ \gamma_{12} \end{bmatrix} = \begin{bmatrix} S_{11} & S_{12} & S_{12} & 0 & 0 & 0 \\ S_{12} & S_{11} & S_{12} & 0 & 0 & 0 \\ S_{12} & S_{12} & S_{11} & 0 & 0 & 0 \\ 0 & 0 & 0 & S_{44} & 0 & 0 \\ 0 & 0 & 0 & 0 & S_{44} & 0 \\ 0 & 0 & 0 & 0 & 0 & S_{44} \end{bmatrix} \begin{bmatrix} \sigma_{11} \\ \sigma_{22} \\ \sigma_{33} \\ \tau_{23} \\ \tau_{13} \\ \tau_{12} \end{bmatrix}$$

$$\begin{bmatrix} \sigma_{11} \\ \sigma_{22} \\ \sigma_{33} \\ \tau_{23} \\ \tau_{13} \\ \tau_{12} \end{bmatrix} = \begin{bmatrix} C_{11} & C_{12} & C_{12} & 0 & 0 & 0 \\ C_{12} & C_{11} & C_{12} & 0 & 0 & 0 \\ C_{12} & C_{12} & C_{11} & 0 & 0 & 0 \\ 0 & 0 & 0 & C_{44} & 0 & 0 \\ 0 & 0 & 0 & 0 & C_{44} & 0 \\ 0 & 0 & 0 & 0 & 0 & C_{44} \end{bmatrix} \begin{bmatrix} \epsilon_{11} \\ \epsilon_{22} \\ \epsilon_{33} \\ \gamma_{23} \\ \gamma_{13} \\ \gamma_{12} \end{bmatrix}$$

Introducing the Compliance Tensor

- The earlier expression introduced the elastic stiffness tensor
- The compliance tensor is the inverse of the elastic stiffness tensor
- May also write Hooke's law as

$$\begin{bmatrix} \epsilon_{11} \\ \epsilon_{22} \\ \epsilon_{33} \\ \gamma_{23} \\ \gamma_{13} \\ \gamma_{12} \end{bmatrix} = \begin{bmatrix} S_{11} & S_{12} & S_{13} & S_{14} & S_{15} & S_{16} \\ S_{21} & S_{22} & S_{23} & S_{24} & S_{25} & S_{26} \\ S_{31} & S_{32} & S_{33} & S_{34} & S_{35} & S_{36} \\ S_{41} & S_{42} & S_{43} & S_{44} & S_{45} & S_{46} \\ S_{51} & S_{52} & S_{53} & S_{54} & S_{55} & S_{56} \\ S_{61} & S_{62} & S_{63} & S_{64} & S_{65} & S_{66} \end{bmatrix} \begin{bmatrix} \sigma_{11} \\ \sigma_{22} \\ \sigma_{33} \\ \tau_{23} \\ \tau_{13} \\ \tau_{12} \end{bmatrix}$$

where the S_{ij} terms are the elastic compliance terms

Introduction to Anisotropic Linear Elasticity

- Crystal deformation is strongly dependent on crystal orientation
- More complete theory must therefore include anisotropy effects
- Can therefore rewrite simple Hooke's law as

$$\begin{bmatrix} \sigma_{11} \\ \sigma_{22} \\ \sigma_{33} \\ \tau_{23} \\ \tau_{13} \\ \tau_{12} \end{bmatrix} = \begin{bmatrix} C_{11} & C_{12} & C_{13} & C_{14} & C_{15} & C_{16} \\ C_{21} & C_{22} & C_{23} & C_{24} & C_{25} & C_{26} \\ C_{31} & C_{32} & C_{33} & C_{34} & C_{35} & C_{36} \\ C_{41} & C_{42} & C_{43} & C_{44} & C_{45} & C_{46} \\ C_{51} & C_{52} & C_{53} & C_{54} & C_{55} & C_{56} \\ C_{61} & C_{62} & C_{63} & C_{64} & C_{65} & C_{66} \end{bmatrix} \begin{bmatrix} \epsilon_{11} \\ \epsilon_{22} \\ \epsilon_{33} \\ \gamma_{23} \\ \gamma_{13} \\ \gamma_{12} \end{bmatrix}$$

where the C_{ij} terms are the elastic stiffness terms

Summary of Elastic Stiffness and Compliance Coefficients for Cubic Systems

Material	(10^{10} Pa)				$(10^{-11} \text{ Pa}^{-1})$			
	c_{11}	c_{12}	c_{33}	c_{44}	s_{11}	s_{12}	s_{13}	s_{44}
<i>Cubic</i>								
Aluminum	10.82	6.13		2.85	1.57	-0.57		3.51
Copper	16.84	12.14		7.54	1.50	-0.63		1.33
Gold	18.60	15.70		4.20	2.33	-1.07		2.38
Iron	23.70	14.10		11.60	0.80	-0.28		0.86
Lithium fluoride	11.2	4.56		6.32	1.16	-0.34		1.58
Magnesium oxide	29.3	9.2		15.5	0.401	-0.096		0.648
Molybdenum ^b	46.0	17.6		11.0	0.28	-0.08		0.91
Nickel	24.65	14.73		12.47	0.73	-0.27		0.80
Sodium chloride ^b	4.87	1.26		1.27	2.29	-0.47		7.85
Spinel (MgAl_2O_4)	27.9	15.3		15.3	0.585	-0.208		0.654
Titanium carbide ^b	51.3	10.6		17.8	0.21	-0.036		0.561
Tungsten	50.1	19.8		15.14	0.26	-0.07		0.66
Zinc sulfide	10.79	7.22		4.12	2.0	-0.802		2.43
	c_{11}	c_{12}	c_{33}	c_{44}	s_{11}	s_{12}	s_{13}	s_{44}

Estimating Moduli in Different Directions

- The moduli in different directions may be estimated from the stiffness coefficients
- For cubic crystals, the Young's modulus in a given hkl direction is given by

$$\frac{1}{E_{hkl}} = S_{11} - 2 \left[(S_{11} - S_{12}) - \frac{S_{44}}{2} \right] (l_1^2 l_2^2 + l_2^2 l_3^2 + l_3^2 l_1^2)$$

where l_1 , l_2 and l_3 are the direction cosines of the angle between the direction and the (x,y,z) axes, respectively

Summary of Elastic Direction Cosines and Anisotropy Ratios

Direction Cosines

Direction	l_1	l_2	l_3
$\langle 100 \rangle$	1	0	0
$\langle 110 \rangle$	$1/\sqrt{2}$	$1/\sqrt{2}$	0
$\langle 111 \rangle$	$1/\sqrt{3}$	$1/\sqrt{3}$	$1/\sqrt{3}$

Anisotropy Ratios

Metal	Relative Degree of Anisotropy		E_{100} (10^6 psi)	E_{111} (10^6 psi)	$\left[\frac{E_{111}}{E_{100}} \right]$
	$\left[\frac{2(s_{11} - s_{12})}{s_{44}} \right]$	E_{111} (10^6 psi)			
Aluminum	1.219	11.0	9.2	11.0	1.19
Copper	3.203	27.7	9.7	27.7	2.87
Gold	2.857	16.9	6.2	16.9	2.72
Iron	2.512	39.6	18.1	39.6	2.18
Magnesium oxide	1.534	50.8	36.2	50.8	1.404
Spinel ($MgAl_2O_4$)	2.425	52.9	24.8	52.9	2.133
Titanium carbide	0.877	62.2	69.1	62.2	0.901
Tungsten	1.000	55.8	55.8	55.8	1.00

Estimating the Young's Moduli of Polycrystalline Materials

- Simplest approaches simply use a rule of mixture approach (limited accuracy)

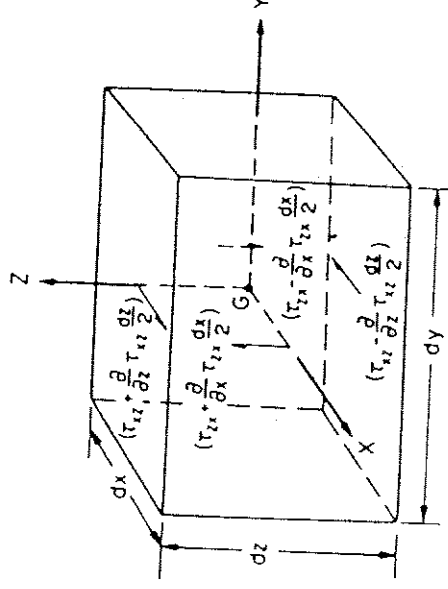
$$E = \sum_{i=1}^n v_i E_i$$

- Bounds in the elastic properties can also be estimated using statistical mechanics or composite theories
- Energy-based approaches provide the best results
- Isotropic elasticity is a true idealization – only true in a few cases e.g. tungsten

Introduction to the Equilibrium Conditions

- The relationships between stress and strain may be determined by the conditions for equilibrium of a cubic element
- Equilibrium can be achieved by considering all the forces that are needed to keep the element suspended in space
- These consist of normal/shear stresses and body forces such as gravitational or centrifugal forces

Equilibrium in the Cartesian Coordinate System



Equilibrium in the Cartesian Coordinate System

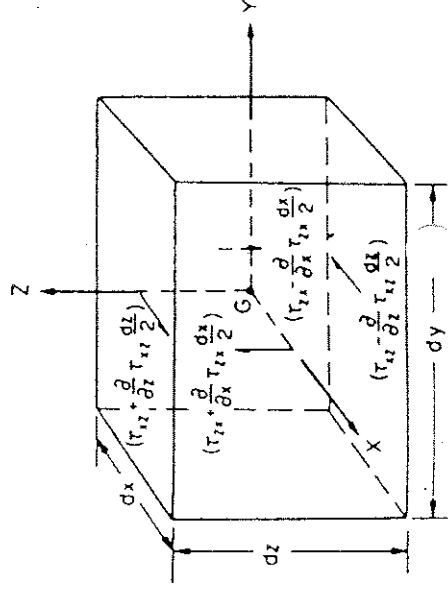
- Considering the gradients in stress across the cube and force equilibrium in the (x,y,z) directions gives:

$$\frac{\partial \sigma_{xx}}{\partial x} + \frac{\partial \tau_{xy}}{\partial y} + \frac{\partial \tau_{xz}}{\partial z} + F_x = \rho \frac{\partial^2 x}{\partial t^2}$$

$$\frac{\partial \tau_{yx}}{\partial x} + \frac{\partial \sigma_{yy}}{\partial y} + \frac{\partial \tau_{yz}}{\partial z} + F_y = \rho \frac{\partial^2 y}{\partial t^2}$$

$$\frac{\partial \tau_{zx}}{\partial x} + \frac{\partial \tau_{zy}}{\partial y} + \frac{\partial \sigma_{zz}}{\partial z} + F_z = \rho \frac{\partial^2 z}{\partial t^2}$$

State of Stress on an Element



Solving for Stress in Equilibrium

- There are 6 stress components and 3 equilibrium equations
- Hence, solving for stress terms requires stress-strain relationships in most problems in linear elasticity
- This can be achieved by
 - simplifying or idealizing the stress states (plane stress versus plane strain)
 - or by solving full 3-D elasticity problem

Equilibrium in Spherical Coordinate System

- Similarly equilibrium conditions may be derived for spherical coordinate system
- Summing forces components in the r , θ and ϕ directions and neglecting body forces gives

$$\frac{\partial \sigma_{rr}}{\partial r} + \frac{1}{r \sin \phi} \frac{\partial \tau_{r\theta}}{\partial \theta} + \frac{1}{r} \frac{\partial \tau_{r\phi}}{\partial \phi} + \frac{2\sigma_{rr} - \sigma_{\theta\theta} - \sigma_{\phi\phi} + \tau_{r\theta} \cot \phi}{r} + F_r = 0$$

$$\frac{\partial \tau_{r\theta}}{\partial r} + \frac{1}{r} \frac{\partial \sigma_{\theta\theta}}{\partial \theta} + \frac{1}{r} \frac{\partial \tau_{\phi\theta}}{\partial \phi} + \frac{3\tau_{\phi\theta}}{r} \cot \phi + F_\theta = 0$$

$$\frac{\partial \tau_{r\phi}}{\partial r} + \frac{1}{r} \frac{\partial \tau_{\phi\theta}}{\partial \theta} + \frac{1}{r} \frac{\partial \sigma_{\phi\phi}}{\partial \phi} + \frac{\sigma_{\phi\phi} - \sigma_{\theta\theta}}{r} \cot \phi + \frac{3\tau_{r\phi}}{r} + F_\phi = 0$$

Equilibrium for Cylindrical and Spherical Coordinate Systems

- For bodies with cylindrical and spherical symmetry, cylindrical and spherical coordinate systems may be used
- Similarly, the equilibrium equations may be derived from force equilibrium
- For cylindrical bodies equilibrium conditions are given by

$$\frac{\partial \sigma_{rr}}{\partial r} + \frac{1}{r} \frac{\partial \tau_{r\theta}}{\partial \theta} + \frac{\partial \tau_{rz}}{\partial z} + \frac{\sigma_{rr} - \sigma_{\theta\theta}}{r} = 0$$

$$\frac{\partial \tau_{r\theta}}{\partial r} + \frac{1}{r} \frac{\partial \sigma_{\theta\theta}}{\partial \theta} + \frac{\partial \tau_{z\theta}}{\partial z} + \frac{2\tau_{r\theta}}{r} = 0$$

$$\frac{\partial \tau_{rz}}{\partial r} + \frac{1}{r} \frac{\partial \tau_{z\theta}}{\partial \theta} + \frac{\partial \sigma_{zz}}{\partial z} + \frac{\tau_{zr}}{r} = 0$$

Equilibrium in Cylindrical Coordinate System

Equilibrium for Stationary Bodies in the Absence of Body Forces

- In the absence of body forces $F_x = F_y = F_z = 0$
- For stationary bodies in equilibrium – acceleration terms

$$\rho \frac{\partial^2 x}{\partial t^2} = \rho \frac{\partial^2 y}{\partial t^2} = \rho \frac{\partial^2 z}{\partial t^2} = 0$$

- Hence, equilibrium equations in Cartesian coordinate system become

$$\frac{\partial \sigma_{xx}}{\partial x} + \frac{\partial \tau_{xy}}{\partial y} + \frac{\partial \tau_{xz}}{\partial z} = 0$$

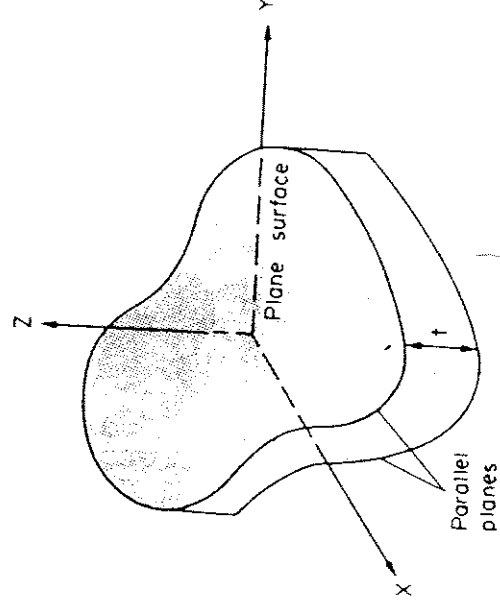
$$\frac{\partial \tau_{yx}}{\partial x} + \frac{\partial \sigma_{yy}}{\partial y} + \frac{\partial \tau_{yz}}{\partial z} = 0$$

$$\frac{\partial \tau_{zx}}{\partial x} + \frac{\partial \tau_{zy}}{\partial y} + \frac{\partial \sigma_{zz}}{\partial z} = 0$$

Plane Stress and Plane Strain Conditions

- In many problems, it is possible to achieve considerable simplification by assuming biaxial stress (plane stress) or biaxial strain (plane strain) conditions
- Plane stress conditions often apply to problems where the thickness is small, while plane strain often applies to thick sections
- However, there are examples in which constraint can lead to plane strain conditions

Schematic of a Plane Element



Introduction to Plane Stress Problems

- In plane stress problems, all the z components of stress are zero

$$\sigma_{zz} = \tau_{xz} = \tau_{yz} = 0$$

- The value of σ_{zz} may also be reduced to zero by imposing an equal stress of opposite sign in the z-direction
- The equations of equilibrium reduce to

$$\frac{\partial \sigma_{xx}}{\partial x} + \frac{\partial \tau_{xy}}{\partial y} + F_x = 0$$

$$\frac{\partial \tau_{xy}}{\partial x} + \frac{\partial \sigma_{yy}}{\partial y} + F_y = 0$$

Equilibrium Conditions Under Plane Strain Conditions

- Under plane strain conditions the strains in the z-direction are zero

$$\epsilon_{zz} = \epsilon_{xz} = \epsilon_{yz} = 0$$

- The equilibrium equations are now given by

$$\frac{\partial \sigma_{xx}}{\partial x} + \frac{\partial \tau_{xy}}{\partial y} + F_x = 0$$

$$\frac{\partial \tau_{xy}}{\partial x} + \frac{\partial \sigma_{yy}}{\partial y} + F_y = 0$$

$$\frac{\partial \sigma_{zz}}{\partial z} + F_z = 0$$

The Relationships Between Stress and Strain Under Plane Stress Conditions

- Under plane stress conditions, the relationships between stress and strain are given by (Hooke's Law):

$$\epsilon_{xx} = \frac{\sigma_{xx}}{E} - \nu \frac{\sigma_{yy}}{E}$$

$$\epsilon_{yy} = \frac{\sigma_{yy}}{E} - \nu \frac{\sigma_{xx}}{E}$$

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

The Relationships Between Strain and Stress Under Plane Strain Conditions

- Under plane strain conditions, the relationships between strain and stress are given by

$$\epsilon_{xx} = \frac{(1-\nu^2)}{E} \left[\sigma_{xx} - \frac{\nu}{(1-\nu)} \sigma_{yy} \right]$$

$$\epsilon_{yy} = \frac{(1-\nu^2)}{E} \left[\sigma_{yy} - \frac{\nu}{(1-\nu)} \sigma_{xx} \right]$$

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

Plane Stress/Strain Relationships Between Stress and Strain

- For plane stress, we have

$$\sigma_{xx} = \frac{E}{(1-\nu^2)} [\epsilon_{xx} + \nu\epsilon_{yy}]$$

$$\sigma_{yy} = \frac{E}{(1-\nu^2)} [\epsilon_{yy} + \nu\epsilon_{xx}]$$

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

- For plane strain, we have

$$\sigma_{xx} = \frac{E(1-\nu)}{(1+\nu)(1-2\nu)} \left[\epsilon_{xx} + \frac{\nu}{(1-\nu)} \epsilon_{yy} \right]$$

$$\sigma_{yy} = \frac{E(1-\nu)}{(1+\nu)(1-2\nu)} \left[\epsilon_{yy} + \frac{\nu}{(1-\nu)} \epsilon_{xx} \right]$$

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

Generalized Three-Dimensional State of Stress

- Thermal stresses may be included on the equilibrium equations by treating them as body forces

$$\frac{\partial \sigma_{xx}}{\partial x} + \frac{\partial \tau_{xy}}{\partial y} + \frac{\partial \tau_{xz}}{\partial z} + \alpha_x \Delta T = 0$$

$$\frac{\partial \tau_{yx}}{\partial x} + \frac{\partial \sigma_{yy}}{\partial y} + \frac{\partial \tau_{yz}}{\partial z} + \alpha_y \Delta T = 0$$

$$\frac{\partial \tau_{zx}}{\partial x} + \frac{\partial \tau_{zy}}{\partial y} + \frac{\partial \sigma_{zz}}{\partial z} + \alpha_z \Delta T = 0$$

- Similar expressions may be written for cylindrical and spherical coordinate systems
- However, equilibrium equations cannot be solved without satisfying compatibility conditions

Compatibility Conditions and Stress Functions

- The compatibility conditions are needed to ensure single-valued displacements
- These are derived in most texts on elasticity
- They are given by equations of the form (for subscripts i and j between 1 and 3):

$$\frac{\partial^2 \varepsilon_{ii}}{\partial x_j^2} + \frac{\partial^2 \varepsilon_{jj}}{\partial x_i^2} = \frac{\partial^2 \gamma_{ij}}{\partial x_i \partial x_j}$$

$$\frac{2\partial^2 \varepsilon_{ii}}{\partial x_j \partial x_k} = \frac{\partial}{\partial x_i} \left(-\frac{\partial \gamma_{jk}}{\partial x_i} + \frac{\partial \gamma_{ki}}{\partial x_j} + \frac{\partial \gamma_{ij}}{\partial x_k} \right)$$

Solving Problems in Linear Elasticity

- Solving problems in linear elasticity can be difficult due to number and type of equations that have to be satisfied
- It is common to use trial and error methods to obtain solutions
- One class of trial functions developed by Sir George Airy is known as Airy functions

$$\sigma_{xx} = \frac{\partial^2 \chi}{\partial y^2}$$

$$\sigma_{yy} = \frac{\partial^2 \chi}{\partial x^2}$$

$$\tau_{xy} = -\frac{\partial^2 \chi}{\partial x \partial y}$$

The Compatibility Condition in Terms of the Airy Stress Function

-
- In Cartesian coordinates, the compatibility condition may be expressed in terms of the Airy stress function

$$\frac{\partial^4 \chi}{\partial x^4} + \frac{2\partial^4 \chi}{\partial x^2 \partial y^2} + \frac{\partial^4 \chi}{\partial y^4} = 0$$

- For stress fields with polar symmetry, the Airy stress functions are given by

$$\sigma_{rr} = \frac{1}{r} \frac{\partial \chi}{\partial r} + \frac{1}{r^2} \frac{\partial^2 \chi}{\partial \theta^2}$$

$$\sigma_{\theta\theta} = \frac{\partial^2 \chi}{\partial r^2}$$

$$\sigma_{r\theta} = -\frac{\partial}{\partial r} \left(\frac{1}{r} \frac{\partial \chi}{\partial \theta} \right)$$

Satisfying the Compatibility Conditions with the Airy Stress Functions

- The compatibility conditions are satisfied with the Airy stress function when Laplace's equation and the biharmonic equation are satisfied

$$\nabla^2 c = 0$$

and

$$\nabla^4 c = 0$$

where ∇^2 is the operator given by

$$\nabla^4 c = 0$$

Summary and Concluding Remarks

- Introduction to elasticity theory presented along with physical basis for elasticity
- Physical basis for elasticity is the stretching and rotation of bonds
- Hooke's law presented for 1D and 3D
- Basic introduction to anisotropic linear elasticity was presented for single crystals and polycrystals
- Equilibrium expressions described along with methods for solving elasticity problems