

# *Texture Components and Euler Angles*

27-731 (normally, 27-750)

Texture, Microstructure & Anisotropy

A.D. (Tony) Rollett

*Last revised: 4<sup>th</sup> Jan. 2020*



# *Lecture Objectives*

- Show how to convert from a description of a crystal orientation based on *Miller indices* to *matrices* to *Euler angles*.
- Give examples of standard named components and their associated Euler angles.
- The overall aim is to be able to describe a ***texture component*** by a ***single point*** (in *orientation space*, which is parameterized with some set of coordinates such as Euler angles) instead of needing to draw the crystal embedded in a reference frame.
- Homework exercises will be converting Miller indices to Euler angles.

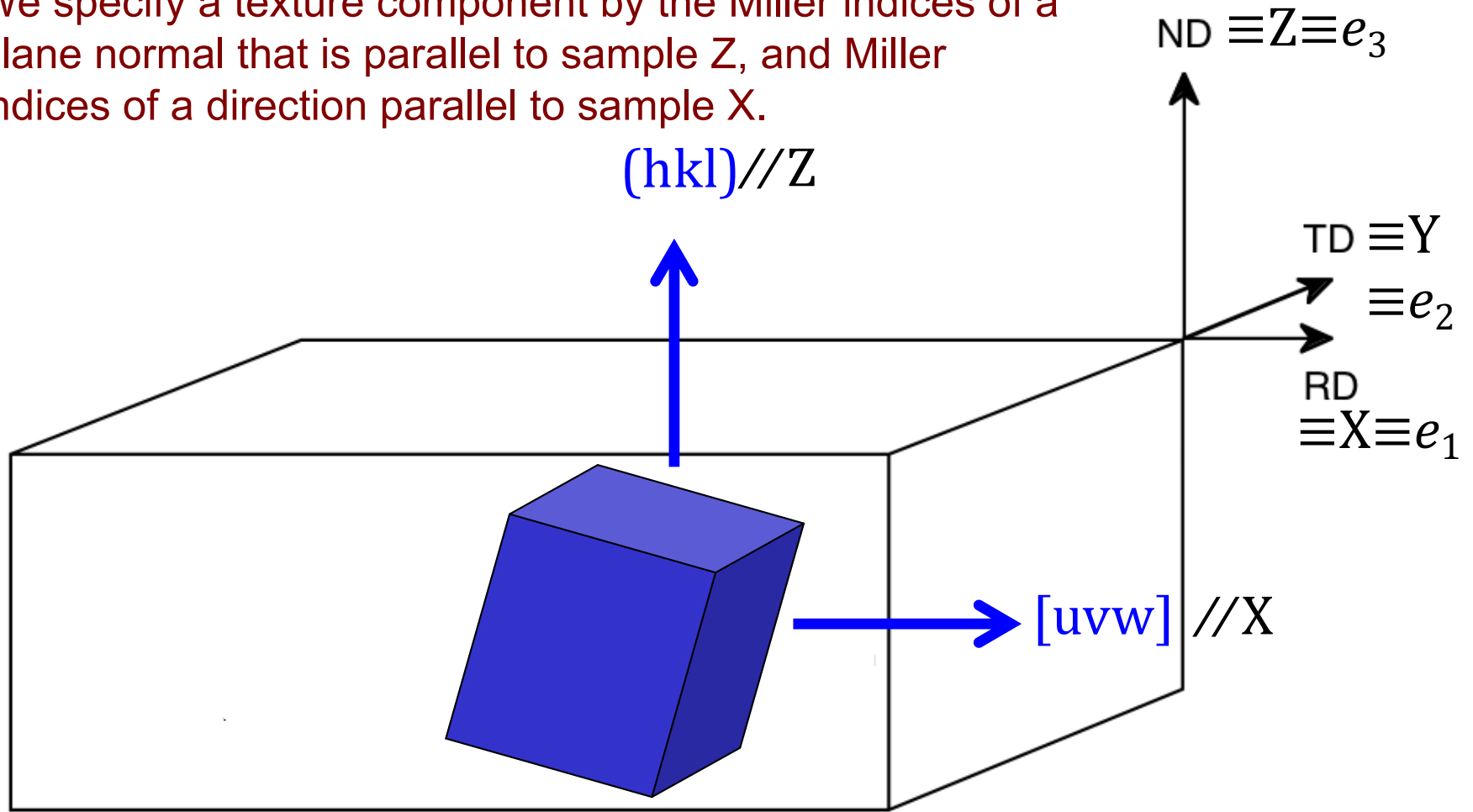
**Obj/notation** AxisTransformation Matrix EulerAngles Components

# *(hkl)[uvw], Miller Index Definition of a Crystal Orientation*

- We use a set of three orthogonal (mutually perpendicular) directions as the reference frame. Mathematicians set up a basis set of **unit vectors** called  $e_1$ ,  $e_2$  and  $e_3$ . All directions can then be described as (linear) combinations of the three basis vectors. See the Supplemental Slides for explanations of *vectors* and *unit vectors*.
- In many cases we use the metallurgical names Rolling Direction (RD) //  $e_1$ , Transverse Direction (TD) //  $e_2$ , and Normal Direction (ND) //  $e_3$ .
- We then identify a crystal (or plane normal) parallel to 3<sup>rd</sup> axis (ND) and a crystal direction parallel to the 1<sup>st</sup> axis (RD), written as  $(hkl)[uvw]$ . The second axis is then completely specified by the other two. *This situation is sometimes referred to as “biaxial texture”, to distinguish it from “uniaxial texture” in which only one crystal axis of each grain is aligned with the specimen.*
- In thin films, we must still use a reference frame with 3 axes but very often, only the normal direction (ND) or film-plane-normal matters because only one crystal direction is aligned with the ND, which is the uniaxial texture referred to above. Be aware, however, that an epitaxial thin film can be biaxially aligned to the substrate in which case all 3 axes are relevant. This latter case is important to High Temperature Superconductor processing, for example.

# Orientation specification via Miller indices: $(hkl)[uvw]$

We specify a texture component by the Miller indices of a plane normal that is parallel to sample Z, and Miller indices of a direction parallel to sample X.



**Obj/notation** AxisTransformation Matrix EulerAngles Components

# *Active versus Passive Rotations*

- Before we discuss the details of how to calculate orientation matrices, it is a good idea to summarize the difference between “active” and “passive” rotations, as mathematicians know them.
- In materials science, we are mostly concerned with describing anisotropic properties of crystals and the aggregate anisotropy of polycrystalline materials, for which it is convenient to use tensors to describe those properties.
- For tensor quantities, we commonly need their coefficients in either the sample frame or the crystal frame. For this we use “transformations of axes”, which are “passive rotations”, in the sense that the two frames share a common origin and differ by only a (proper) rotation. The tensor quantities do not rotate in real space, however.
- In solid mechanics, however, it is more typical to need to describes the motions of objects. Certain motions are just rotations and one can think of rotating a vector, for example, about the origin, in which case one is describing an “active rotation”. Some object is rotated about the origin and moves through the frame.
- When we deal with Miller indices, remember to divide by the magnitude of the vector so as to obtain a unit vector.
- **For all work in texture we will consistently use axis transformations, a.k.a. passive rotations.**

# Why do we need Axis Transformations?

- The reason that we need axis transformations is because in order to compute the anisotropic properties of a polycrystal, we need to be able to transfer (single crystal) properties from the each crystal frame to the sample frame. Equally, we need to transfer stresses, strains and other fields from the sample frame to the crystal frame.
- In order for a quantity to “qualify” as a *tensor* it has to obey the *axis transformation rule*, as discussed in the previous slides.
- The *tensor transformation rule* defines relationships between transformed and untransformed tensors of various ranks.

Vector:

$$V'_i = a_{ij}V_j$$

2<sup>nd</sup> rank

$$T'_{ij} = a_{ik}a_{jl}T_{kl}$$

3<sup>rd</sup> rank

$$T'_{ijk} = a_{il}a_{jm}a_{kn}T_{lmn}$$

4<sup>th</sup> rank

$$T'_{ijkl} = a_{im}a_{jn}a_{ko}a_{lp}T_{mnop}$$

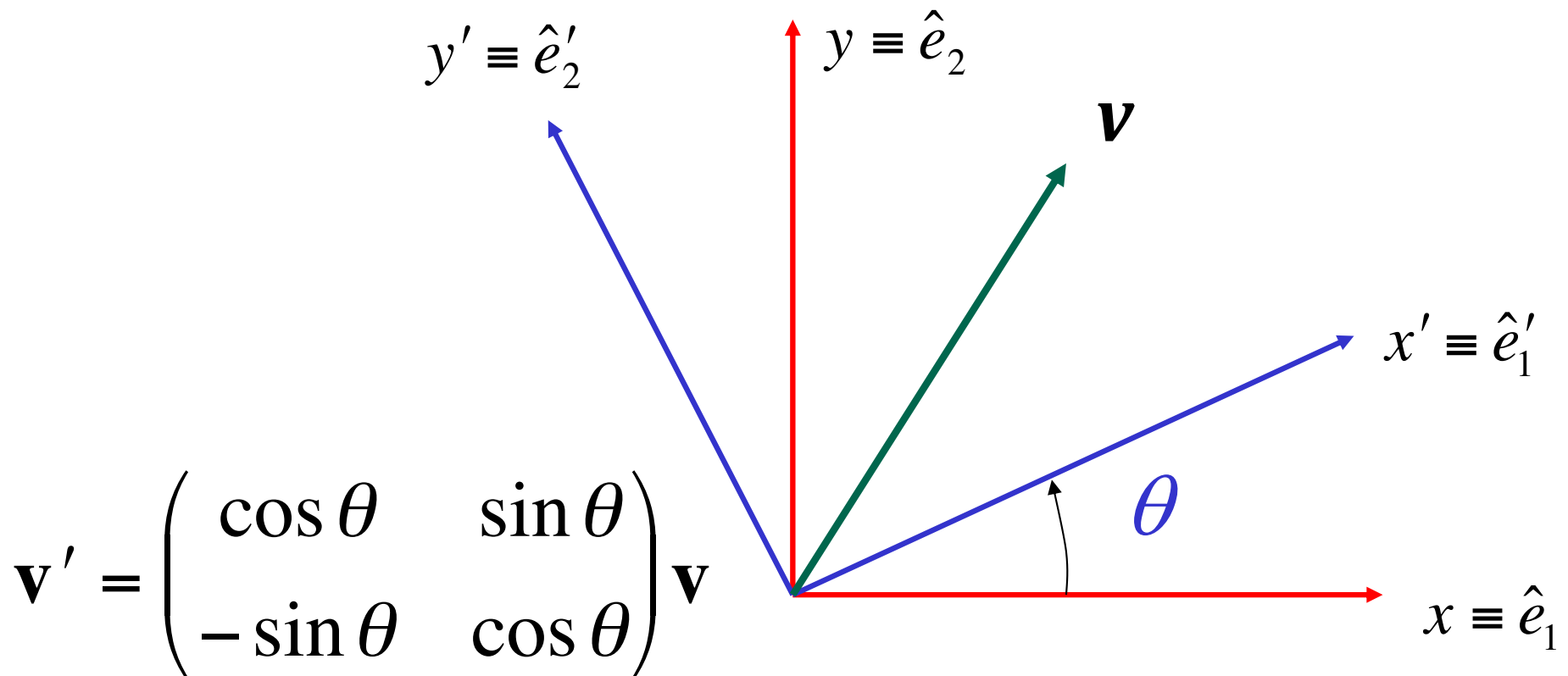
and so on and so forth (without limit in the rank)

7

# Rotation of axes in the (2D) plane:

$x, y = \text{old axes}; x', y' = \text{new axes}$

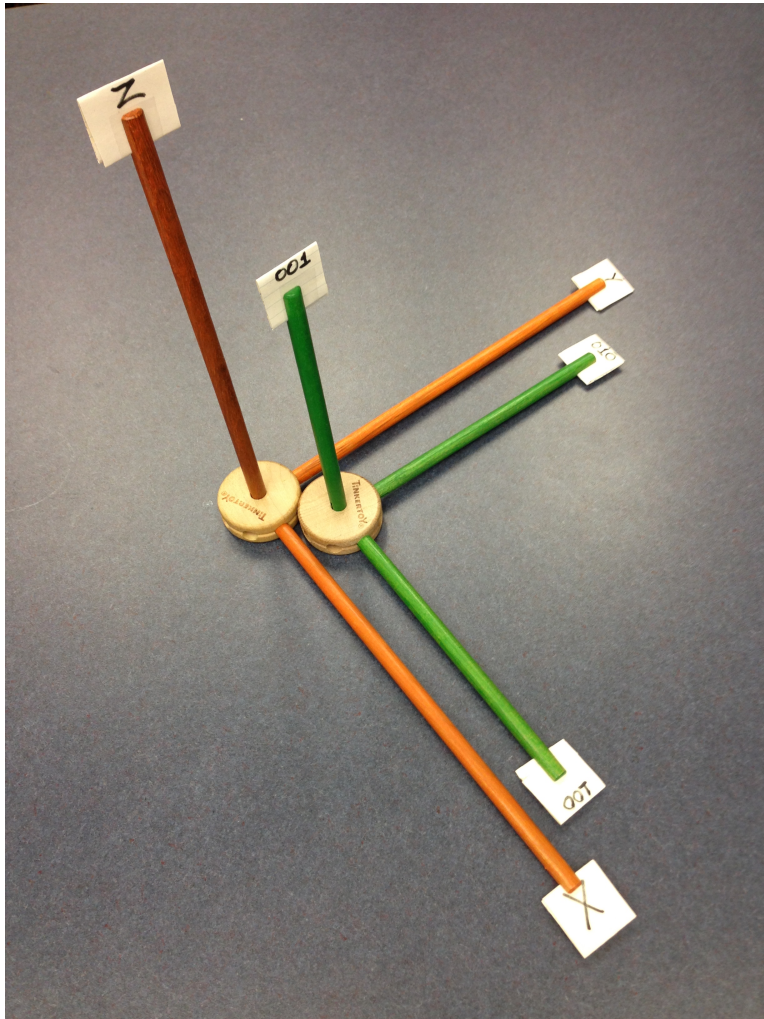
We transform the coefficients of, e.g., a vector,  $\mathbf{v}$ , from one set of axes to another; note that the vector does *not* change position in real space



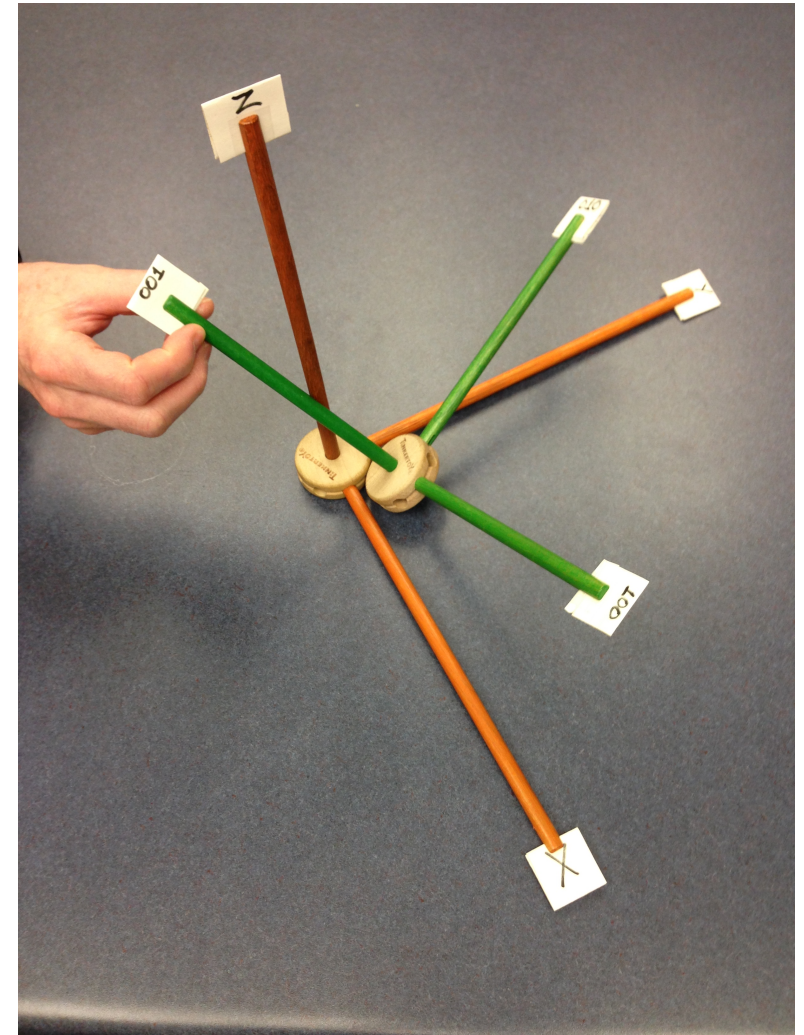
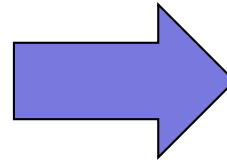
N.B. *Passive Rotation/ Transformation of Axes*

Obj/notation **AxisTransformation** Matrix EulerAngles Components

# Reference Frames



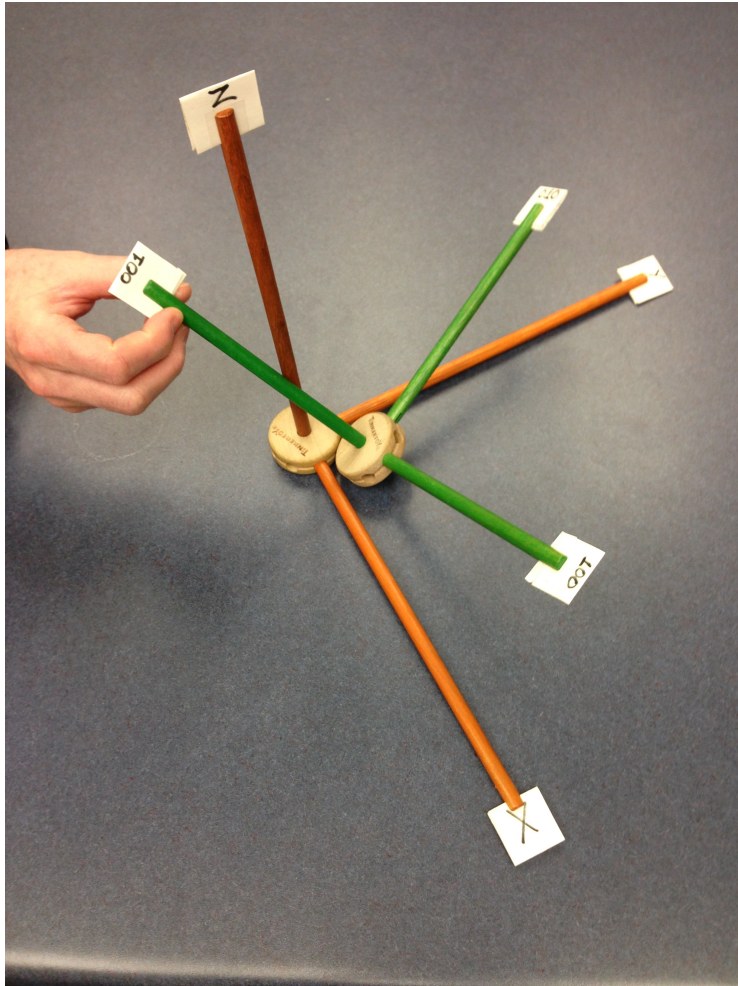
Initial position, crystal frame (green,  $K_B$  in Bunge) aligned with reference (sample,  $K_A$  in Bunge) frame (orange)



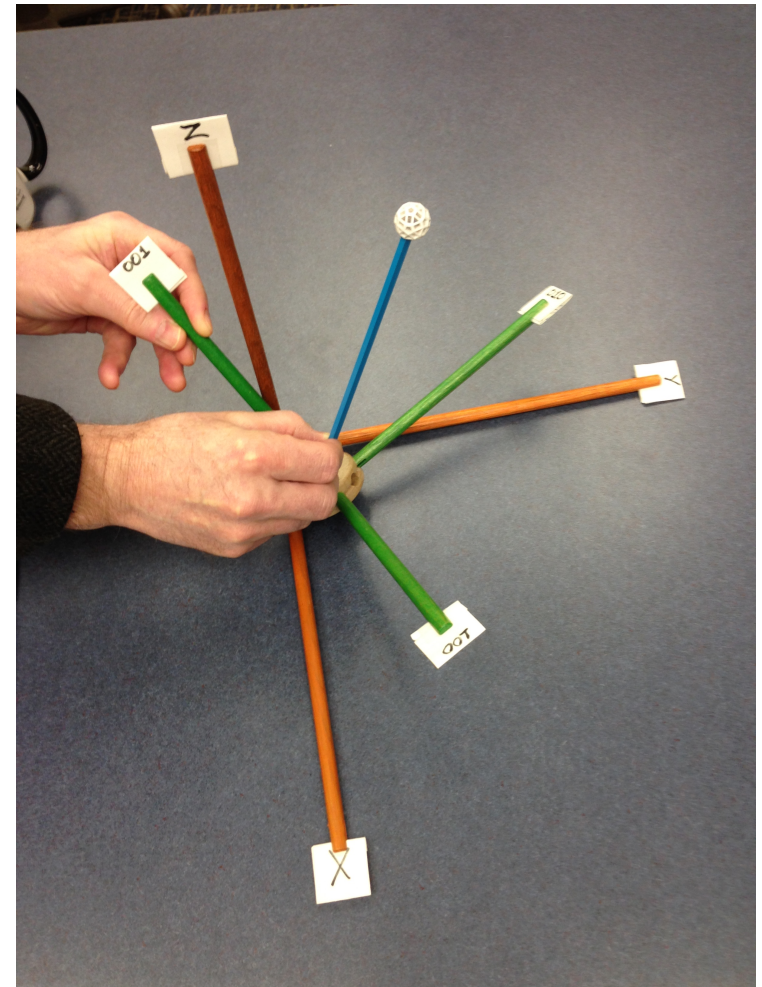
New position, crystal frame (green) rotated with respect to the reference (sample) frame (orange)



# *Vector Represented in Two Different Frames*



New position, crystal frame (green,  $K_B$  in Bunge) rotated with respect to the reference (sample,  $K_A$  in Bunge) frame (orange)



Now we add a vector (the blue stick) that we wish to represent in two different frames (sample vs. crystal)

# Definition of an Axis Transformation:

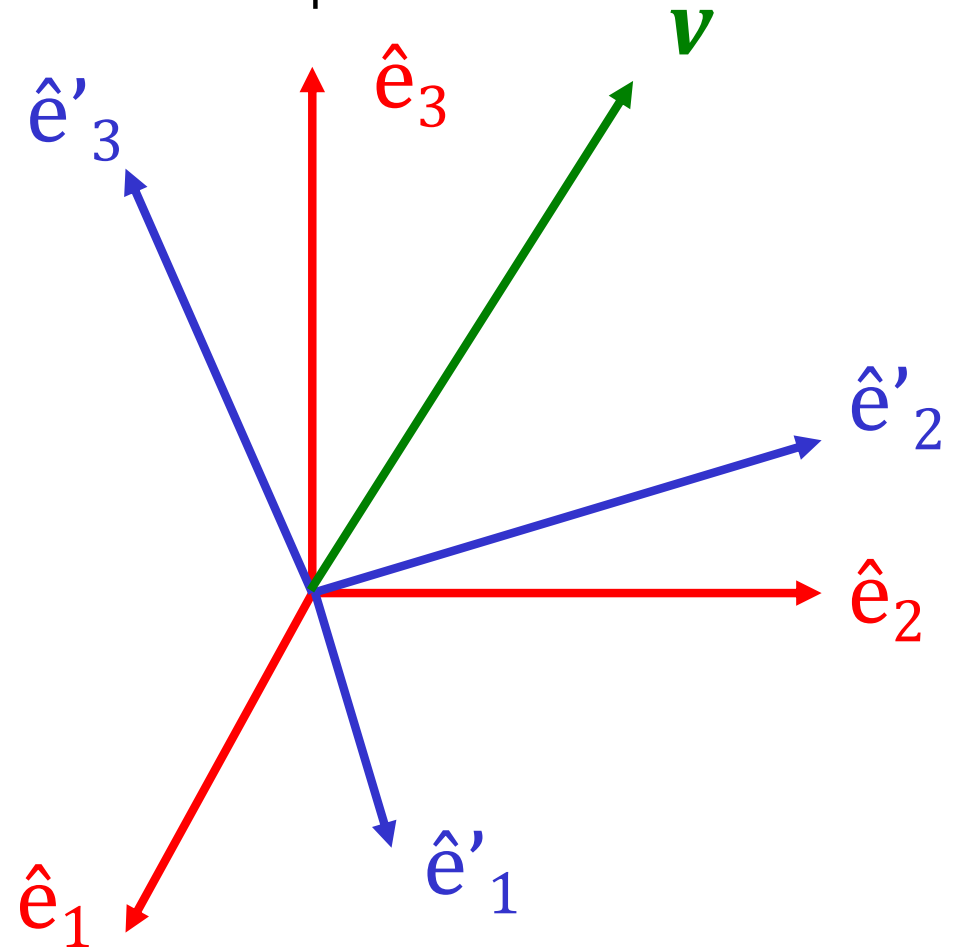
$\hat{e}$  = old axes;  $\hat{e}'$  = new axes

We transform the coefficients of, e.g., a vector,  $v$ , from one set of axes to another; note that the vector does *not* change position in real space

From Sample to  
Crystal (primed)

$$a_{ij} = \hat{e}'_i \bullet \hat{e}_j$$

$$= \begin{pmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{pmatrix}$$

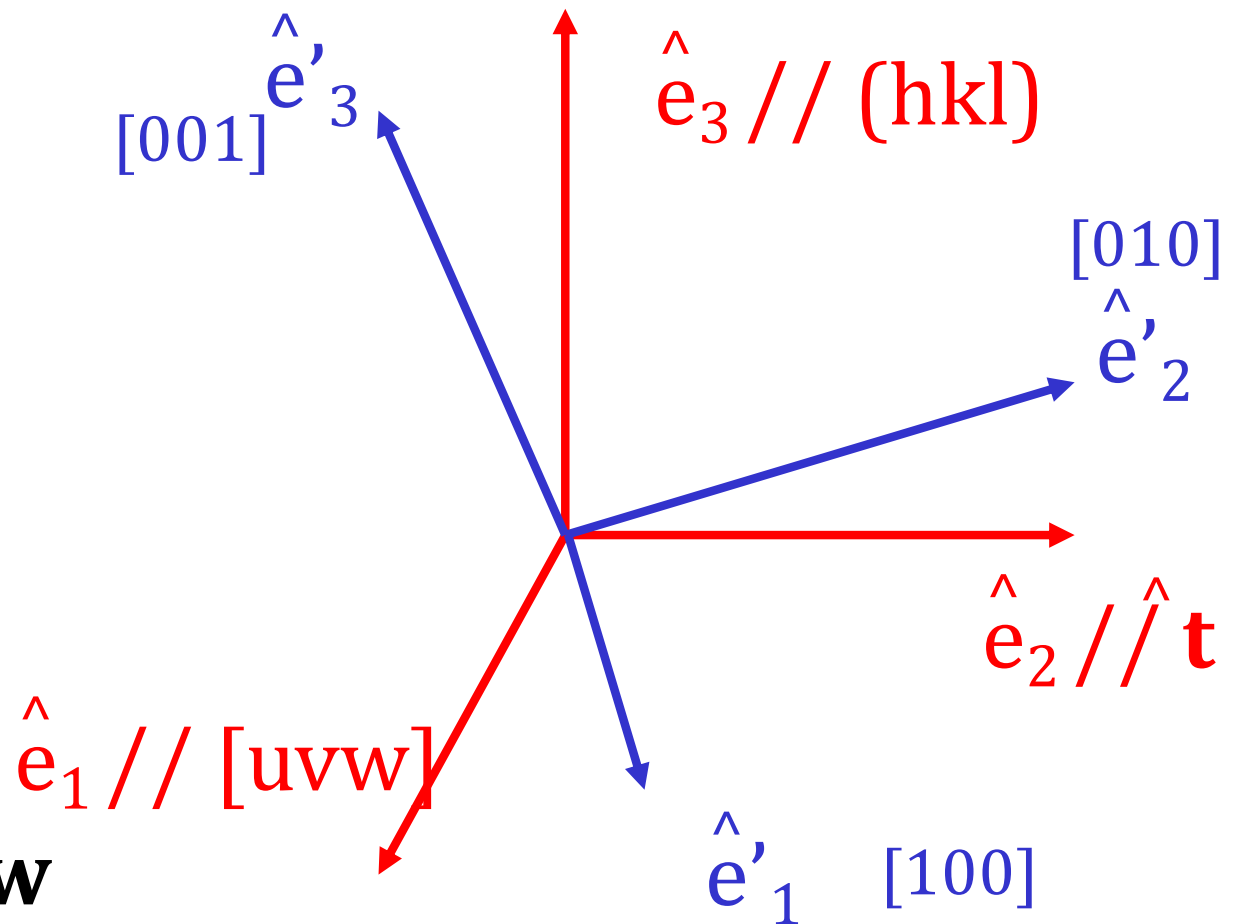


Obj/notation **AxisTransformation** Matrix EulerAngles Components

# Geometry of $\{hkl\}\langle uvw\rangle$

## Sample to Crystal (primed)

Miller index notation of texture component specifies direction cosines of crystal directions // to sample axes. Form the second axis from the cross-product of the 3<sup>rd</sup> and 1<sup>st</sup> axes.



$$\mathbf{t} = \mathbf{hkl} \times \mathbf{uvw}$$

Obj/notation **AxisTransformation** Matrix EulerAngles Components

# Form matrix from Miller Indices

Basic idea: we can construct the complete rotation matrix from two known, easy to determine columns of the matrix. Knowing that we have columns rather than rows is a consequence of the sense of rotation, which is equivalent to the direction in which the axis transformation is carried out.

$$\hat{\mathbf{n}} = \frac{(h, k, l)}{\sqrt{h^2 + k^2 + l^2}}$$

$$\hat{\mathbf{b}} = \frac{(u, v, w)}{\sqrt{u^2 + v^2 + w^2}} \quad \text{Sample}$$

$$\hat{\mathbf{t}} = \frac{\hat{\mathbf{n}} \times \hat{\mathbf{b}}}{|\hat{\mathbf{n}} \times \hat{\mathbf{b}}|} \quad a_{ij} = \text{Crystal}$$

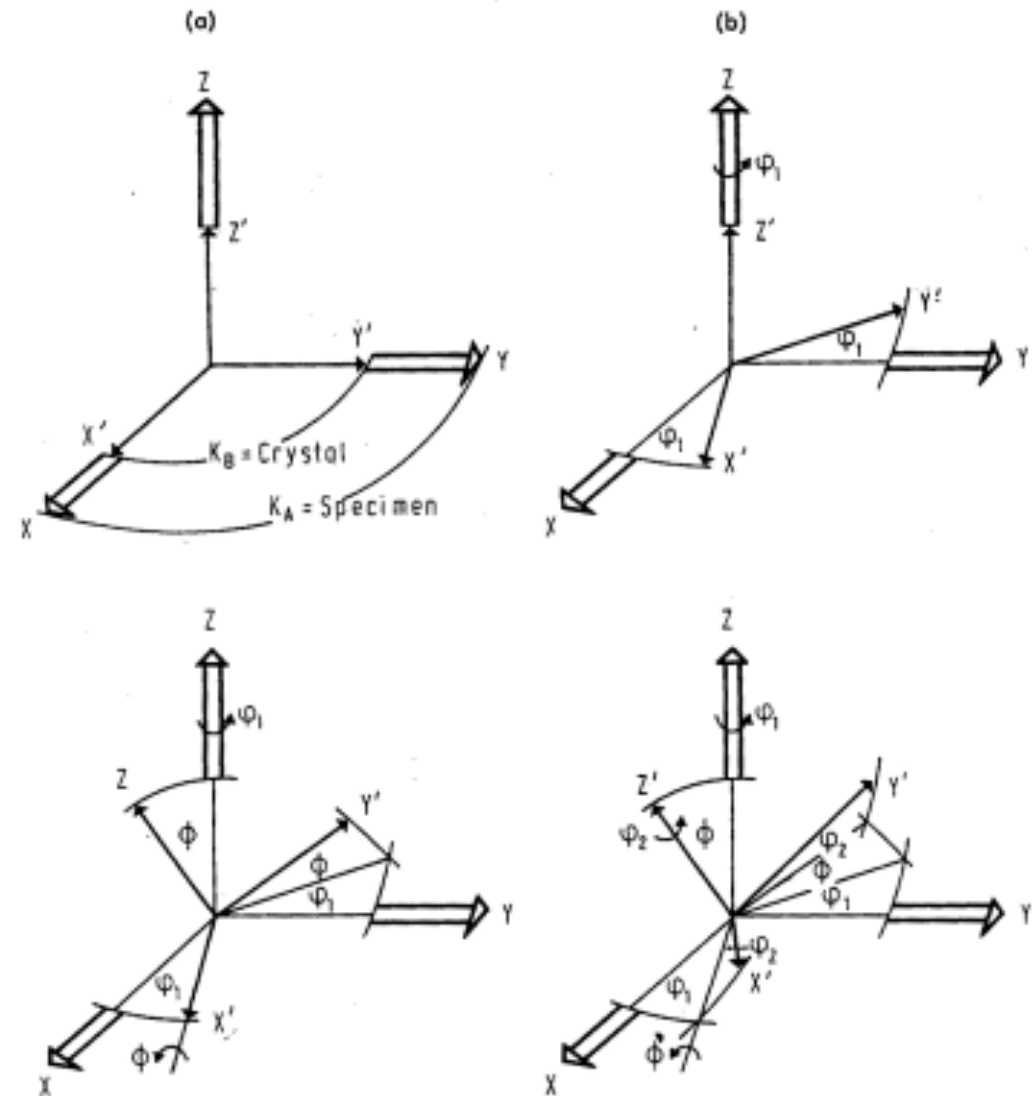
$$\begin{pmatrix} b_1 & t_1 & n_1 \\ b_2 & t_2 & n_2 \\ b_3 & t_3 & n_3 \end{pmatrix}$$

Obj/notation AxisTransformation **Matrix** EulerAngles Components

# (Bunge) Euler Angle Definition

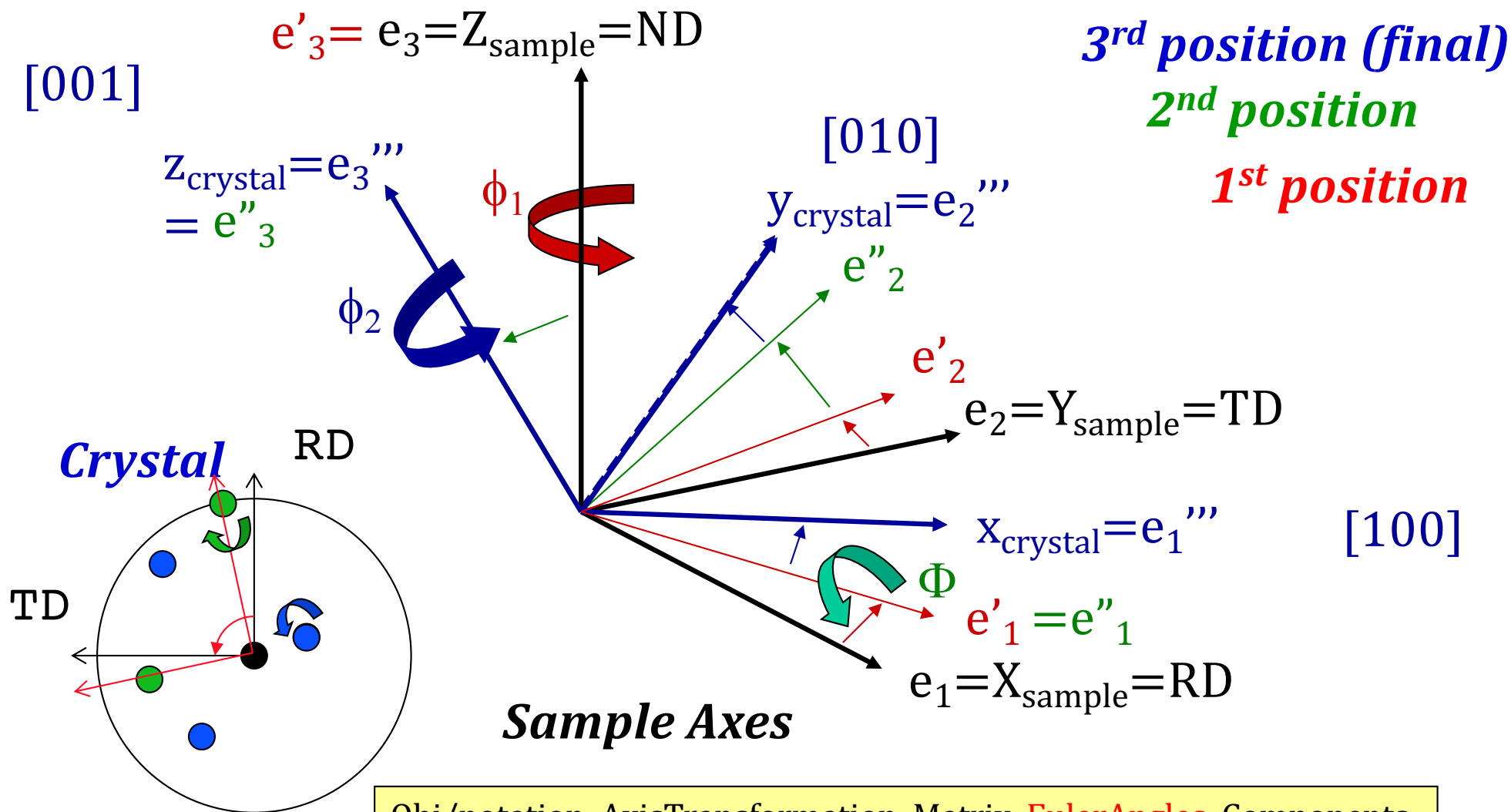
- The three reference axes are labeled as **X, Y & Z**; also commonly known as ND, RD, and TD.
- The three crystal axes are labeled as **X', Y' & Z'**; also commonly known as [100], [010] & [001] in cubic crystals.
- The reference (sample) frame is labeled as " $K_A$ " in the figure, and the crystal frame as " $K_B$ ".
- Each diagram shows successive rotations, more properly thought of as transformations of axes.

[Bunge]



Obj/notation AxisTransformation Matrix EulerAngles Components

# Euler Angles, Animated



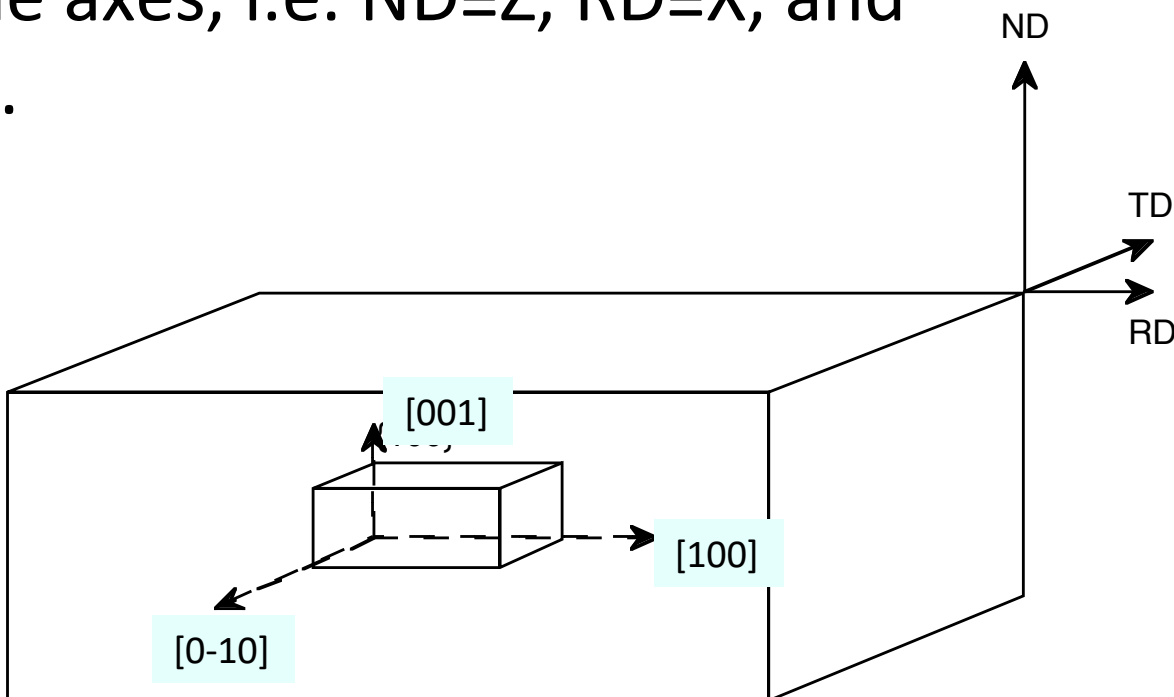
Obj/notation AxisTransformation Matrix EulerAngles Components

# *Named Texture Components*

- In metals, especially rolled fcc metals it is common to give proper names to particular texture components that commonly occur. Thus, you will encounter “cube”, “brass”, “copper”, “S” and the like. Each component has particular Euler angles, as we will now explain. Along with these names go names of so-called fibers such as alpha and beta.
- In rolled bcc metals, it is typical to refer to fibers only, such as the alpha fiber in which the orientations all have  $\langle 110 \rangle // \text{RD}$ , and the gamma fiber for which  $\langle 111 \rangle // \text{ND}$ . One exception to this is the “Goss” component, named after the person who identified the textures critical to transformer steels.
- In hexagonal metals, the texture is mostly evident in the basal pole figure and so the names generally refer to the deviation of the  $\langle 0001 \rangle$  with respect to a particular sample direction such as ND. This is described in more detail later on.
- In thin films of all kinds, unless an epitaxial texture is present, the texture is described in terms of which crystal direction is most commonly found parallel to the normal of the film plane.

# *Cube Texture (001)[100]: cube-on-face*

- Observed in recrystallization of *fcc* metals
- The three crystal axes,  $\langle 001 \rangle$ , are parallel to the three sample axes, i.e.  $ND \equiv Z$ ,  $RD \equiv X$ , and  $TD \equiv Y$  directions.

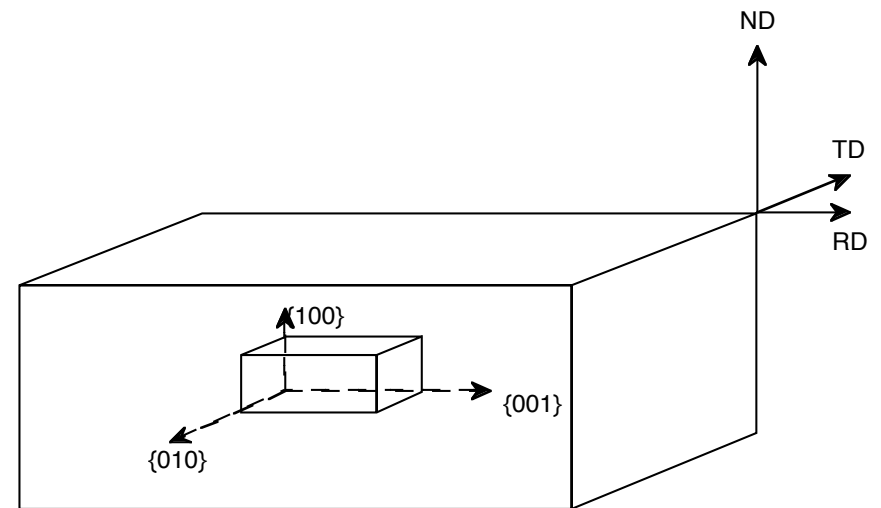
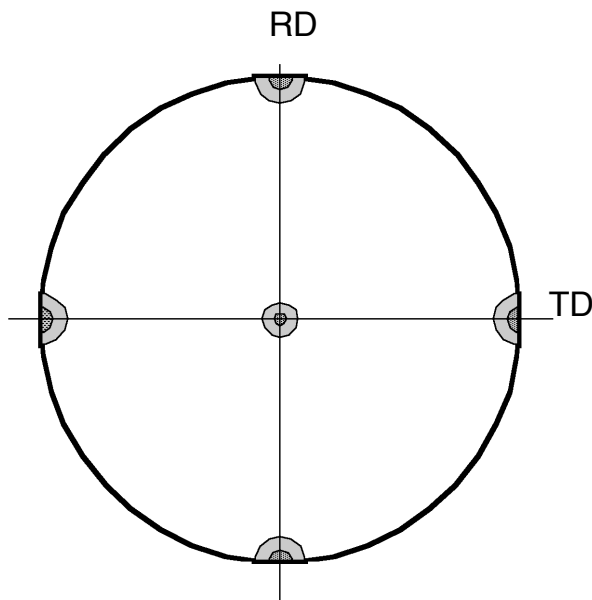


Obj/notation AxisTransformation Matrix EulerAngles **Components**



# Sharp Texture (Recrystallization)

- Figure on the left shows the  $\{001\}$  pole figures (PFs) for this texture component: maxima correspond to  $\{100\}$  poles in the standard stereographic projection aligned with the sample axes. Note that this pole figure could not be measured experimentally unless transmission and reflection methods were to be combined because the reflections on the edge of the PF cannot be detected (see lecture on PF measurement).



Obj/notation AxisTransformation Matrix EulerAngles **Components**

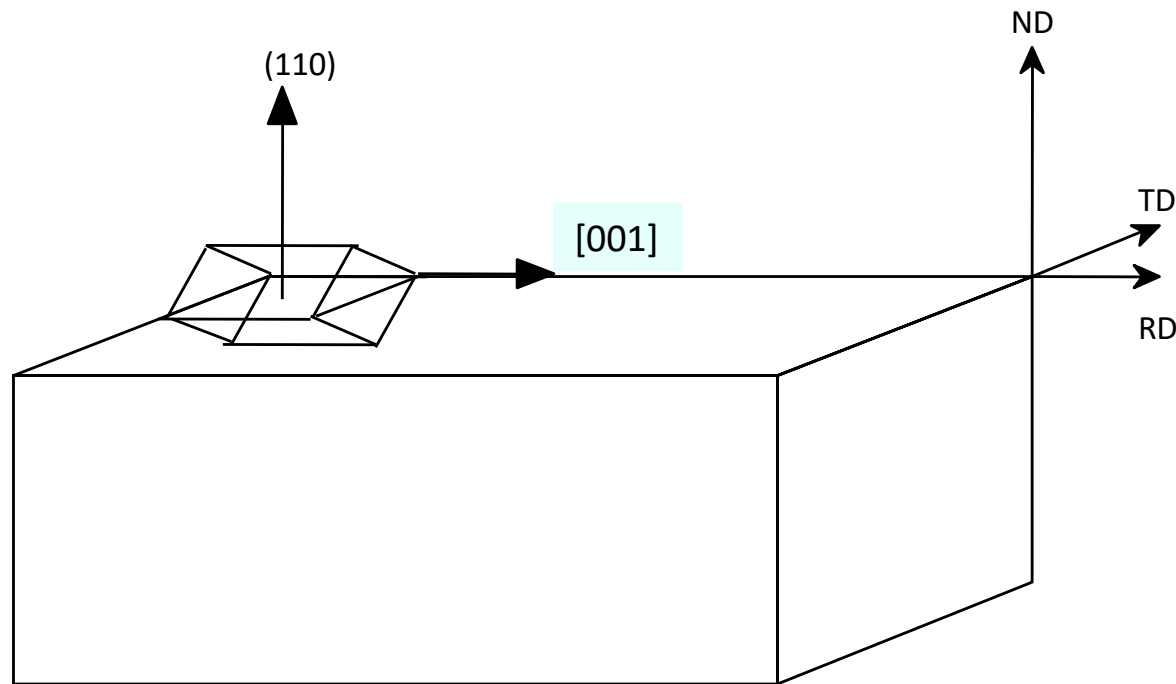
# *Euler angles of Cube component*

- The Euler angles for this component are simple, and yet not so simple!
- The crystal axes align exactly with the specimen axes, therefore all three Euler angles are exactly zero:  
 $(\phi_1, \Phi, \phi_2) = (0^\circ, 0^\circ, 0^\circ)$ .
- Orientation Matrix: 
$$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$
- Rodrigues vector:  $[0,0,0]$
- Unit quaternion:  $[0,0,0,1]$
- As an introduction to the effects of crystal symmetry: consider aligning  $[100]//TD$ ,  $[010]//-RD$ ,  $[001]//ND$ . This is evidently still the cube orientation, but the Euler angles are  
 $(\phi_1, \Phi, \phi_2) = (90^\circ, 0^\circ, 0^\circ)$ !
- This is a first illustration of the confusing (but real) situation where the same physical orientation has multiple mathematical/numerical descriptions. This leads eventually to the discussion of *fundamental zones*.

Obj/notation AxisTransformation Matrix EulerAngles **Components**

# $\{110\}\langle 001\rangle$ : the Goss Component

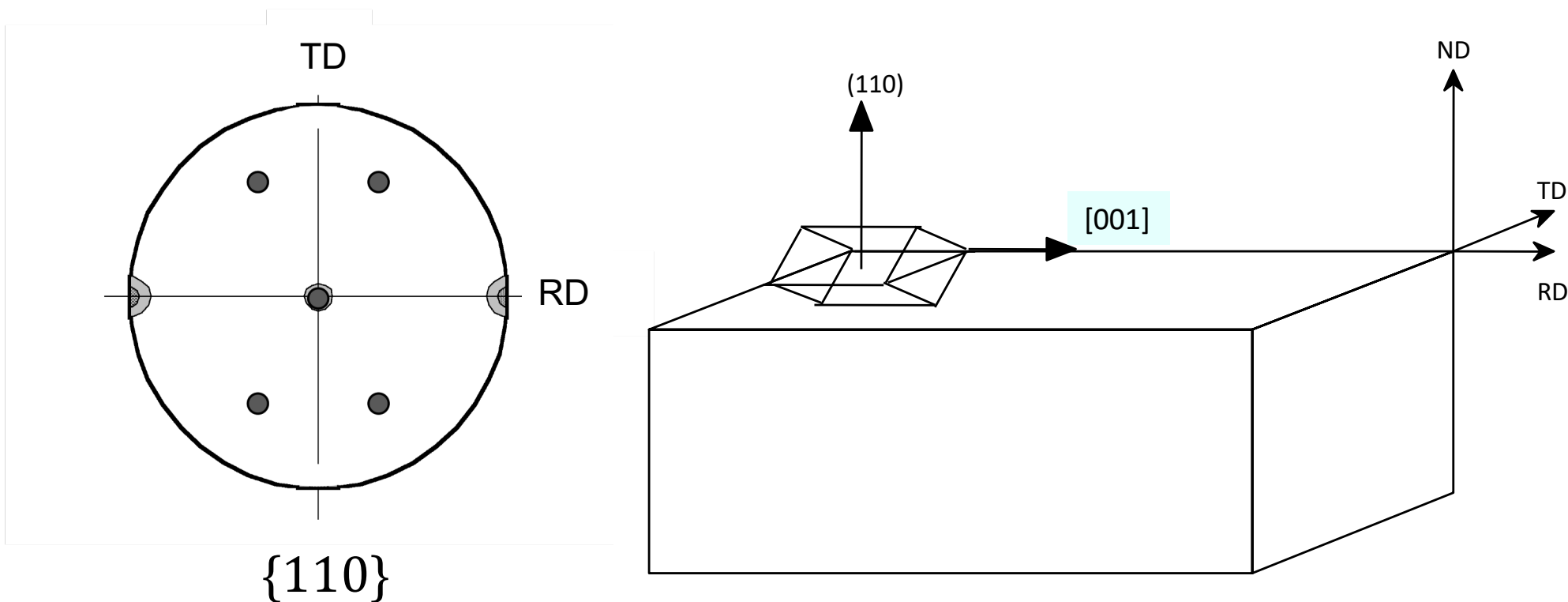
- This type of texture is known as Goss Texture and occurs as a Recrystallization texture for FCC materials such as Brass, ...
- In this case the (011) plane is oriented towards the ND and the [001] inside the (011) plane is along the RD.



Obj/notation AxisTransformation Matrix EulerAngles **Components**

# $\{110\}\langle 001\rangle$ : *cube-on-edge*

- In the 011 pole figure, one of the poles is oriented parallel to the ND (center of the pole figure) but the other ones will be at  $60^\circ$  or  $90^\circ$  angles but tilted  $45^\circ$  from the RD! (Homework: draw the (111) pole figure by hand.)



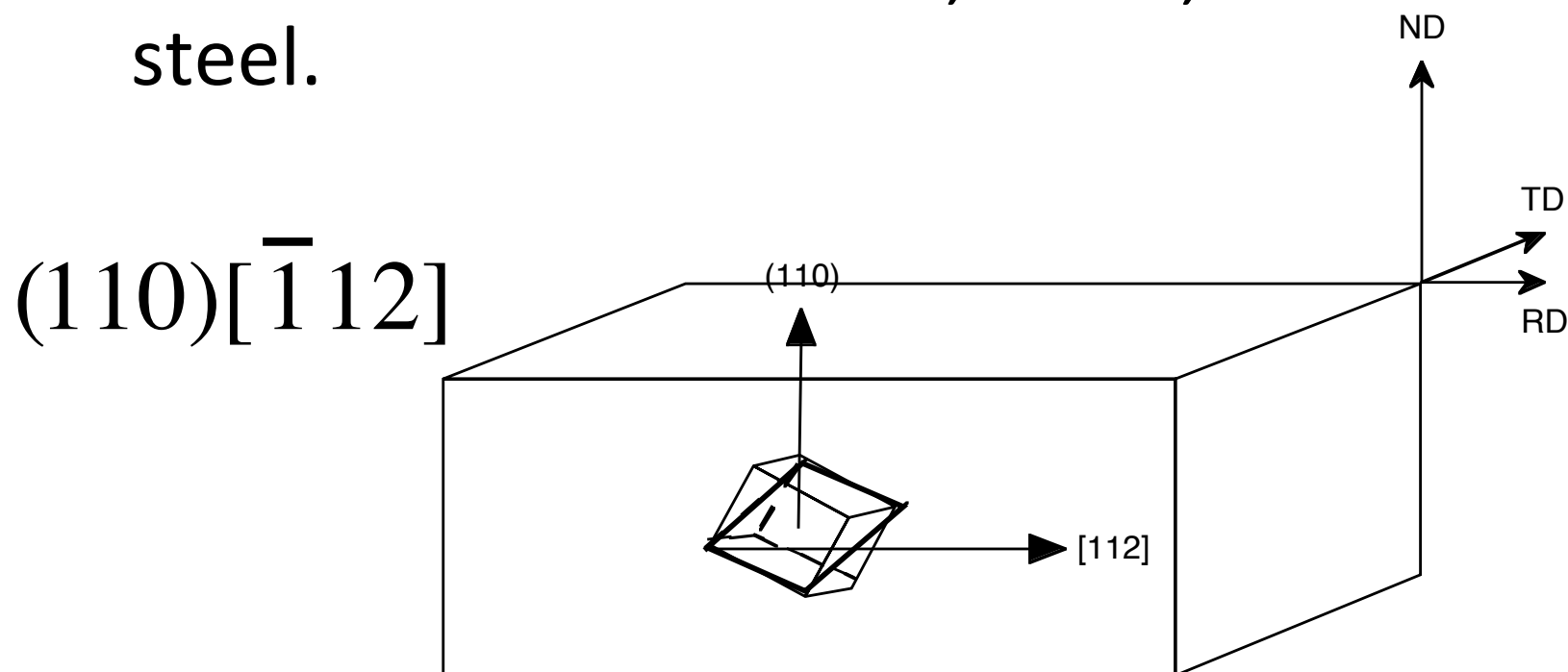
# *Euler angles of Goss component*

- The Euler angles for this component are simple, and yet other variants exist, just as for the cube component.
- Only one rotation of  $45^\circ$  is needed to rotate the crystal from the reference position (i.e. the cube component) to (011)[100]; this happens to be accomplished with the 2nd Euler angle.
- $(\phi_1, \Phi, \phi_2) = (0^\circ, 45^\circ, 0^\circ)$ .
- Other variants will be shown when *symmetry* is discussed.
- Matrix: 
$$\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1/\sqrt{2} & 1/\sqrt{2} \\ 0 & -1/\sqrt{2} & 1/\sqrt{2} \end{pmatrix}$$
- Rodrigues vector:  
[  $\tan(22.5^\circ)$ , 0, 0 ]
- Unit quaternion ( $q_1, q_2, q_3, q_4$ ):  
[  $\sin(22.5^\circ)$ , 0, 0,  $\cos(22.5^\circ)$  ]
- Note that, since there is only one non-zero Euler angle, the rotation axis is obvious by inspection, i.e. the x-axis. For more general cases, the rotation axis has to be calculated.

Obj/notation AxisTransformation Matrix EulerAngles **Components**

## *Brass component*

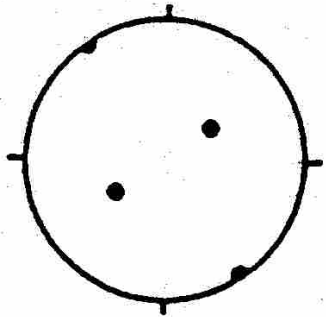
- This type of texture is known as Brass Texture and occurs as a rolling texture component for materials such as Brass, Silver, and Stainless steel.



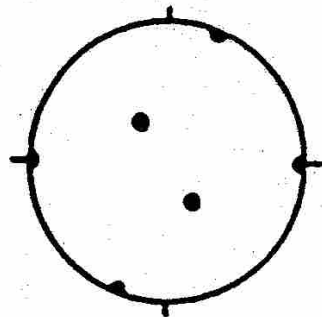
Obj/notation AxisTransformation Matrix EulerAngles **Components**

## *Brass component, contd.*

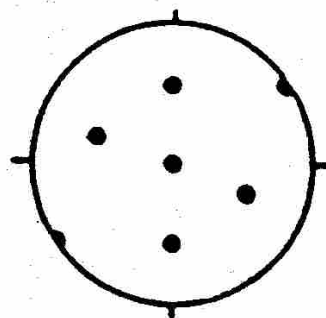
- The associated (110) pole figure is very similar to the Goss texture pole figure except that it is rotated about the ND. In this example, the crystal has been rotated in only one sense (anticlockwise).



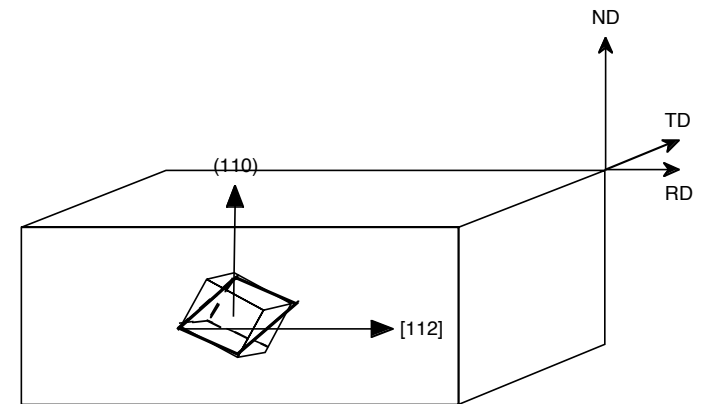
(100)



(111)



(110)

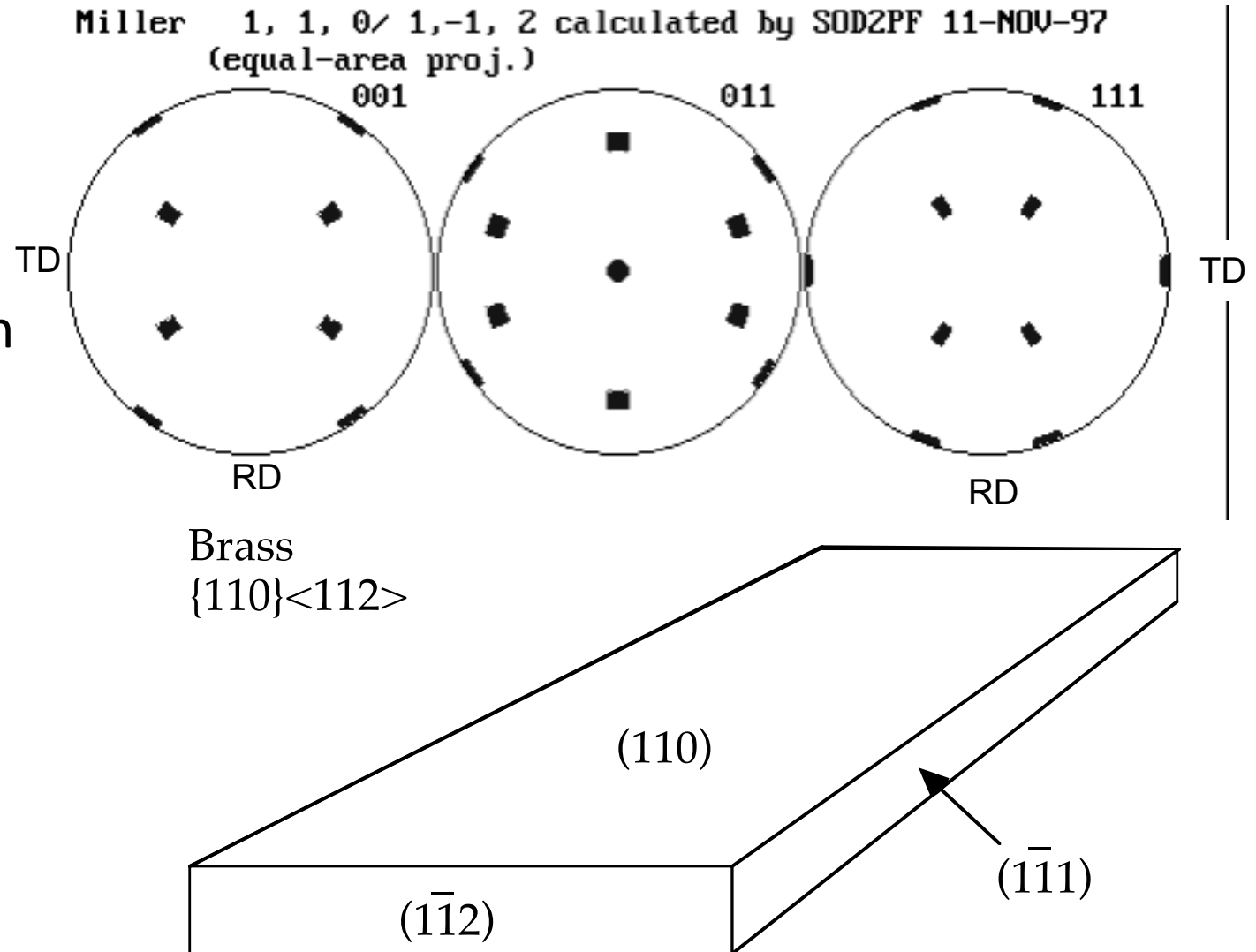


(110)[ $\bar{1}12$ ]

Obj/notation AxisTransformation Matrix EulerAngles **Components**

# $\{110\}\langle 112\rangle$ Brass component

Think of rotating the Goss component around the ND. In this example, the xtal has been rotated in both clockwise and anti-clockwise (two variants).





## *Brass component: Euler angles*

- The brass component is convenient because we can think about performing two successive rotations:
- 1st about the ND, 2nd about the new position of the [100] axis.
- 1st rotation is  $35^\circ$  about the ND; 2nd rotation is  $45^\circ$  about the [100].
- $(\phi_1, \Phi, \phi_2) = (35^\circ, 45^\circ, 0^\circ)$ .

# *Euler Angle Conventions*

- An inconvenient fact is that the definition of Euler angles that we have given so far is *not* unique.
- Many other variants not only exist but have names and are in regular use! Euler angles are used in physics, aeronautics, robotics, to name but a few other fields.
- The differences between the conventions lie in the choice of the rotation axes (generally only the second axis) and the sense of rotation.
- Some of the commonly used conventions are *Bunge*, *Kocks*, *Roe* and *Canova*\* (in approximate order of decreasing popularity).
- **For the purposes of this course we will only use the Bunge convention. At this point in time, almost all codes assume this. Nevertheless, it is wise to check what is being used. The VPSC code, e.g., requires that the input list of orientations has a letter in column 1 on line 4 to define the convention; use “B” in this position to select the Bunge convention.**

\*You will find Canova angles inside the computer code LApp

# Meaning of Bunge Euler angles

- In the *Bunge* convention, the first two angles,  $\phi_1$  and  $\Phi$ , tell you the position of the [001] crystal direction relative to the specimen axes.
- Think of rotating the crystal about the ND (1st angle,  $\phi_1$ ); then rotate the crystal out of the plane (about the [100] axis,  $\Phi$ );
- Finally, the 3<sup>rd</sup> angle ( $\phi_2$ ) tells you how much to rotate the crystal about [001].

# *Meaning of Roe Euler angles*

- In the *Roe* convention, the first two angles,  $\Psi$  and  $\Theta$ , tell you the position of the [001] crystal direction relative to the specimen axes.
- Then think of rotating the crystal about the ND or 001 or z-axis (1st angle,  $\Psi$ ); then rotate the crystal out of the plane (about [010], or the y-axis,  $\Theta$ );
- Finally, the 3<sup>rd</sup> angle ( $\Phi$ ) tells you how much to rotate the crystal about [001].

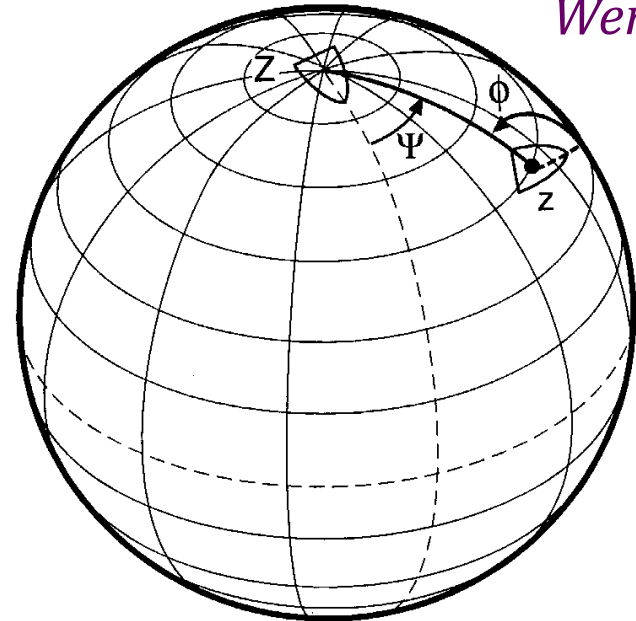
# Meaning of Kocks Euler angles

- In the *Kocks* convention, which is almost the same as the *Roe* convention with the exception of the third angle, the first two angles,  $\Psi$  and  $\Theta$ , tell you the position of the [001] crystal direction relative to the specimen axes.
- Then think of rotating the crystal about the ND or 001 or z-axis (1st angle,  $\Psi$ ); then rotate the crystal out of the plane (about [010], or the y-axis,  $\Theta$ );
- Finally, the 3<sup>rd</sup> angle ( $\phi$ ) tells you how much to rotate the crystal *anti-clockwise* about [001]. Here's the difference between *Roe* and *Kocks* - the last angle is a negative or clockwise rotation about 001, instead of a positive or counter-clockwise rotation.

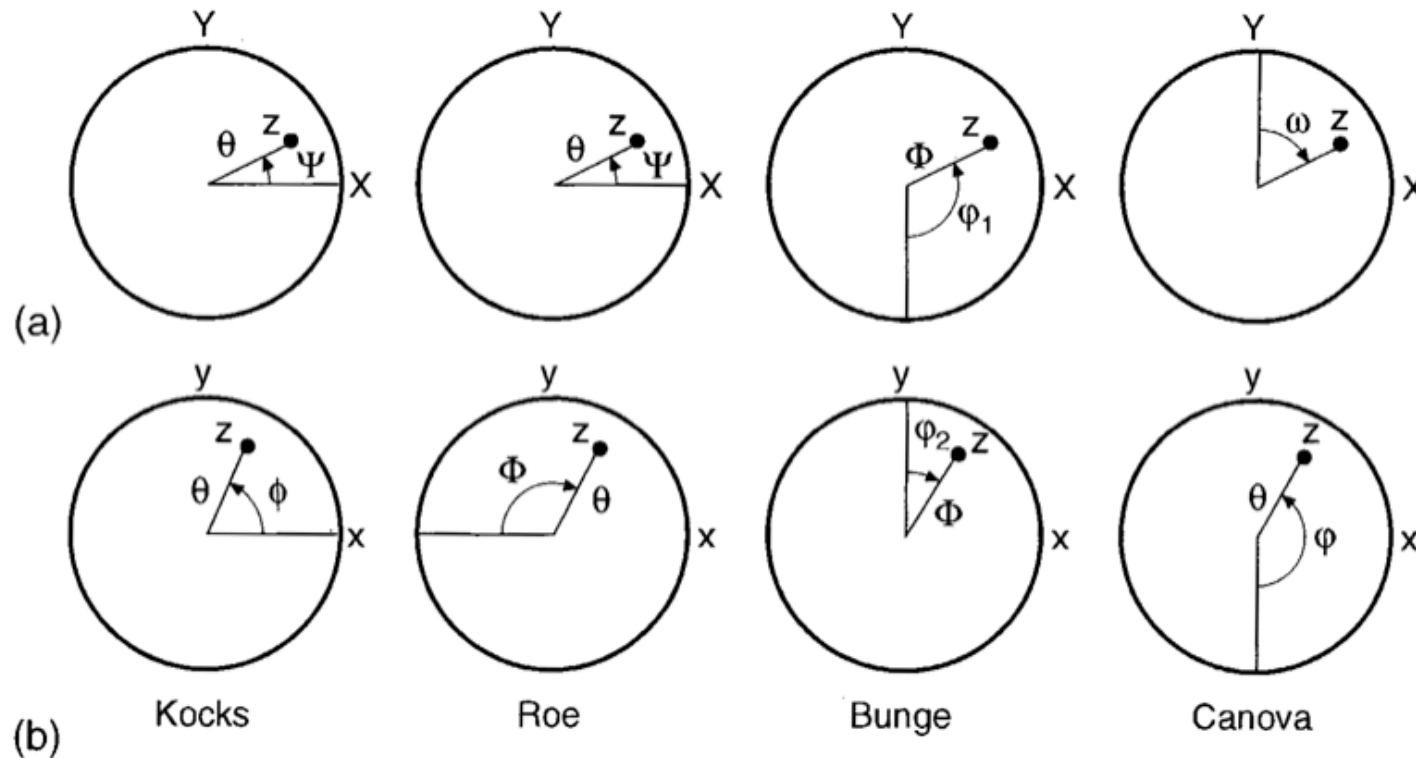
# Kocks *Euler Angles:* *Ship Analogy*

[Kocks,  
Tomé,  
Wenk]

Analogy: position and the heading of a boat with respect to the globe. *Latitude* ( $\theta$ ) and *longitude* ( $\psi$ ) describe the position of the boat; third angle describes the *heading* ( $\phi$ ) of the boat relative to the line of longitude that connects the boat to the North Pole.



# Euler Angle Definitions



[Kocks,  
Tomé,  
Wenk]

*Bunge* and *Canova* are inverse to one another  
*Kocks* and *Roe* differ by sign of third angle  
*Bunge* rotates about  $x'$ , *Roe/Kocks* about  $y'$  (2nd angle)

Obj/notation AxisTransformation Matrix EulerAngles Components

# *Euler Angle Conversions*

Convention	1 <sup>st</sup>	2 <sup>nd</sup>	3 <sup>rd</sup>	2 <sup>nd</sup> angle about axis:
Kocks (symmetric)	$\Psi$	$\Theta$	$\phi$	y
Bunge	$\phi_1 - \pi/2$	$\Phi$	$\pi/2 - \phi_2$	x
Matthies	$\alpha$	$\beta$	$\pi - \gamma$	y
Roe	$\Psi$	$\Theta$	$\pi - \Phi$	y

*[Kocks, Tomé, Wenk]*

Obj/notation AxisTransformation Matrix **EulerAngles** Components



# Complete orientations in the Pole

## Figure

Note the loss of information in a diffraction experiment if each set of poles from a single component cannot be related to one another.

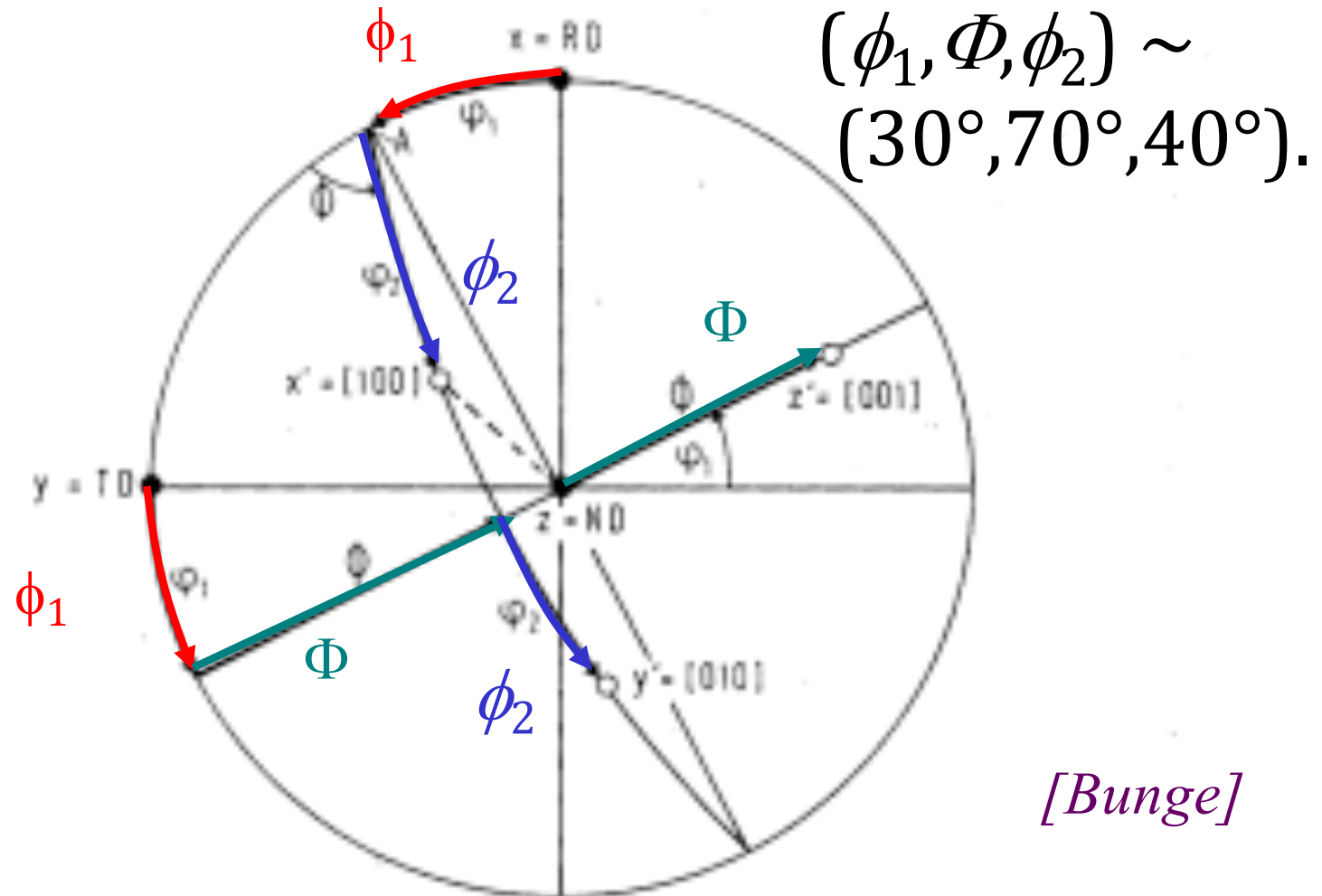
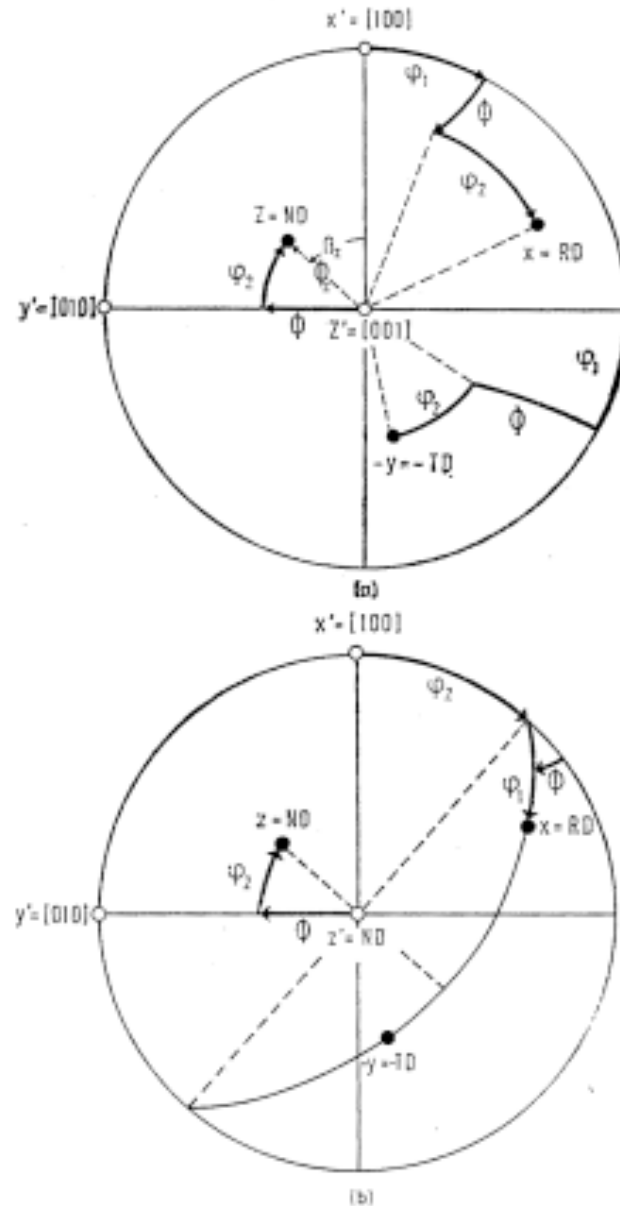


Figure 2.22 The orientation of the crystal directions  $X' = [100]$ ,  $Y' = [010]$ ,  $Z' = [001]$  in the stereographic projection in the sample coordinate system (pole figure) expressed by the Euler angles  $\varphi_1, \Phi, \varphi_2$

Obj/notation AxisTransformation Matrix EulerAngles Components

# Complete orientations in the Inverse Pole Figure

Think of yourself as an observer standing on the crystal axes, and measuring where the sample axes lie in relation to the crystal axes.



[Bunge]

Figure 2.23 The orientation of the sample direction  $X = RD$ ,  $Y = TD$ ,  $Z = ND$  in the stereographic projection in the crystal coordinate system (inverse pole figure) expressed by the EULER angles  $\varphi_1, \varphi_2$ .

Obj/notation AxisTransformation Matrix EulerAngles Components

# Summary

- Conversion between different forms of description of *texture components* described.
- Physical picture of the meaning of Euler angles as rotations of a crystal given.
- Miller indices are descriptive (and typical in physical metallurgy), but *matrices* are useful for computation, and *Euler angles* are useful for mapping out textures (to be discussed). Later on we will describe *Rodrigues vectors* and (unit) *quaternions*: the latter are the most useful representation for computations.

# *In-Class Exercises: 1*

1. Draw three orthogonal axes to represent the reference frame. Draw a unit cube to represent a crystal such that the  $\langle 100 \rangle$  directions are aligned with the  $\{x,y,z\}$  axes. Explain that there are several alternative labels for such sets of axes or frames-of-reference, such as  $\{RD,TD,ND\}$ . Distinguish between the use of  $\{ \}$  to denote a family of plane normals and  $\{ \}$  to denote a set of objects (which may, or may not, constitute a group). Explain that this alignment is called the “cube component”.
2. Starting with the reference frame, also called the sample frame, sketch the 3 successive rotations that correspond to the (Bunge) Euler angles,  $\{\phi_1, \Phi, \phi_2\}$ . Label the axes of rotation for each one. Explain that it is permissible to describe this process in terms of rotations about the crystal z axis, then the x axis, then the z axis again.
3. Draw the sample frame, with labels, and then draw a unit cube with one edge parallel to the sample x axis and rotated by  $45^\circ$  about this axis. State that this is a graphical representation of the texture component known as “Goss”. Identify the Miller indices of the three crystal directions parallel to  $\{RD, TD, ND\}$  as  $\{100,011,0-11\}$ .

## *In-Class Exercises: 2*

1. For the Goss component shown previously, explain that the Euler angles that parameterize this texture component are  $\{0,45,0\}$  (in degrees). Refer to the previous explanation of the definition of the Bunge Euler angles to do this.
2. For the same Goss component, draw the  $\{110\}$  pole figure. Hint: a simple way to do this is to find a diagram of the stereographic projection for cubic materials with 110 in the center of the figure; rotate it until a  $\langle 100 \rangle$  direction is aligned pointing to the right. In all our work we will align the x axis of a Cartesian reference frame pointing to the right, just as in standard mathematical plots, and the y axis pointing up.
3. For the same Goss component, introduce the rotation matrix. Setting aside its properties as an orthogonal matrix for the time being, explain how one can construct it as a  $2 \times 2$   $45^\circ$  rotation matrix, which can then be embedded in a  $3 \times 3$  matrix to represent the  $45^\circ$  rotation about the x axis.

## *In-Class Exercises: 3*

1. Introduce the Brass component by using the same 110 stereographic projection as before but rotating it so that, instead of having, say, 001//RD, now we have -112//RD. Show that the angle required to rotate from Goss to Brass is approximately  $35^\circ$ . Show that this rotation is about the center of the projection, i.e. 110.
2. Picking up from the previous exercise, show that the combination of the  $45^\circ$  rotation about x to arrive at the Goss component, combined with the  $35^\circ$  rotation about 110 means that we can describe the Brass component with Euler angles  $\{35,45,0\}$ .
3. Show, again with the aid of the projection, that the 3 crystal directions parallel to the sample frame are  $\{-112, 1-11, 110\}$ . Check that these form a right-handed set!
4. Draw the definition of the Roe Euler angles. Explain the differences between this definition and the Bunge definition.
5. Draw the definition of the Kocks convention for Euler angles. Explain the differences between this definition and the Bunge definition.
6. List the conversions between the different definitions of Euler angles.

## *In-Class Exercises: 4*

1. Setting aside the precise definition of a pole figure for the time being, sketch a pole figure with the sample X pointing to the right, the Y up and the Z in the center. (Later on we will discuss the confusing practice of drawing pole figures with RD up and TD to the right.) Sketch the successive rotations that correspond to the 3 (Bunge) Euler angles.
2. Sketch an inverse pole figure with the crystal x (i.e. 100) pointing to the right, the y (010) up and the z (001) in the center. Sketch the successive rotations that correspond to the 3 (Bunge) Euler angles. You will find it easier to start with the last angle,  $\phi_2$ , and work backwards.
3. Anticipating the next lecture, discuss how to use the information about  $hkl//Z$  and  $uvw//X$  to construct a set of 3 mutually orthogonal unit vectors as a preliminary to obtaining an orientation matrix.

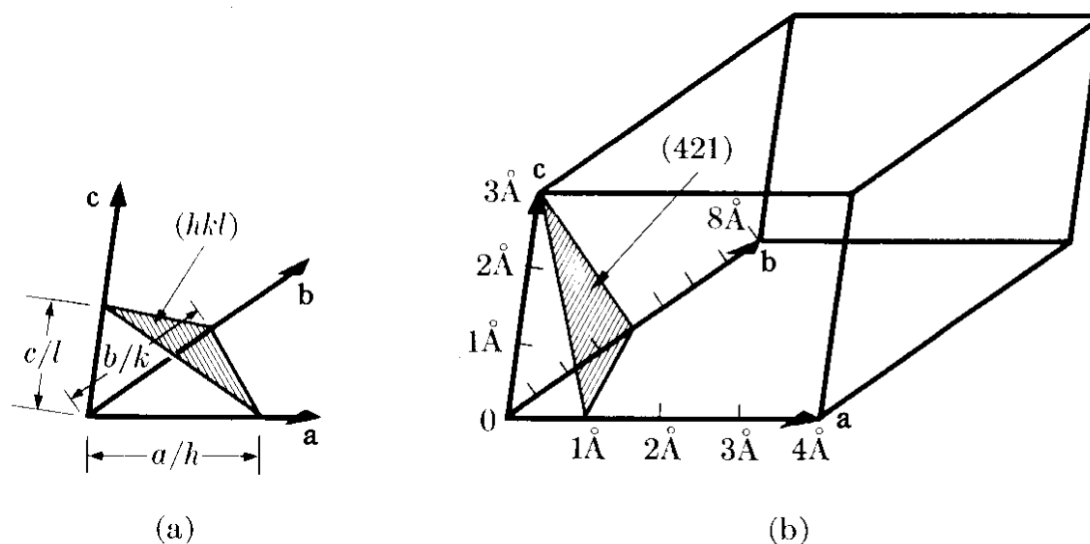
## *Supplementary Slides*

- The following slides provide supplementary information on Miller indices, the dot (scalar) product and direction cosines.



# Miller Indices

- Cubic system: directions,  $[uvw]$ , are equivalent to, and parallel to plane normals with the same indices,  $(hkl)$ .
- Miller indices for a plane specify reciprocals of intercepts on each axis.



**Fig. 2-9** Plane designation by Miller indices.

## Miller $\leftrightarrow$ vectors

- Miller indices [integer representation of direction cosines] can be converted to a *unit vector*,  $\mathbf{n}$ , by dividing by the square root of the sum of the squares: {similar for [uvw]}. This is known as *normalization*.

$$\hat{\mathbf{n}} = \frac{(h, k, l)}{\sqrt{h^2 + k^2 + l^2}}$$

# *Miller Index Definition of a Texture Component*

- The commonest method for specifying a texture component is the plane-direction.
- Specify the crystallographic plane normal that is parallel to the specimen normal (e.g. the ND $\equiv$ Z) and a crystallographic direction that is parallel to the long direction (e.g. the RD $\equiv$ X).

$(hkl) // \text{ND}, [uvw] // \text{RD}, \text{ or } (hkl)[uvw]$

# *Dot Product*

- Given two vectors, **a** and **b**, the dot product, **a•b** is a scalar quantity that is equal to the product of the magnitudes (lengths) of the vectors, multiplied by the cosine of the angle between them:

$$\mathbf{a} \cdot \mathbf{b} = a b \cos \theta$$

- If both vectors are unit vectors then the dot product is equal to the cosine of the angle.
- In index form,  $\mathbf{a} \cdot \mathbf{b} = a_i b_i$ .
- Given a set of unit vectors defining an axis system,  $e_x, e_y, e_z$ , a vector can be defined on that system by taking the dot product with each axis vector in turn, e.g.:

$$a_x = \mathbf{a} \cdot e_x$$

# *Direction Cosines*

- Definition of *direction cosines*: In analytic geometry, the direction cosines of a vector are the cosines of the angles between the vector and the three coordinate axes. Equivalently, they are the contributions of each component of the basis to a unit vector in that direction [[en.wikipedia.org/wiki/Direction\\_cosine](https://en.wikipedia.org/wiki/Direction_cosine)].
- The components of a unit vector are equal to the cosines of the angle between the vector and each (orthogonal, Cartesian) reference axis.
- We can use *axis transformations* to describe vectors in different reference frames (room, specimen, crystal, slip system....)